

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041122\
 Data File : VY008272.D
 Acq On : 11 Apr 2022 14:56
 Operator : KP/MD
 Sample : N2340-01
 Misc : 5.08g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 SB-04-14.5-15.0

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\82Y040822S.M

Title : SW846 8260

Signal : TIC: VY008272.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.716	958	967	973	rBV2	13221	30942	3.75%	0.516%
2	7.789	973	979	988	rVB	24640	55089	6.68%	0.918%
3	8.143	1029	1037	1047	rBV2	12035	25679	3.11%	0.428%
4	8.691	1118	1127	1138	rVB	33285	64543	7.83%	1.075%
5	10.179	1364	1371	1380	rBV	55172	95931	11.64%	1.598%
6	10.904	1484	1490	1497	rBV2	4793	9412	1.14%	0.157%
7	11.038	1505	1512	1516	rVV3	5662	8994	1.09%	0.150%
8	11.093	1516	1521	1526	rVB4	4790	9070	1.10%	0.151%
9	11.270	1548	1550	1562	rVB5	6470	15140	1.84%	0.252%
10	11.380	1562	1568	1573	rBV3	7457	13334	1.62%	0.222%
11	11.490	1580	1586	1595	rVB	49529	81817	9.92%	1.363%
12	11.739	1622	1627	1630	rBV	6650	11483	1.39%	0.191%
13	11.782	1631	1634	1639	rVB3	5309	8519	1.03%	0.142%
14	12.002	1663	1670	1678	rBV7	8586	22307	2.71%	0.372%
15	12.124	1686	1690	1694	rVB2	8963	12232	1.48%	0.204%
16	12.203	1699	1703	1706	rVV5	7895	13308	1.61%	0.222%
17	12.246	1706	1710	1715	rVB3	13341	24008	2.91%	0.400%
18	12.331	1720	1724	1727	rBV2	12604	21203	2.57%	0.353%
19	12.392	1727	1734	1735	rVV5	11486	26275	3.19%	0.438%
20	12.416	1735	1738	1744	rVV	17293	31365	3.80%	0.523%
21	12.483	1744	1749	1754	rVB2	47334	73768	8.95%	1.229%
22	12.672	1770	1780	1784	rVB	21812	52158	6.33%	0.869%
23	12.721	1784	1788	1794	rBV6	6749	14986	1.82%	0.250%
24	12.812	1794	1803	1807	rBV3	23972	58966	7.15%	0.982%
25	12.861	1807	1811	1816	rVB4	12110	19126	2.32%	0.319%
26	12.953	1816	1826	1829	rBV8	21482	53029	6.43%	0.883%
27	13.044	1837	1841	1847	rVB	41538	59560	7.22%	0.992%
28	13.105	1849	1851	1858	rVB7	9090	18479	2.24%	0.308%
29	13.172	1858	1862	1866	rVB5	9452	12810	1.55%	0.213%
30	13.227	1867	1871	1872	rBV2	11002	13802	1.67%	0.230%
31	13.251	1872	1875	1879	rVB2	22630	31504	3.82%	0.525%
32	13.325	1879	1887	1890	rBV5	41000	83174	10.09%	1.386%
33	13.361	1890	1893	1897	rVV3	25871	39593	4.80%	0.660%
34	13.428	1897	1904	1914	rVB	48221	95239	11.55%	1.587%
35	13.568	1922	1927	1928	rBV4	12886	17300	2.10%	0.288%
36	13.593	1929	1931	1934	rVV	19031	25029	3.04%	0.417%

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Peak Location: TOP

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Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y040822S.M
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37	13.629	1934	1937	1941	rVB	21359	28630	3.47%	0.477%
38	13.696	1941	1948	1952	rBV2	22147	53001	6.43%	0.883%
39	13.751	1952	1957	1961	rBV3	58536	102939	12.49%	1.715%
40	13.794	1961	1964	1968	rVB	26714	35130	4.26%	0.585%
41	13.855	1968	1974	1977	rBV4	22521	46069	5.59%	0.768%
42	13.916	1981	1984	1989	rBV6	17281	28274	3.43%	0.471%
43	13.971	1989	1993	1998	rVB	80232	116712	14.16%	1.944%
44	14.020	1998	2001	2006	rVB2	30705	48068	5.83%	0.801%
45	14.081	2006	2011	2016	rVB4	64178	105061	12.74%	1.750%
46	14.148	2016	2022	2027	rBV7	30977	60630	7.35%	1.010%
47	14.202	2027	2031	2033	rBV3	17303	26037	3.16%	0.434%
48	14.263	2033	2041	2045	rBV4	95020	222070	26.94%	3.700%
49	14.337	2050	2053	2057	rVB	32874	44367	5.38%	0.739%
50	14.379	2057	2060	2063	rBV2	32949	45143	5.48%	0.752%
51	14.422	2063	2067	2071	rVV2	47219	63984	7.76%	1.066%
52	14.477	2074	2076	2080	rVB3	16391	21539	2.61%	0.359%
53	14.568	2086	2091	2096	rBV2	55619	117881	14.30%	1.964%
54	14.617	2096	2099	2103	rVB2	35608	43407	5.27%	0.723%
55	14.684	2103	2110	2114	rBV3	168402	353384	42.86%	5.888%
56	14.733	2114	2118	2125	rVB2	139865	241089	29.24%	4.017%
57	14.806	2126	2130	2133	rBV7	20183	33386	4.05%	0.556%
58	14.910	2144	2147	2149	rBV2	22224	27150	3.29%	0.452%
59	14.946	2149	2153	2157	rVV3	73902	118064	14.32%	1.967%
60	14.983	2157	2159	2162	rVV	29621	32220	3.91%	0.537%
61	15.056	2166	2171	2177	rVB2	67647	133247	16.16%	2.220%
62	15.214	2191	2197	2203	rBV4	93031	168357	20.42%	2.805%
63	15.294	2206	2210	2214	rVB2	45778	66632	8.08%	1.110%
64	15.349	2215	2219	2223	rBV3	46481	86880	10.54%	1.447%
65	15.531	2242	2249	2256	rBV3	71581	181309	21.99%	3.021%
66	15.611	2257	2262	2266	rVB4	49910	86605	10.50%	1.443%
67	15.690	2268	2275	2278	rVV5	48745	121244	14.71%	2.020%
68	15.733	2278	2282	2291	rVB	80877	157568	19.11%	2.625%
69	15.891	2302	2308	2314	rBV3	40821	98337	11.93%	1.638%
70	16.038	2326	2332	2336	rBV3	62290	122171	14.82%	2.035%
71	16.159	2348	2352	2357	rBV3	39502	68497	8.31%	1.141%
72	16.233	2360	2364	2368	rVV3	58722	124989	15.16%	2.082%
73	16.294	2368	2374	2388	rVB	400318	824417	100.00%	13.735%
74	16.489	2399	2406	2420	rVB	208222	498562	60.47%	8.306%

Sum of corrected areas: 6002224

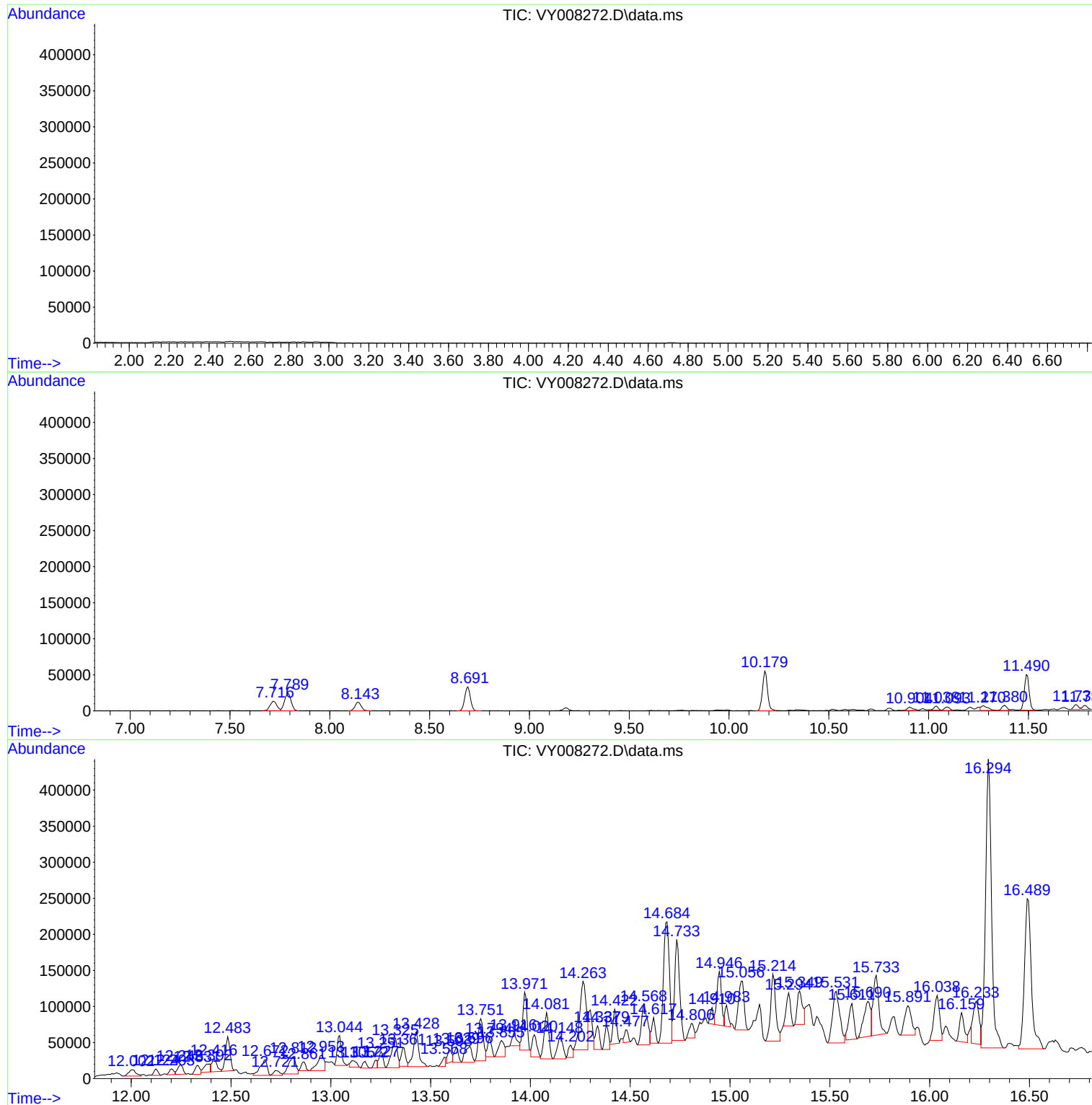
Instrument :
MSVOA_Y
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y040822S.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_Y
ClientSampleId :
SB-04-14.5-15.0

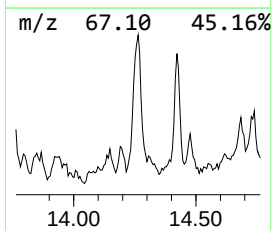
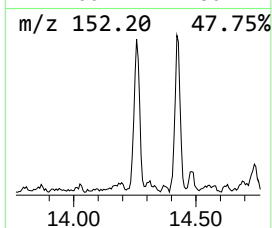
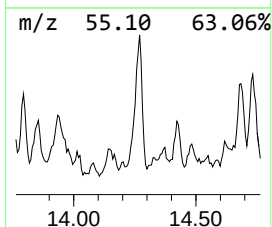
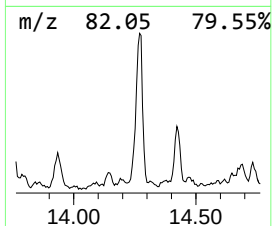
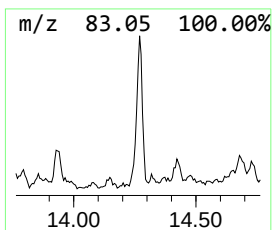
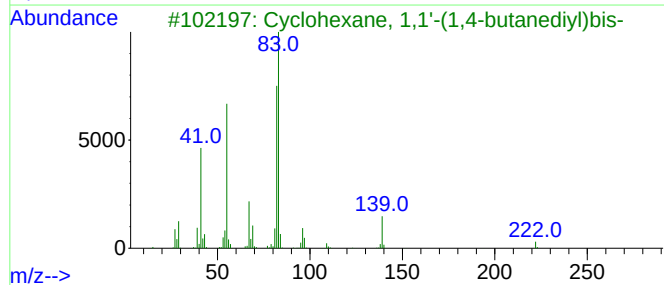
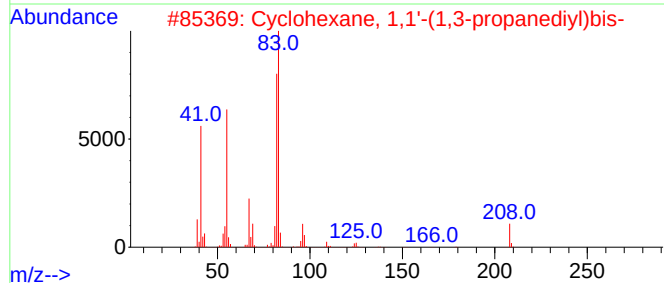
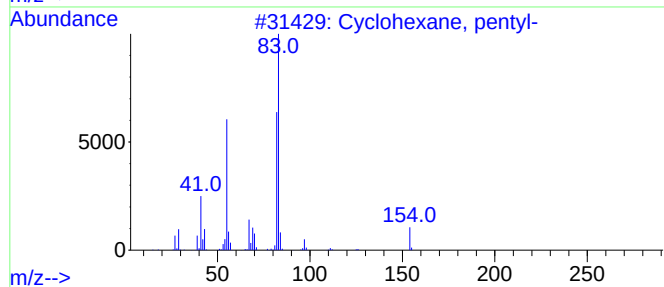
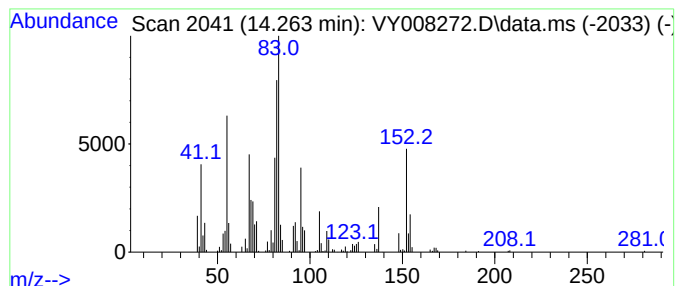
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y040822S.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, pentyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.263	116.59 ug/l	222070	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, pentyl-	154	C11H22	004292-92-6	55
2			Cyclohexane, 1,1'-(1,3-propanedi...	208	C15H28	003178-24-3	53
3			Cyclohexane, 1,1'-(1,4-butanedi...	222	C16H30	006165-44-2	43
4			Hexane, 1,6-dicyclohexyl-	250	C18H34	001610-23-7	43
5			Cyclohexane, octyl-	196	C14H28	001795-15-9	43



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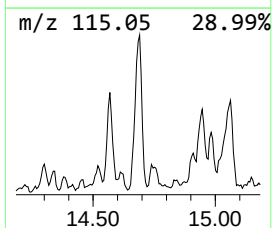
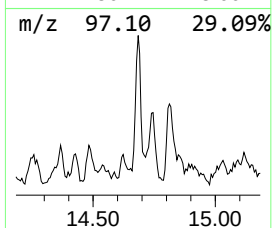
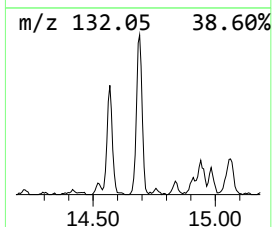
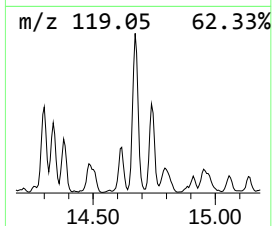
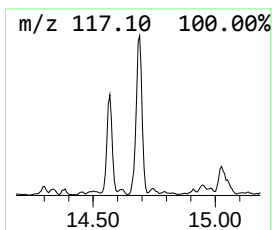
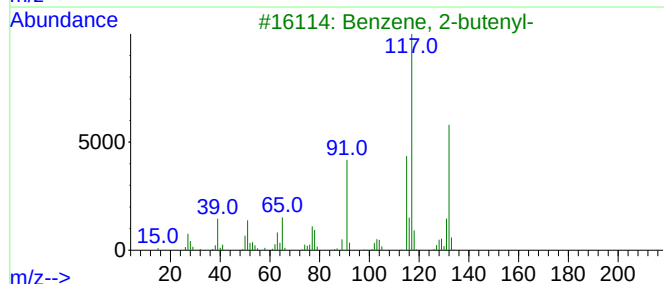
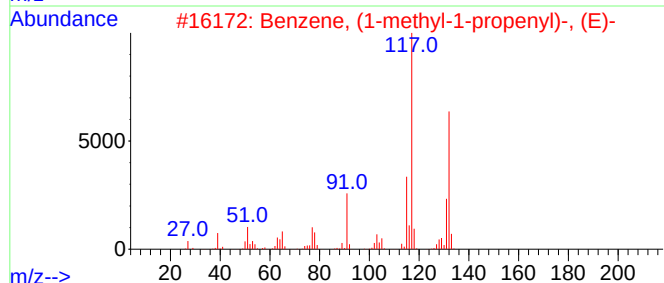
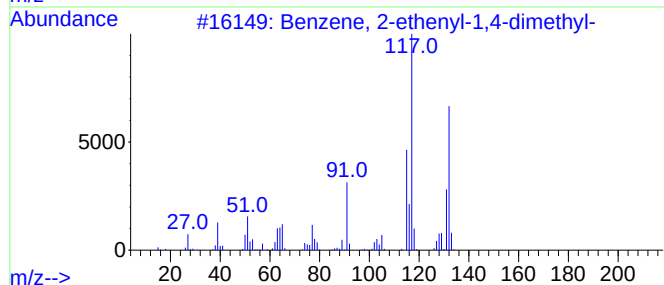
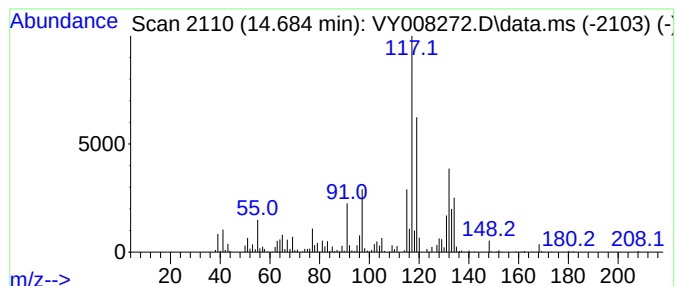
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y040822S.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	185.53 ug/l	353384	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	90
2			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	86
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
4			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64
5			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	64



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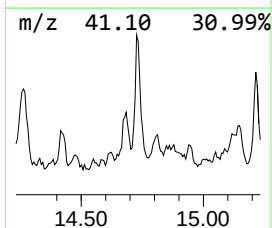
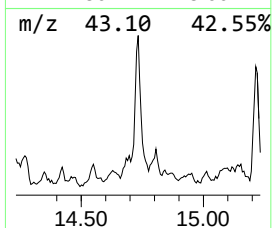
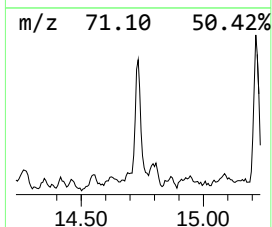
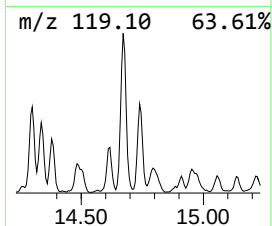
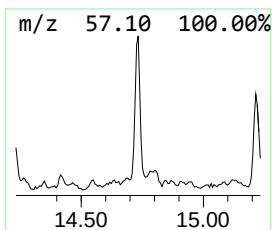
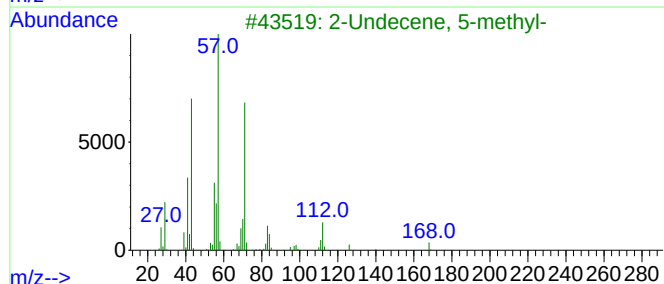
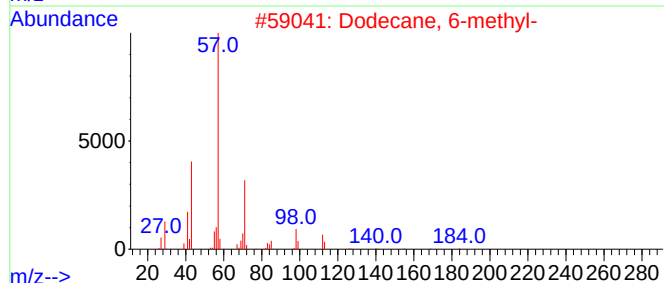
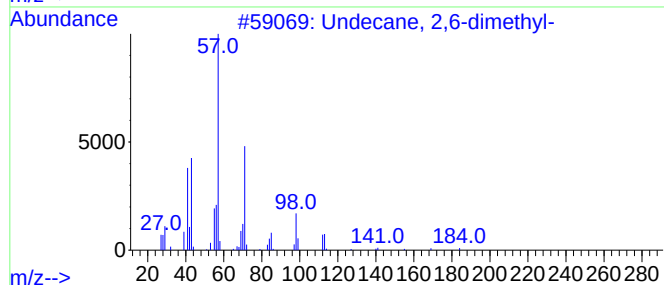
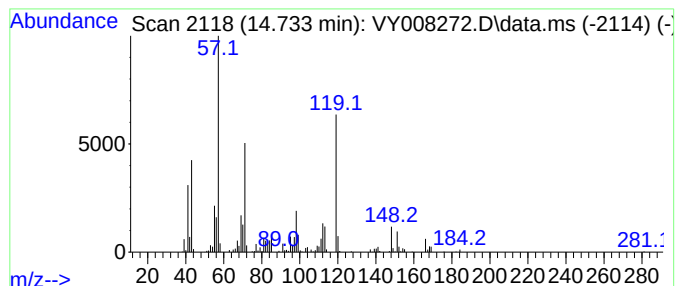
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Undecane, 2,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.733	126.57 ug/l	241089	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	60
2			Dodecane, 6-methyl-	184	C13H28	006044-71-9	41
3			2-Undecene, 5-methyl-	168	C12H24	056851-34-4	38
4			Undecane, 6-ethyl-	184	C13H28	017312-60-6	38
5			Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	35



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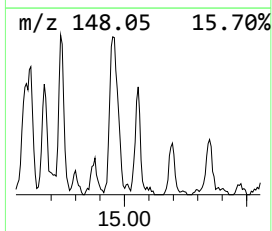
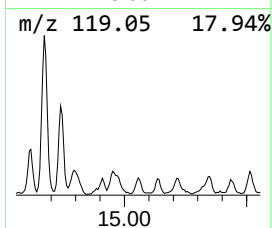
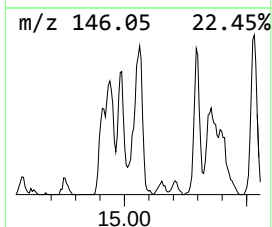
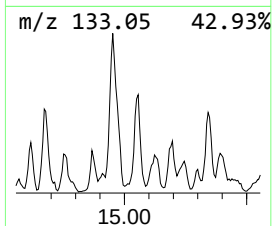
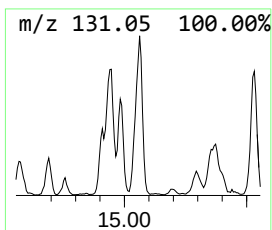
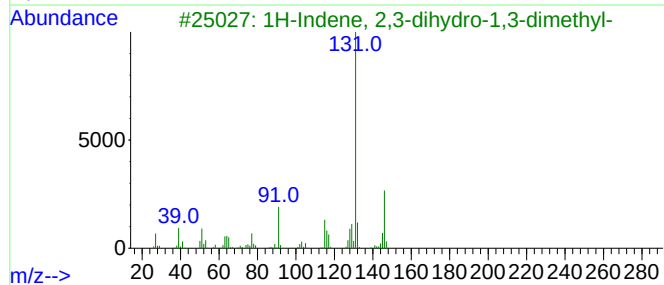
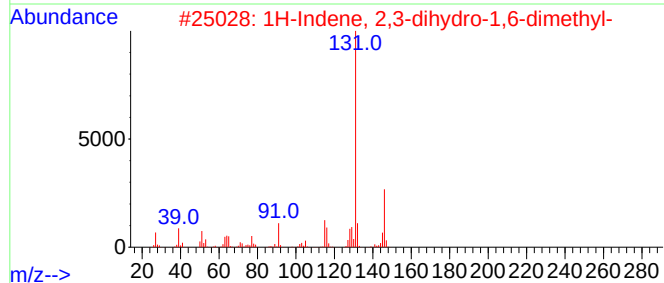
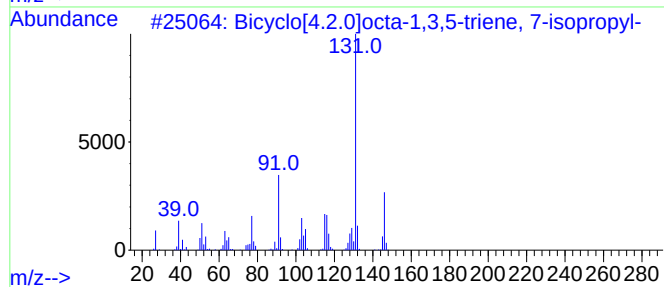
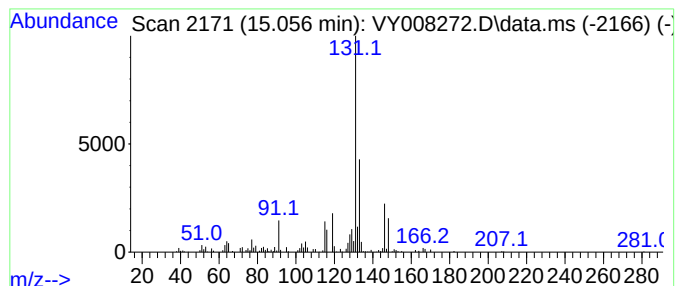
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y040822S.M
Quant Title : SW846 8260

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

Peak Number 4 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.056	69.95 ug/l	133247	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	83
2			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	64
3			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	64
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	64
5			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041122\
Data File : VY008272.D
Acq On : 11 Apr 2022 14:56
Operator : KP/MD
Sample : N2340-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-04-14.5-15.0

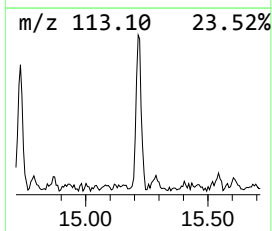
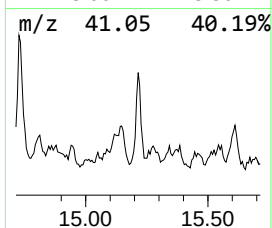
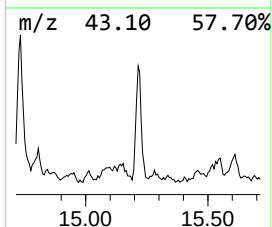
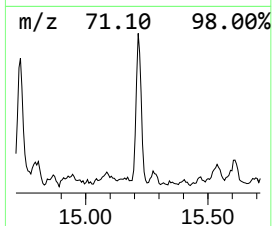
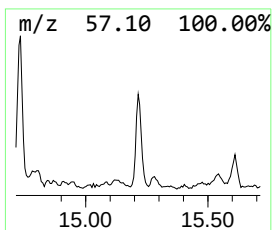
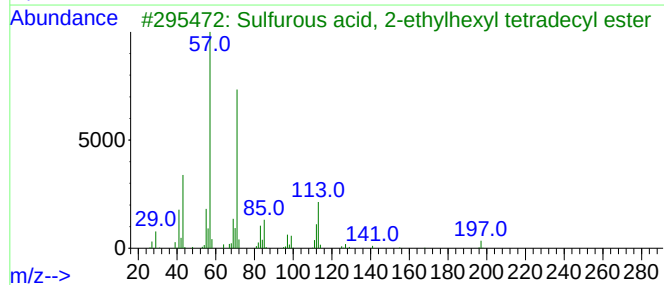
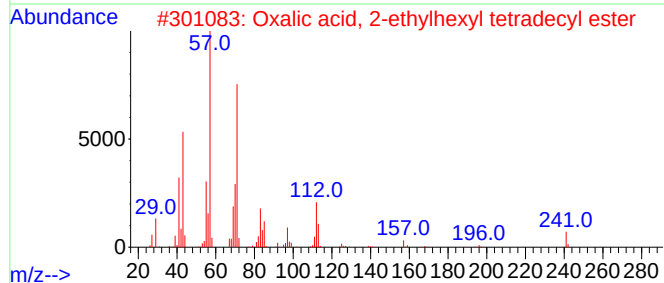
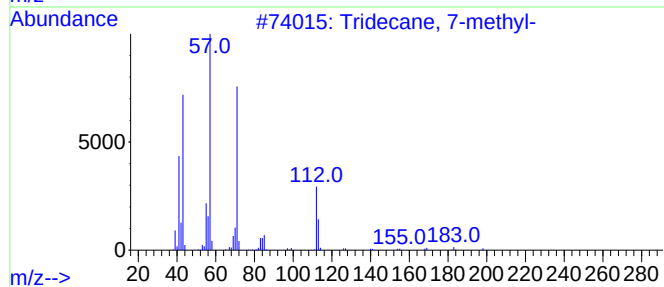
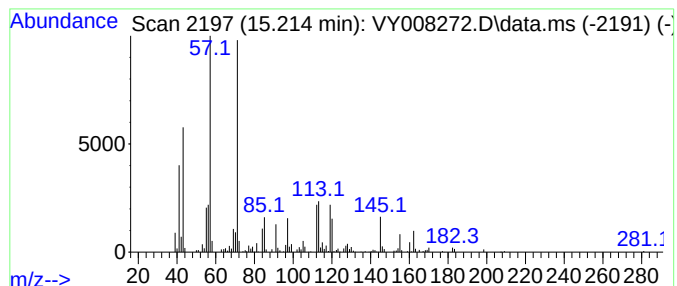
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Quant Title : SW846 8260

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

Peak Number 5 Tridecane, 7-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.214	88.39 ug/l	168357	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 7-methyl-	198	C14H30	026730-14-3	60
2			Oxalic acid, 2-ethylhexyl tetrad...	398	C24H46O4	1000309-39-7	58
3			Sulfurous acid, 2-ethylhexyl tet...	390	C22H46O3S	1000309-19-7	50
4			Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	50
5			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041122\
Data File : VY008272.D
Acq On : 11 Apr 2022 14:56
Operator : KP/MD
Sample : N2340-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-04-14.5-15.0

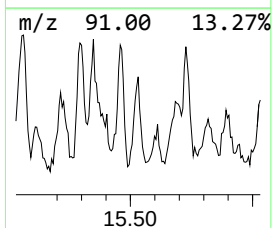
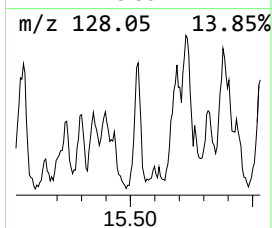
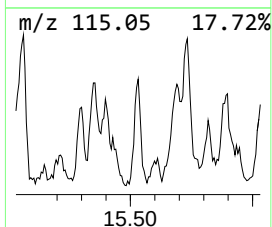
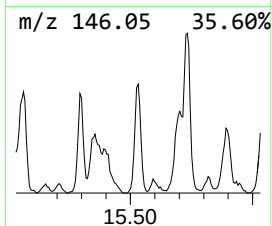
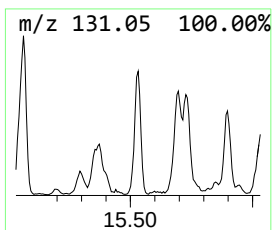
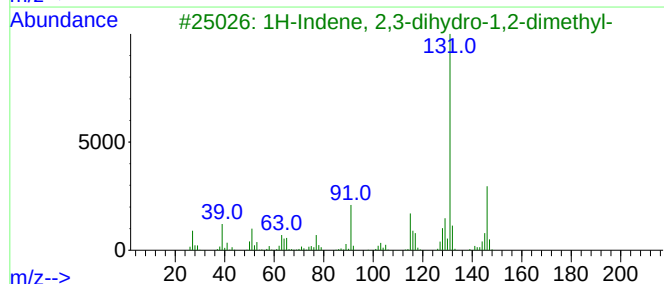
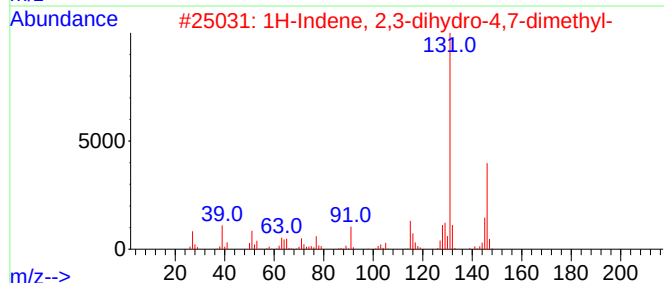
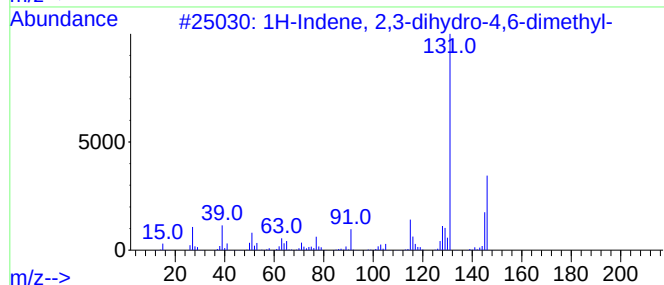
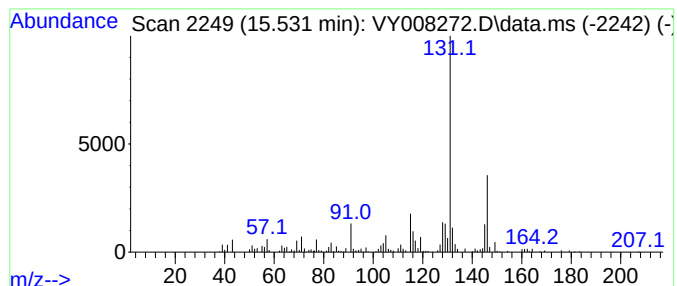
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Quant Title : SW846 8260

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

Peak Number 6 1H-Indene, 2,3-dihydro-4,6-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.531	95.19 ug/l	181309	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	93		
2	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93		
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90		
4	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90		
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041122\
Data File : VY008272.D
Acq On : 11 Apr 2022 14:56
Operator : KP/MD
Sample : N2340-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-04-14.5-15.0

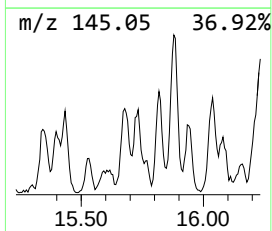
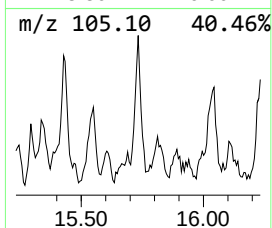
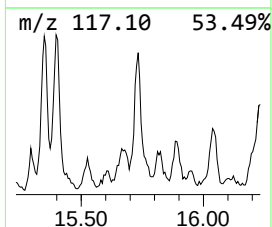
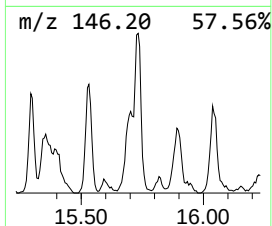
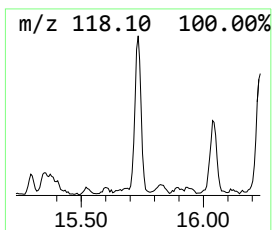
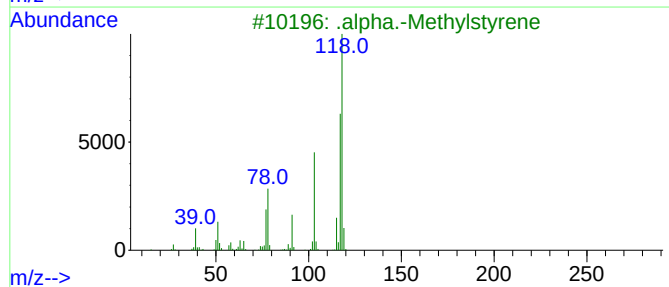
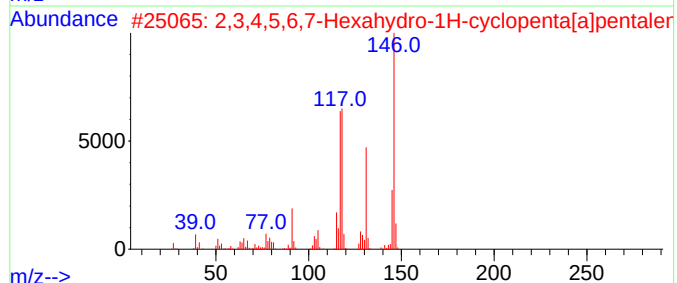
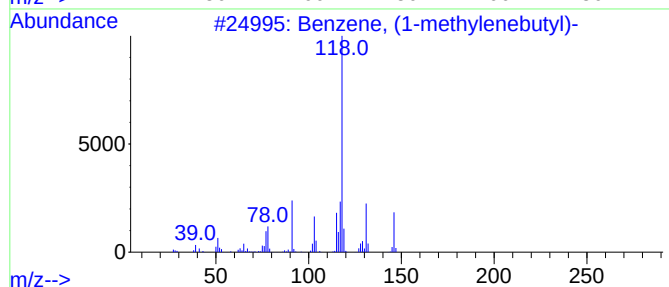
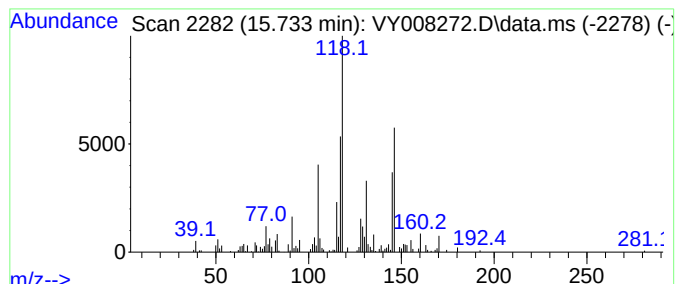
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y040822S.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, (1-methylenebutyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.733	82.72 ug/l	157568	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	59
2			2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	49
3			.alpha.-Methylstyrene	118	C9H10	000098-83-9	43
4			Cyclobutanone, 3-phenylmethyl-	160	C11H12O	055262-02-7	43
5			7-Methylindan-1-one	146	C10H10O	039627-61-7	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041122\
Data File : VY008272.D
Acq On : 11 Apr 2022 14:56
Operator : KP/MD
Sample : N2340-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-04-14.5-15.0

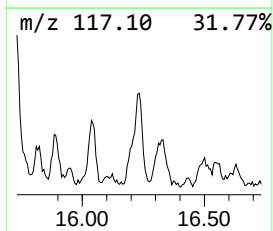
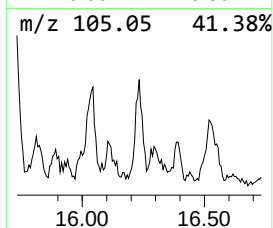
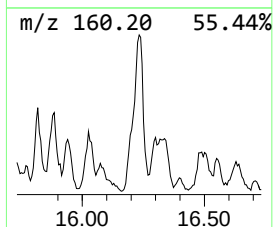
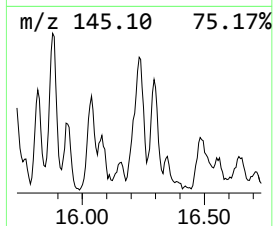
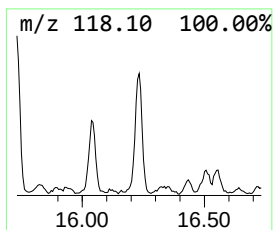
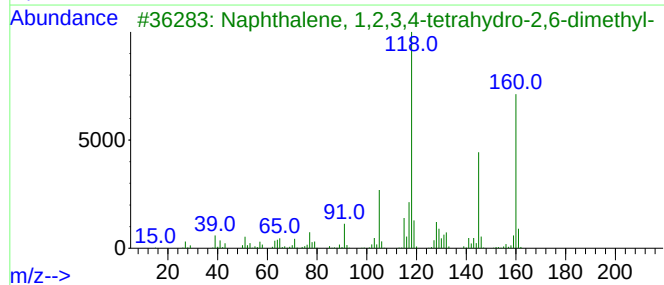
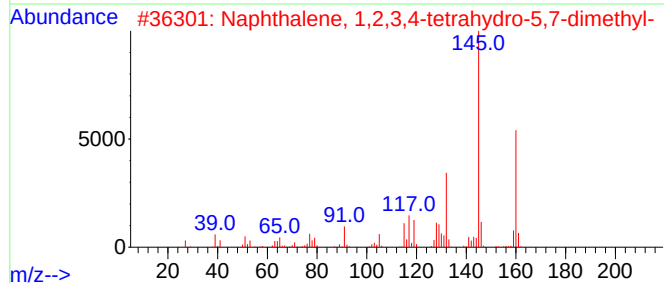
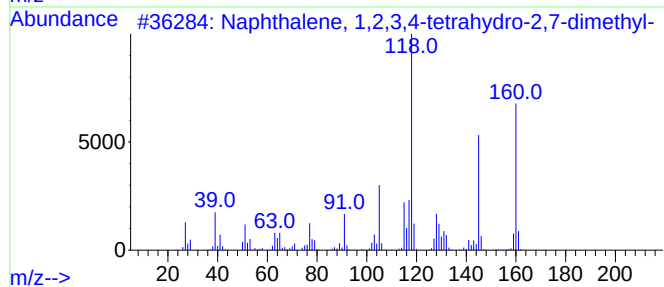
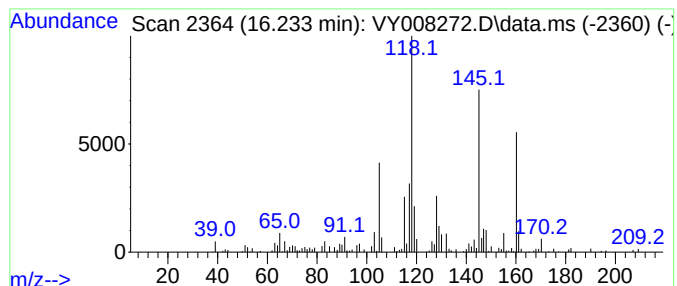
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.233	65.62 ug/l	124989	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	83
2			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021693-54-9	53
3			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	49
4			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001076-61-5	49
5			Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041122\
Data File : VY008272.D
Acq On : 11 Apr 2022 14:56
Operator : KP/MD
Sample : N2340-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-04-14.5-15.0

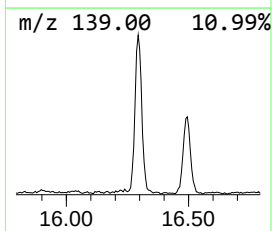
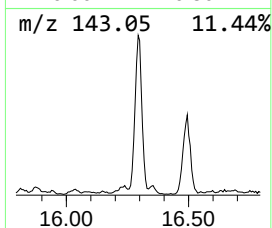
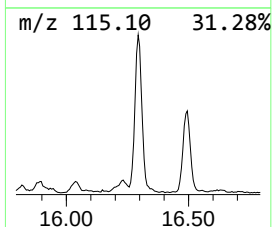
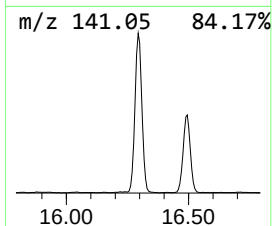
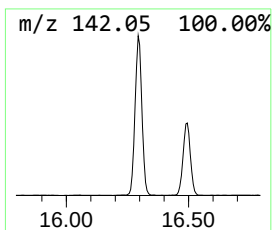
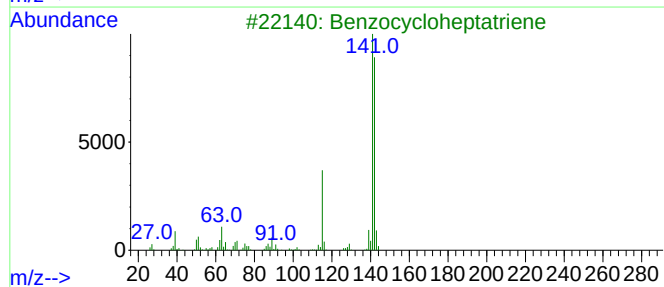
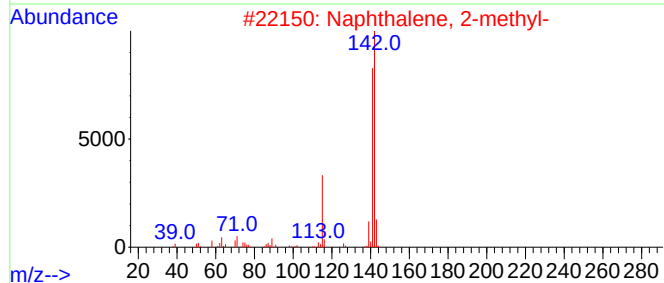
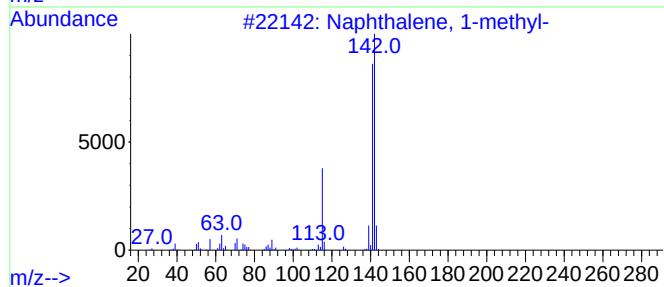
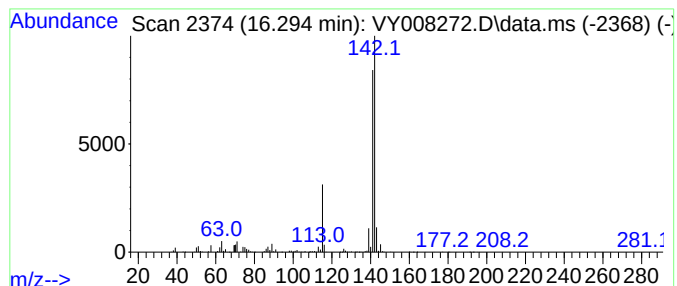
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y040822S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.294	432.82 ug/l	824417	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041122\
Data File : VY008272.D
Acq On : 11 Apr 2022 14:56
Operator : KP/MD
Sample : N2340-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-04-14.5-15.0

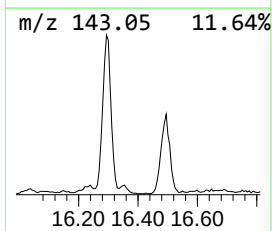
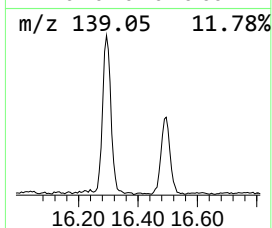
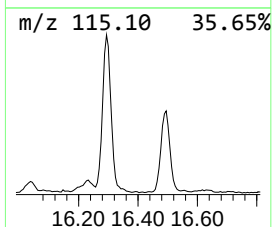
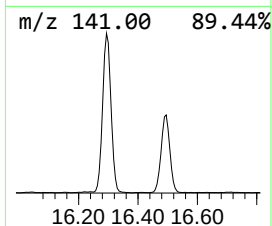
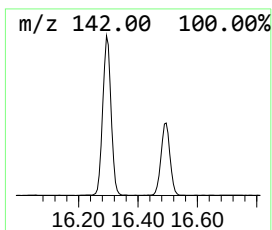
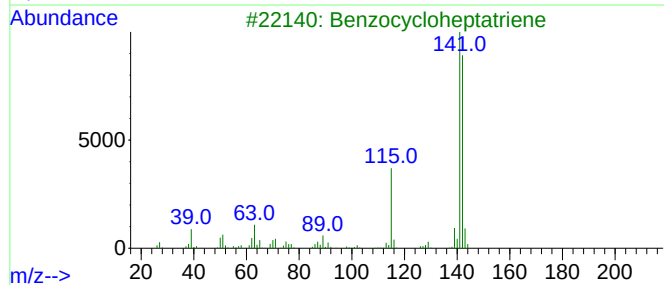
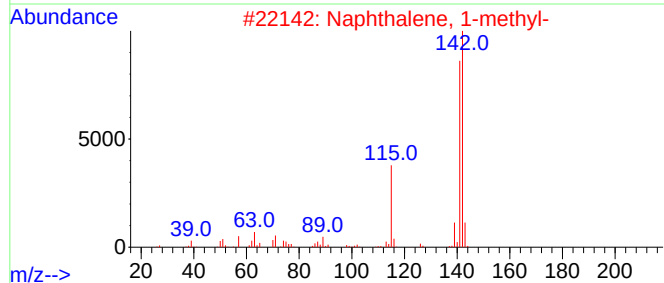
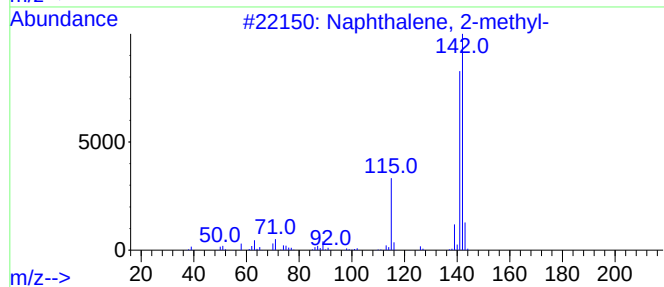
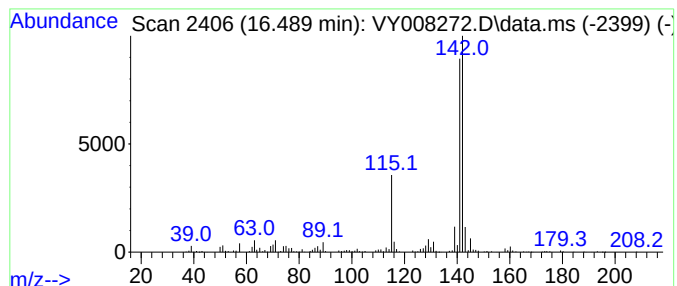
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.489	261.74 ug/l	498562	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
5			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY041122\
Data File : VY008272.D
Acq On : 11 Apr 2022 14:56
Operator : KP/MD
Sample : N2340-01
Misc : 5.08g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-04-14.5-15.0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y040822S.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, 2-ethe...	14.684	185.5	ug/l	353384	4	13.428	95239	50.0
Undecane, 2,6-d...	14.733	126.6	ug/l	241089	4	13.428	95239	50.0
Bicyclo[4.2.0]o...	15.056	70.0	ug/l	133247	4	13.428	95239	50.0
Tridecane, 7-me...	15.214	88.4	ug/l	168357	4	13.428	95239	50.0
1H-Indene, 2,3-...	15.531	95.2	ug/l	181309	4	13.428	95239	50.0
Benzene, (1-met...	15.733	82.7	ug/l	157568	4	13.428	95239	50.0
Naphthalene, 1,...	16.233	65.6	ug/l	124989	4	13.428	95239	50.0
Naphthalene, 1-...	16.294	432.8	ug/l	824417	4	13.428	95239	50.0
Naphthalene, 2-...	16.489	261.7	ug/l	498562	4	13.428	95239	50.0