

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110222\
 Data File : VY011255.D
 Acq On : 02 Nov 2022 12:25
 Operator : KP/MD
 Sample : N5337-03
 Misc : 5.06g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 SB-02-11.5-12.0

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\82Y110122S.M

Title : SW846 8260

Signal : TIC: VY011255.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	6.972	834	845	862	rBV	23499	71969	1.30%	0.126%
2	7.716	957	967	972	rBV	39163	94405	1.71%	0.165%
3	7.783	972	978	988	rVB	60228	134816	2.44%	0.236%
4	8.136	1026	1036	1046	rBV	36797	84938	1.54%	0.149%
5	8.685	1118	1126	1137	rBV	89206	173623	3.15%	0.304%
6	10.179	1363	1371	1385	rBV	126150	221853	4.02%	0.389%
7	11.288	1547	1553	1562	rVB6	31527	88539	1.60%	0.155%
8	11.489	1579	1586	1595	rBV	120508	207265	3.76%	0.363%
9	11.739	1621	1627	1631	rBV5	28607	65903	1.19%	0.116%
10	12.002	1663	1670	1676	rBV4	161611	367254	6.65%	0.644%
11	12.117	1682	1689	1694	rVB	225016	338941	6.14%	0.594%
12	12.197	1694	1702	1705	rBV3	157796	338193	6.13%	0.593%
13	12.233	1705	1708	1713	rVB3	185696	278410	5.04%	0.488%
14	12.319	1713	1722	1725	rBV10	51797	157121	2.85%	0.275%
15	12.379	1725	1732	1734	rBV4	115561	246570	4.47%	0.432%
16	12.410	1734	1737	1743	rVB2	159580	222960	4.04%	0.391%
17	12.477	1743	1748	1752	rVB5	119094	228993	4.15%	0.401%
18	12.520	1752	1755	1759	rBV2	89835	109102	1.98%	0.191%
19	12.568	1759	1763	1767	rBV5	83204	140511	2.55%	0.246%
20	12.642	1771	1775	1782	rVB4	340786	667560	12.09%	1.170%
21	12.727	1782	1789	1794	rBV7	195376	517545	9.38%	0.907%
22	12.806	1795	1802	1806	rBV6	326471	835574	15.14%	1.465%
23	12.855	1808	1810	1816	rVB3	243955	366416	6.64%	0.642%
24	12.910	1816	1819	1821	rBV2	60037	79636	1.44%	0.140%
25	12.946	1821	1825	1829	rBV4	324263	456478	8.27%	0.800%
26	13.001	1829	1834	1836	rBV2	372419	622135	11.27%	1.091%
27	13.038	1836	1840	1847	rVB	1433894	2044776	37.05%	3.584%
28	13.099	1847	1850	1853	rVB2	148050	181214	3.28%	0.318%
29	13.166	1857	1861	1866	rBV2	442058	696769	12.62%	1.221%
30	13.221	1866	1870	1872	rBV3	308730	436479	7.91%	0.765%
31	13.288	1877	1881	1883	rVV3	351556	560680	10.16%	0.983%
32	13.318	1883	1886	1889	rVV2	533968	794352	14.39%	1.392%
33	13.361	1889	1893	1896	rVV3	748113	1182075	21.42%	2.072%
34	13.398	1896	1899	1904	rVB3	381434	524286	9.50%	0.919%
35	13.495	1913	1915	1921	rVB4	247565	319411	5.79%	0.560%
36	13.562	1922	1926	1929	rBV4	278292	480163	8.70%	0.842%

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37	13.684	1941	1946	1949	rBV4	175573	418869	7.59%	0.734%
38	13.745	1950	1956	1961	rVV2	1418173	2805705	50.83%	4.918%
39	13.794	1961	1964	1968	rVV4	588118	1037060	18.79%	1.818%
40	13.855	1970	1974	1979	rVV7	661808	1623834	29.42%	2.846%
41	13.910	1979	1983	1989	rVV4	952917	2394751	43.39%	4.198%
42	13.965	1989	1992	1996	rVV5	690229	1149863	20.83%	2.016%
43	14.013	1996	2000	2005	rVB	856812	1261434	22.85%	2.211%
44	14.074	2005	2010	2015	rBV5	613476	1042612	18.89%	1.828%
45	14.141	2015	2021	2026	rBV5	706174	1347104	24.41%	2.361%
46	14.196	2026	2030	2032	rBV3	354981	524165	9.50%	0.919%
47	14.257	2032	2040	2052	rVB8	1708542	5519643	100.00%	9.675%
48	14.379	2055	2060	2062	rBV2	508662	857331	15.53%	1.503%
49	14.416	2062	2066	2070	rVB3	950886	1300964	23.57%	2.280%
50	14.544	2084	2087	2089	rBV3	180192	229804	4.16%	0.403%
51	14.617	2096	2099	2104	rBV6	282510	445754	8.08%	0.781%
52	14.684	2105	2110	2113	rVV2	757351	1290324	23.38%	2.262%
53	14.727	2113	2117	2124	rVB2	2785783	4434111	80.33%	7.773%
54	14.800	2124	2129	2133	rVB5	415599	751082	13.61%	1.317%
55	14.842	2134	2136	2138	rBV3	215697	267925	4.85%	0.470%
56	14.940	2148	2152	2156	rVB3	1144792	1789698	32.42%	3.137%
57	15.050	2165	2170	2177	rVB2	885412	1706640	30.92%	2.992%
58	15.208	2191	2196	2202	rBV2	1599357	2535990	45.94%	4.445%
59	15.342	2213	2218	2222	rBV4	502572	987858	17.90%	1.732%
60	15.434	2230	2233	2241	rVB6	310666	669787	12.13%	1.174%
61	15.525	2241	2248	2255	rBV2	793622	1954199	35.40%	3.426%
62	15.605	2256	2261	2266	rVB4	647784	1168912	21.18%	2.049%
63	15.879	2301	2306	2311	rVB3	346447	725437	13.14%	1.272%
64	16.025	2326	2330	2335	rVB3	309099	537181	9.73%	0.942%
65	16.153	2347	2351	2355	rBV	279551	432222	7.83%	0.758%
66	16.470	2398	2403	2419	rVB8	337740	1098209	19.90%	1.925%
67	16.592	2419	2423	2437	rVB8	116659	402575	7.29%	0.706%

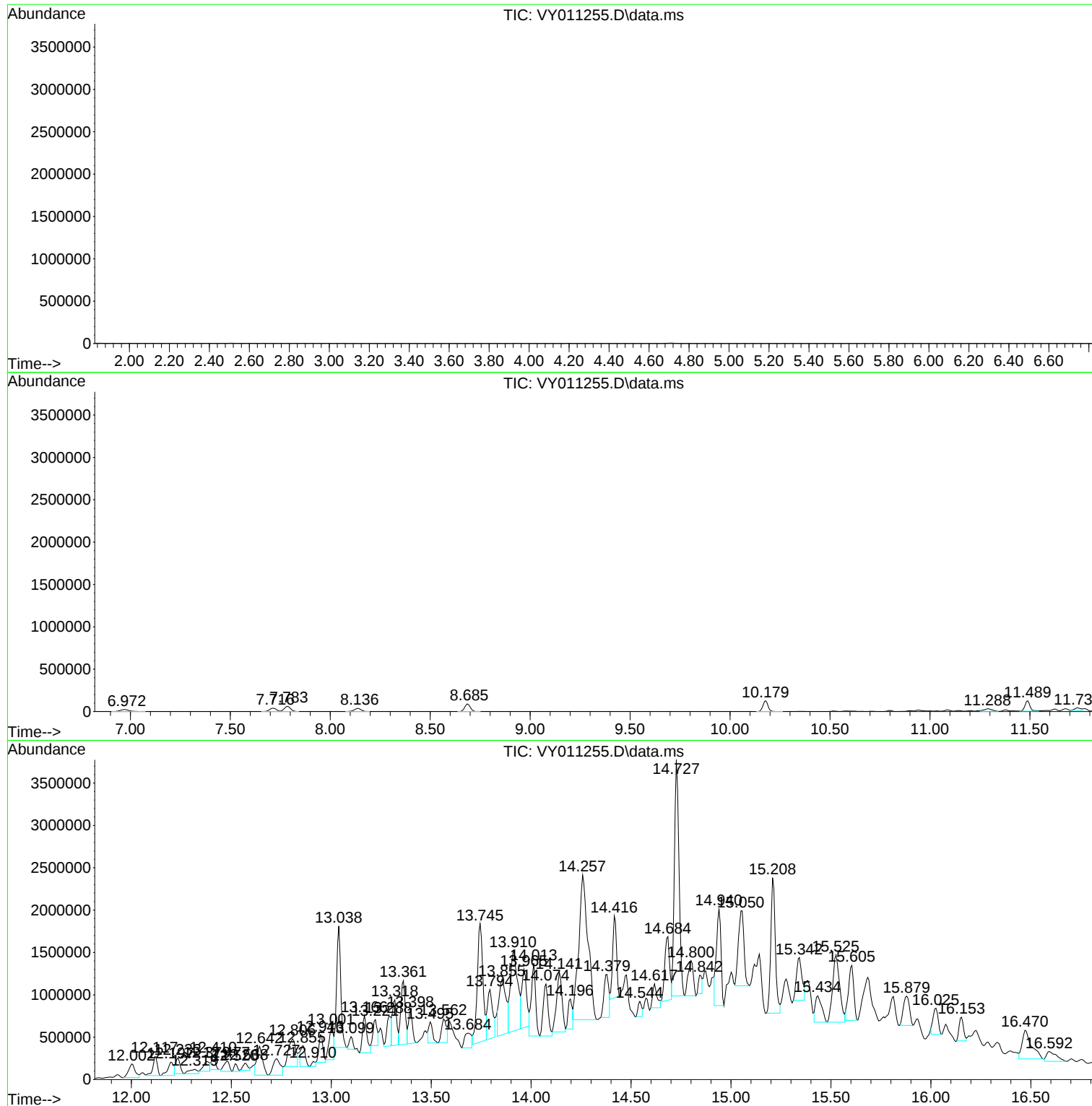
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y110122S.M
Quant Title : SW846 8260

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



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ClientSampleId :
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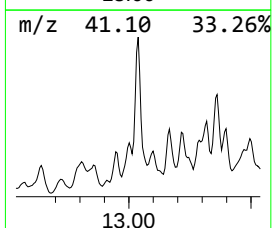
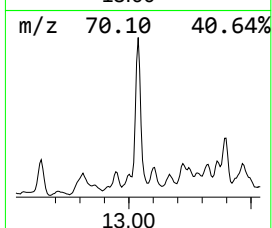
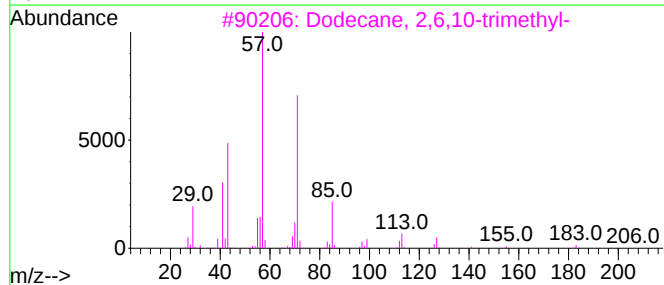
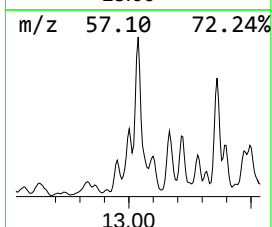
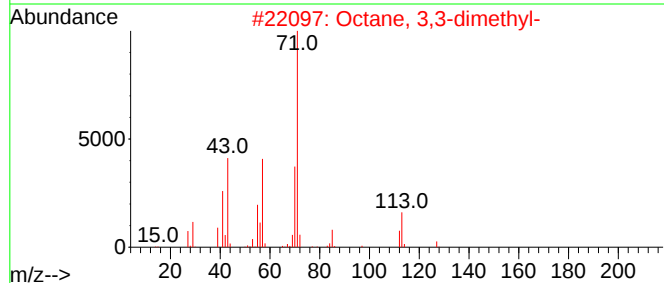
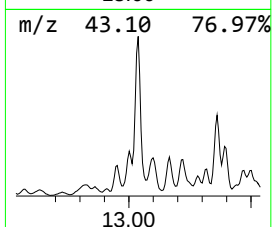
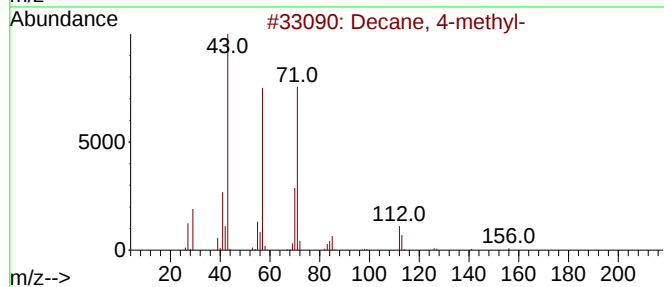
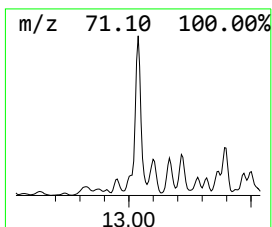
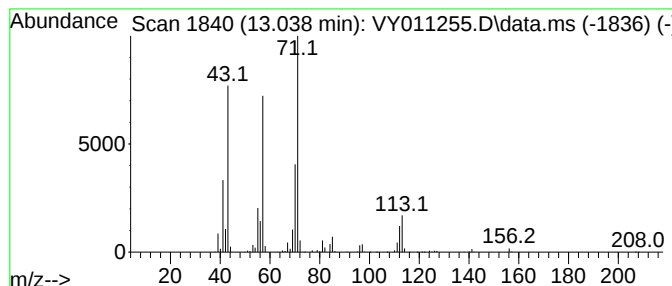
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y110122S.M
Quant Title : SW846 8260

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

Peak Number 1 Decane, 4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.038	195.01 ug/l	2044780	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 4-methyl-	156	C11H24	002847-72-5	83
2			Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	78
3			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	50
4			Sulfurous acid, isobutyl 2-penty...	208	C9H20O3S	1000309-15-2	50
5			Oxalic acid, ethyl neopentyl ester	188	C9H16O4	1000309-72-4	47



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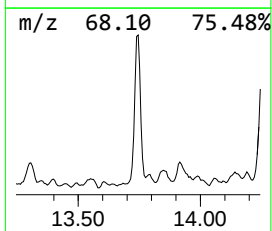
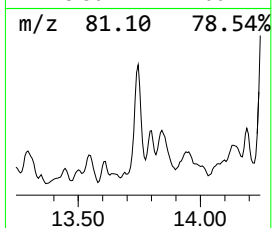
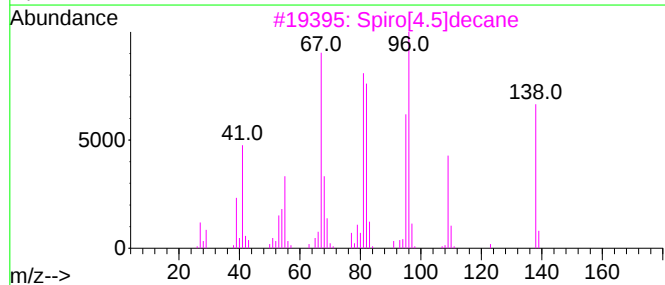
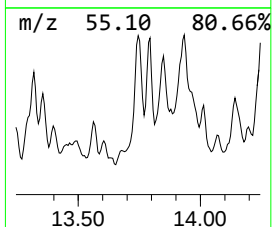
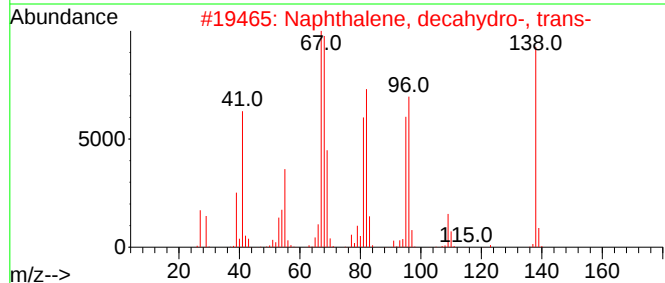
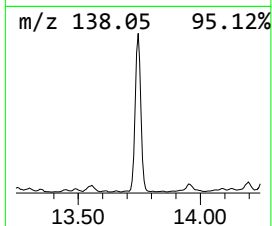
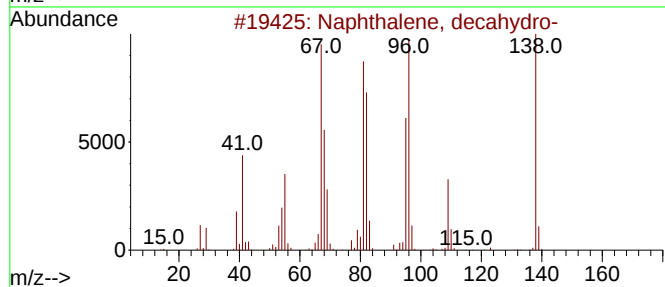
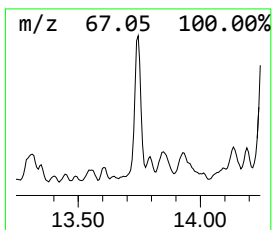
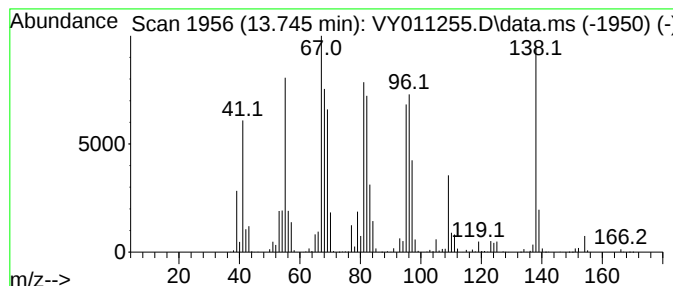
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y110122S.M
Quant Title : SW846 8260

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

Peak Number 2 Naphthalene, decahydro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.745	267.57 ug/l	2805710	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-	138	C10H18	000091-17-8	96
2			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	94
3			Spiro[4.5]decane	138	C10H18	000176-63-6	81
4			Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	76
5			Cyclodecene	138	C10H18	003618-12-0	64



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Instrument :
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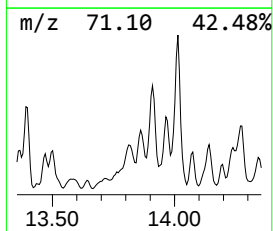
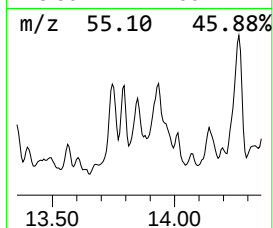
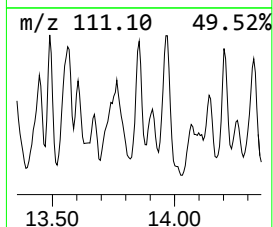
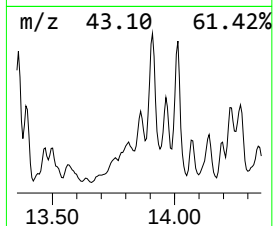
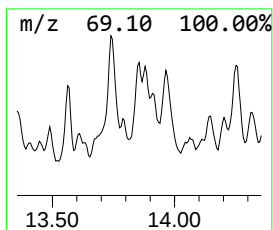
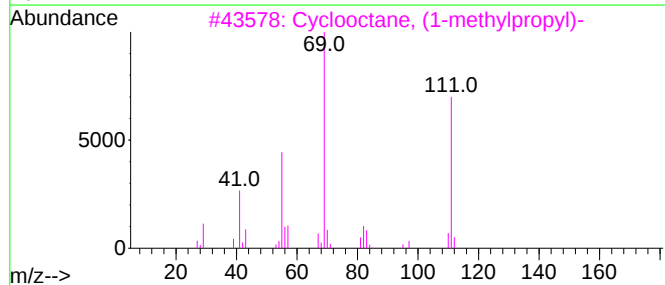
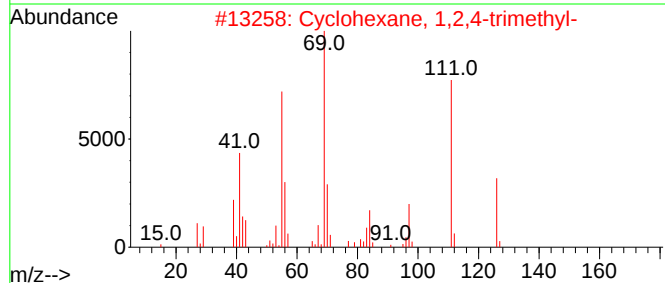
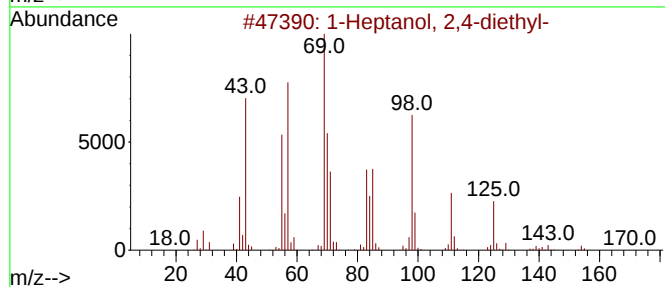
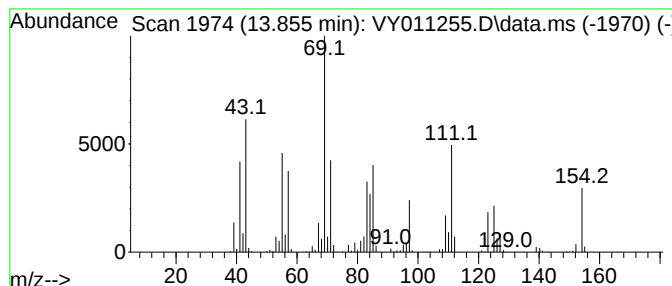
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 1-Heptanol, 2,4-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.855	154.86 ug/l	1623830	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heptanol, 2,4-diethyl-	172	C11H24O	080192-55-8	59
2			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	43
3			Cyclooctane, (1-methylpropyl)-	168	C12H24	016538-89-9	35
4			2-Hexene, 4-ethyl-2,3-dimethyl-	140	C10H20	1000149-19-7	30
5			3-Isopropyl-5-methyl-hex-4-en-2-one	154	C10H18O	077142-85-9	27



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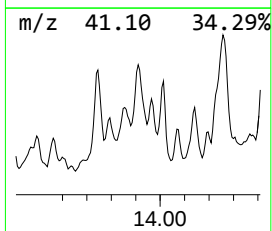
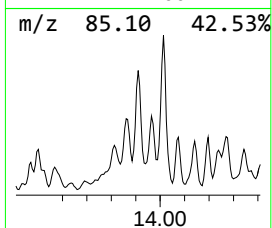
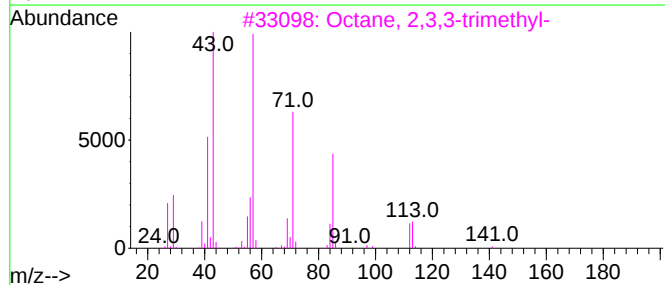
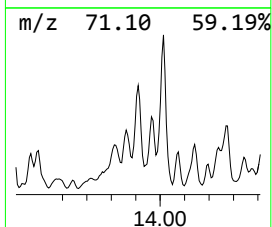
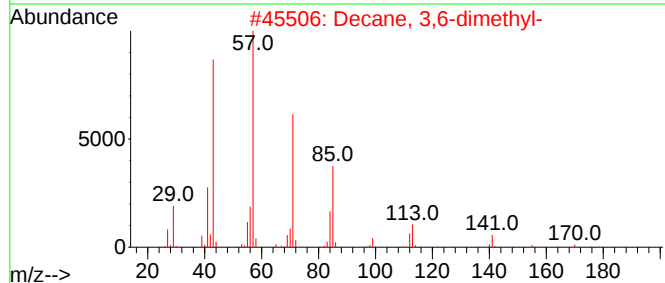
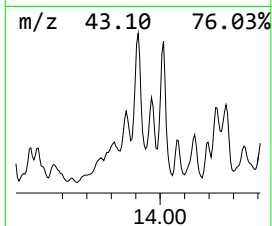
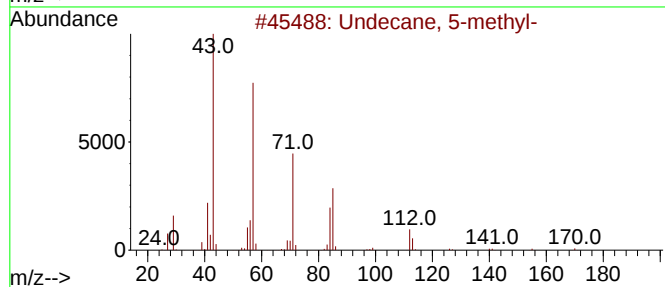
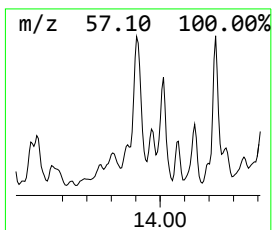
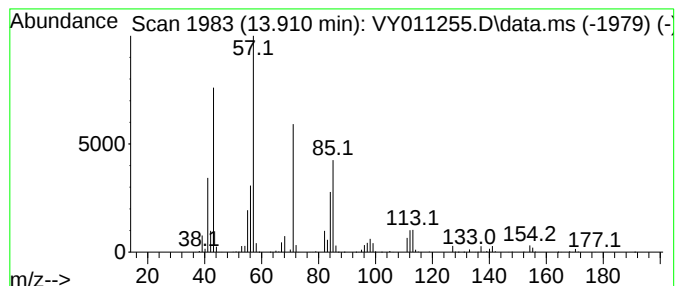
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TIC Integration Parameters: LSCINT.P

Peak Number 4 Undecane, 5-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.910	228.38 ug/l	2394750	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 5-methyl-	170	C12H26	001632-70-8	87
2			Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	81
3			Octane, 2,3,3-trimethyl-	156	C11H24	062016-30-2	72
4			2,6-Dimethyldecane	170	C12H26	013150-81-7	62
5			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110222\
Data File : VY011255.D
Acq On : 02 Nov 2022 12:25
Operator : KP/MD
Sample : N5337-03
Misc : 5.06g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-02-11.5-12.0

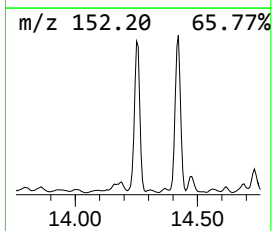
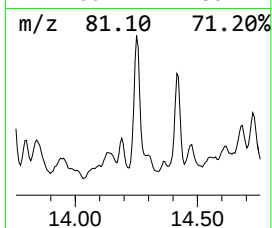
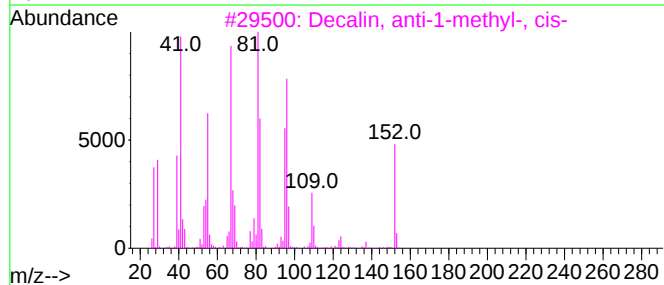
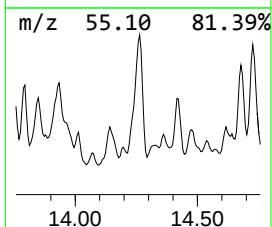
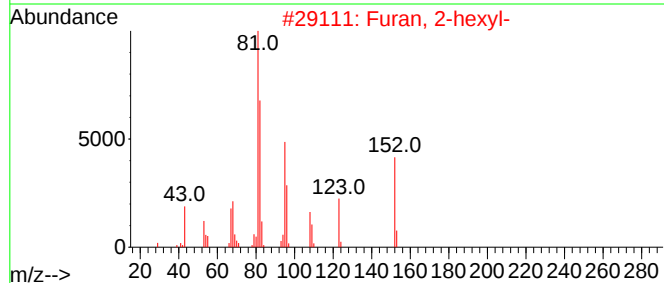
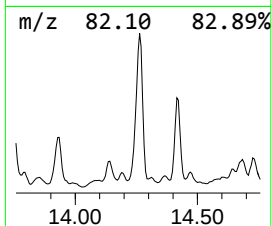
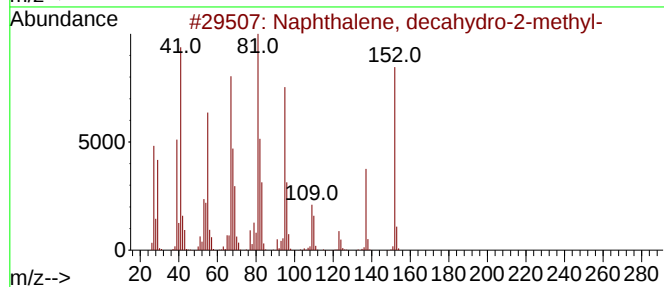
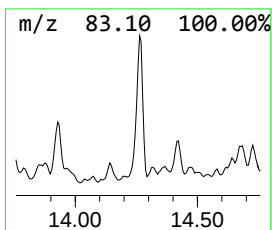
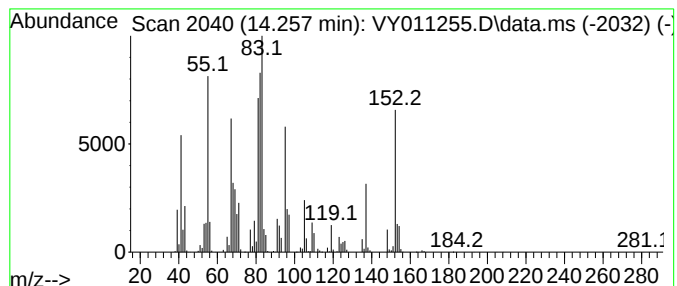
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Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, decahydro-2-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.257	526.40 ug/l	5519640	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	95
2			Furan, 2-hexyl-	152	C10H16O	003777-70-6	46
3			Decalin, anti-1-methyl-, cis-	152	C11H20	1000158-89-0	46
4			Cyclohexane, pentyl-	154	C11H22	004292-92-6	45
5			(Z)-(Z)-Hex-3-en-1-yl 2-methylbu...	182	C11H18O2	084060-80-0	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110222\
Data File : VY011255.D
Acq On : 02 Nov 2022 12:25
Operator : KP/MD
Sample : N5337-03
Misc : 5.06g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-02-11.5-12.0

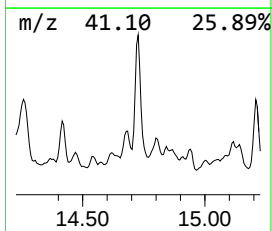
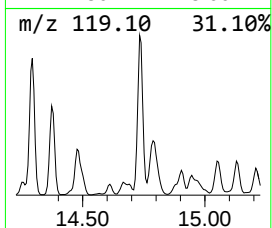
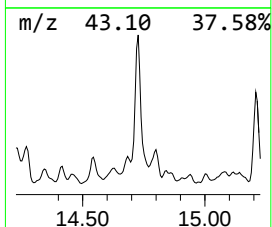
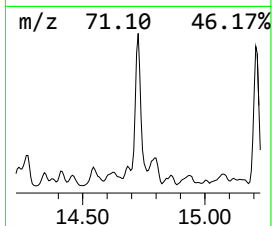
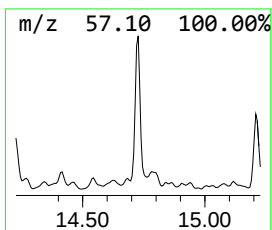
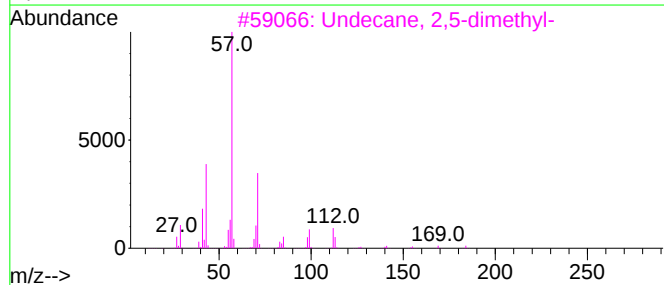
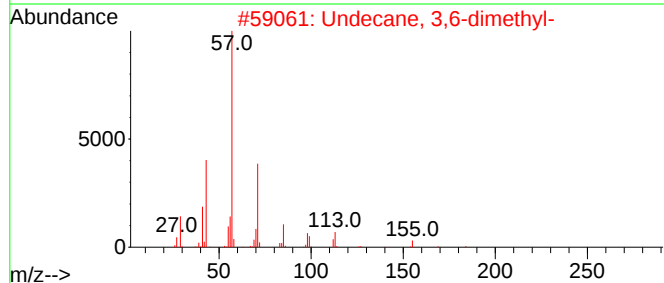
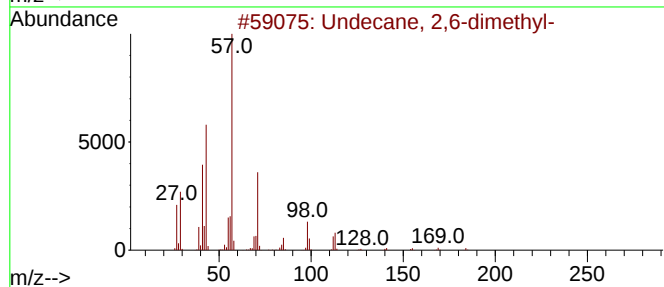
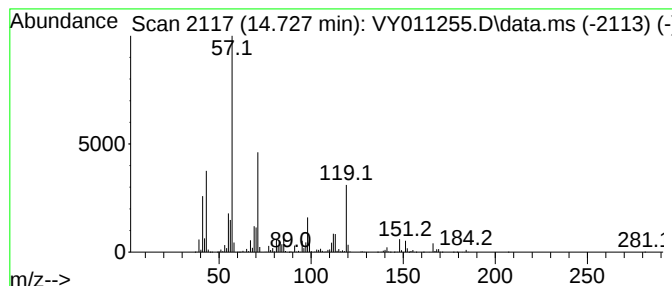
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.727	422.87 ug/l	4434110	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	83		
2	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	53		
3	Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	46		
4	Dodecane, 6-methyl-	184	C13H28	006044-71-9	46		
5	Sulfurous acid, butyl octyl ester	250	C12H26O3S	1000309-17-5	43		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110222\
Data File : VY011255.D
Acq On : 02 Nov 2022 12:25
Operator : KP/MD
Sample : N5337-03
Misc : 5.06g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-02-11.5-12.0

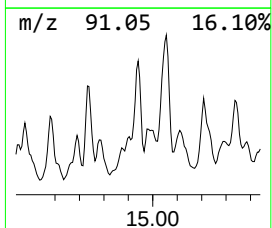
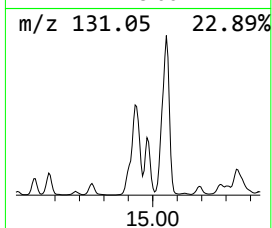
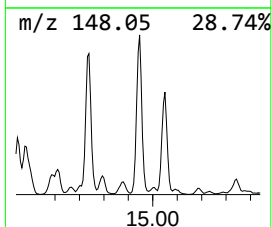
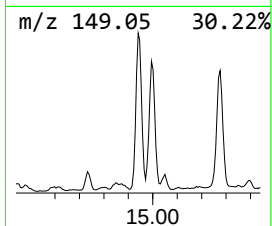
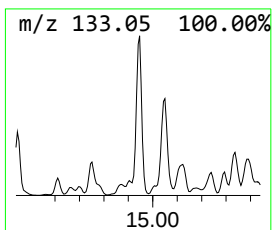
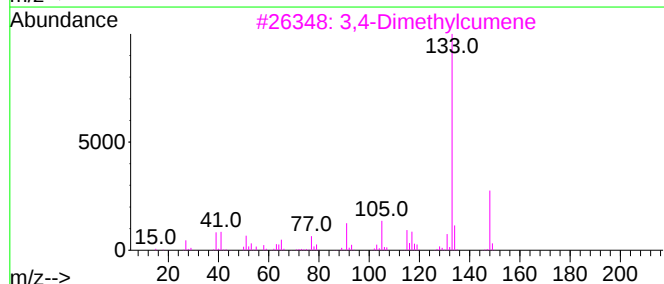
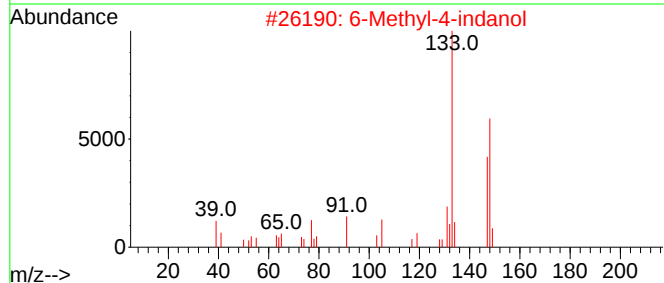
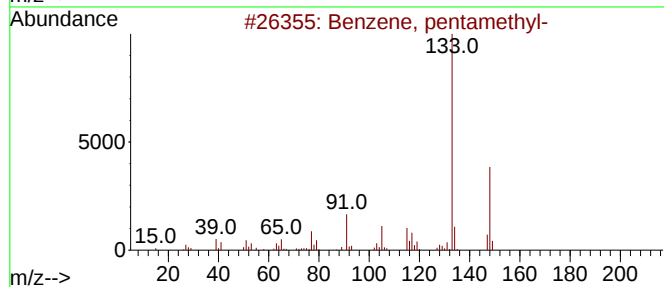
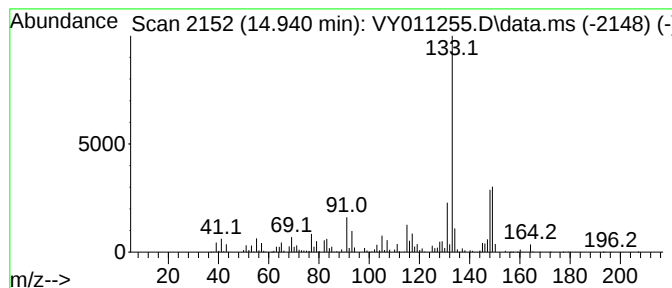
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, pentamethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.940	170.68 ug/l	1789700	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, pentamethyl-	148	C11H16	000700-12-9	91
2			6-Methyl-4-indanol	148	C10H12O	020294-32-0	64
3			3,4-Dimethylcumene	148	C11H16	1000370-34-1	58
4			4-tert-Butyltoluene	148	C11H16	000098-51-1	58
5			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110222\
Data File : VY011255.D
Acq On : 02 Nov 2022 12:25
Operator : KP/MD
Sample : N5337-03
Misc : 5.06g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-02-11.5-12.0

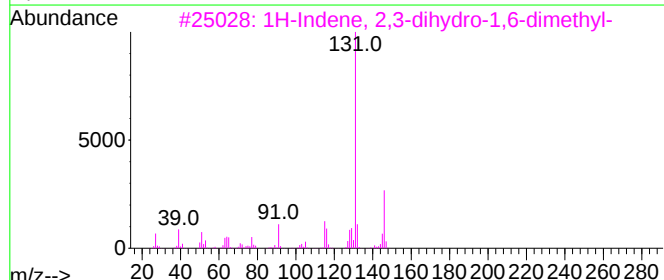
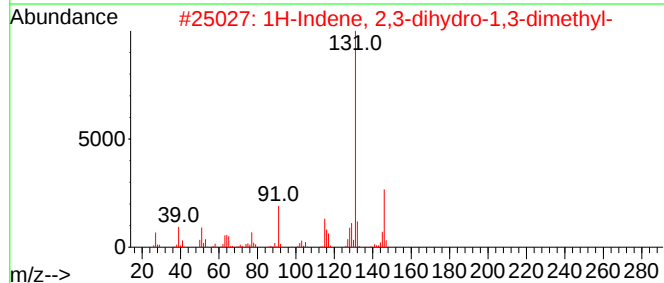
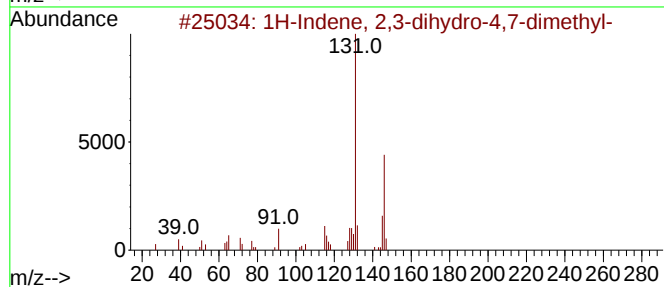
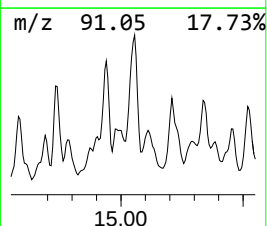
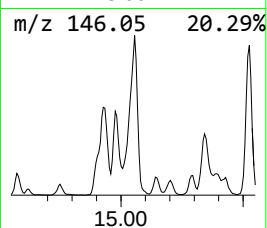
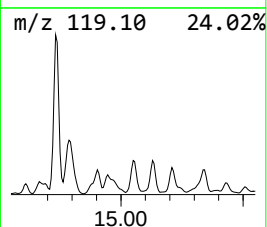
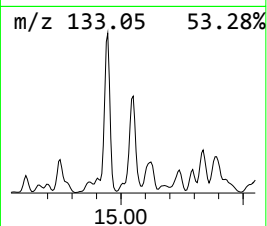
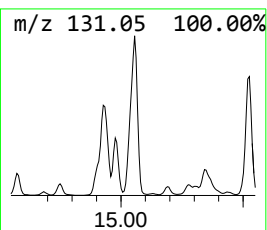
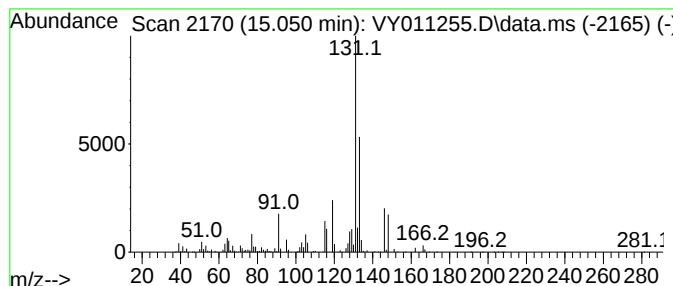
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Quant Title : SW846 8260

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

Peak Number 8 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.050	162.76 ug/l	1706640	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64
2			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	64
3			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	64
4			1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	64
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110222\
Data File : VY011255.D
Acq On : 02 Nov 2022 12:25
Operator : KP/MD
Sample : N5337-03
Misc : 5.06g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-02-11.5-12.0

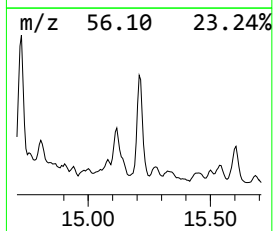
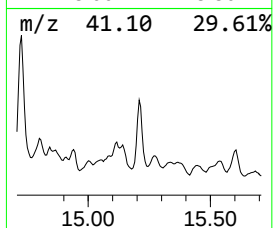
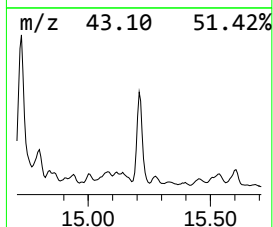
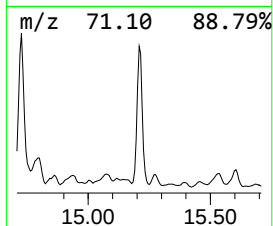
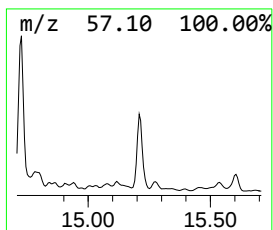
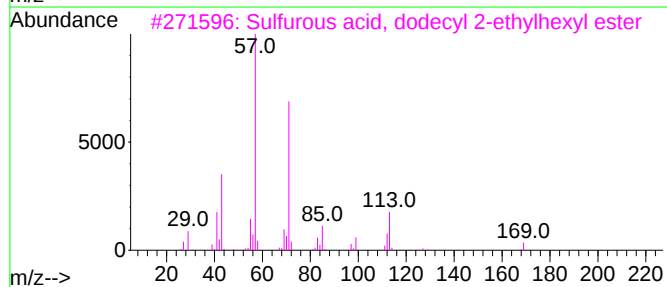
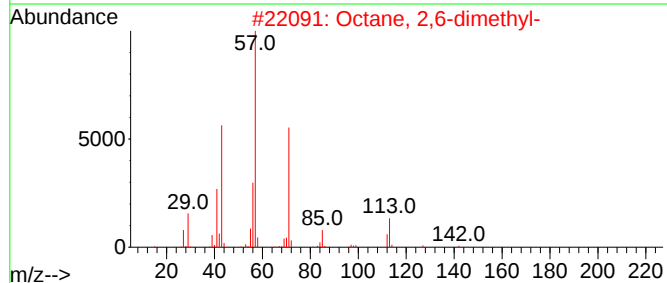
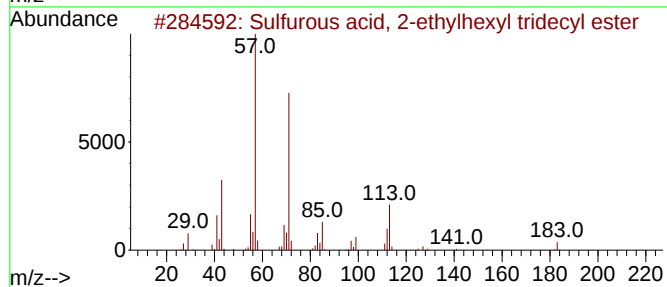
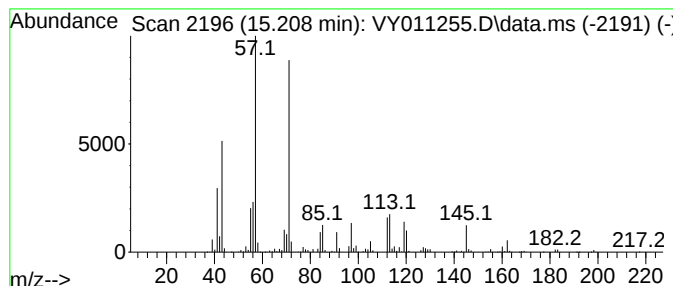
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Quant Title : SW846 8260

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

Peak Number 9 Sulfurous acid, 2-ethylhexy... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.208	241.85 ug/l	2535990	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, 2-ethylhexyl tri...	376	C21H44O3S	1000309-19-6	64
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
3			Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	64
4			Tridecane, 7-methyl-	198	C14H30	026730-14-3	53
5			Decane, 2-methyl-	156	C11H24	006975-98-0	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110222\
Data File : VY011255.D
Acq On : 02 Nov 2022 12:25
Operator : KP/MD
Sample : N5337-03
Misc : 5.06g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-02-11.5-12.0

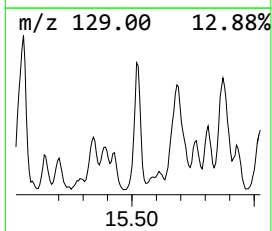
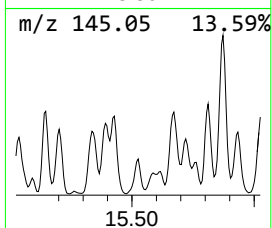
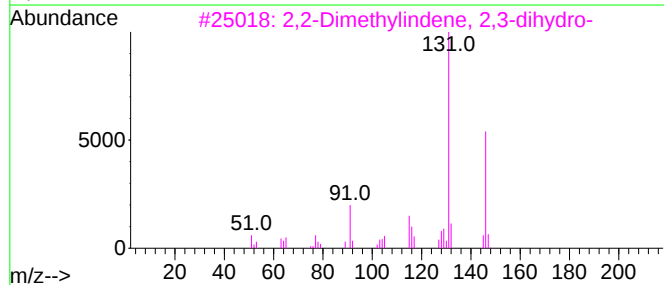
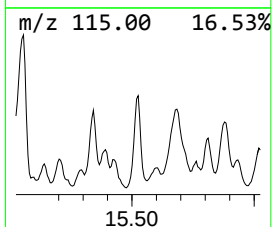
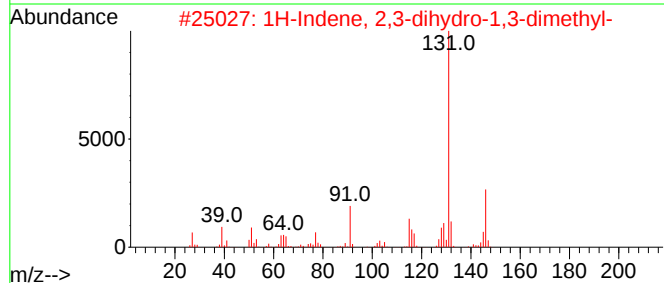
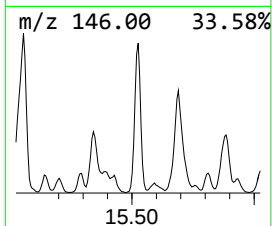
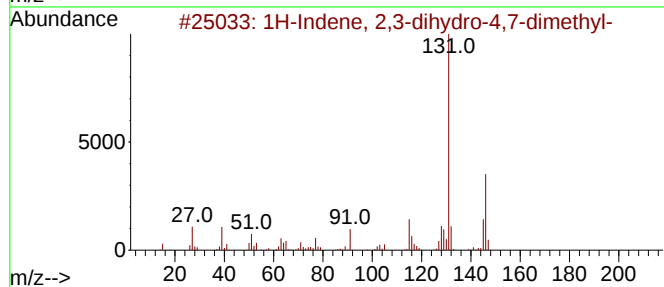
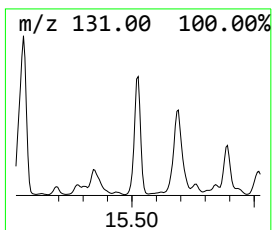
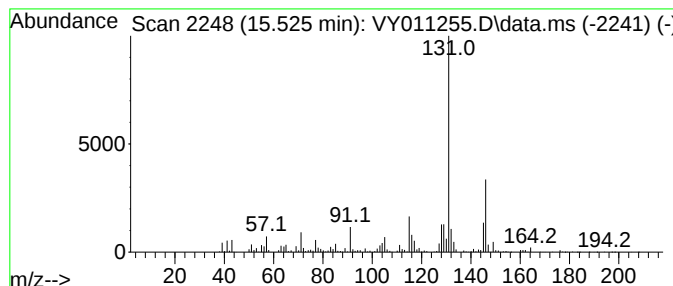
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Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.525	186.37 ug/l	1954200	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94		
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	93		
3	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91		
4	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	91		
5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY110222\
Data File : VY011255.D
Acq On : 02 Nov 2022 12:25
Operator : KP/MD
Sample : N5337-03
Misc : 5.06g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
SB-02-11.5-12.0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y110122S.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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Naphthalene, de...	13.745	267.6	ug/l	2805710	4	13.422	524286	50.0
1-Heptanol, 2,4...	13.855	154.9	ug/l	1623830	4	13.422	524286	50.0
Undecane, 5-met...	13.910	228.4	ug/l	2394750	4	13.422	524286	50.0
Naphthalene, de...	14.257	526.4	ug/l	5519640	4	13.422	524286	50.0
Undecane, 2,6-d...	14.727	422.9	ug/l	4434110	4	13.422	524286	50.0
Benzene, pentam...	14.940	170.7	ug/l	1789700	4	13.422	524286	50.0
1H-Indene, 2,3-...	15.050	162.8	ug/l	1706640	4	13.422	524286	50.0
Sulfurous acid,...	15.208	241.8	ug/l	2535990	4	13.422	524286	50.0
1H-Indene, 2,3-...	15.525	186.4	ug/l	1954200	4	13.422	524286	50.0