

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY022322\
 Data File : VY007606.D
 Acq On : 23 Feb 2022 18:29
 Operator : SY/MD
 Sample : N1584-11
 Misc : 5.12g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 D15-6

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\82Y022322S.M
 Title : SW846 8260

Signal : TIC: VY007606.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.899	167	177	189	rBV3	31341	72773	6.18%	0.791%
2	3.253	227	235	248	rBV3	12535	33263	2.82%	0.361%
3	3.698	300	308	318	rVB3	8605	21432	1.82%	0.233%
4	3.942	336	348	360	rBV3	10662	32501	2.76%	0.353%
5	4.704	459	473	478	rBV4	28584	101498	8.62%	1.103%
6	5.234	548	560	577	rVB2	31693	116143	9.86%	1.262%
7	5.551	602	612	624	rVB4	3851	13389	1.14%	0.145%
8	5.753	633	645	656	rBV2	11039	34695	2.95%	0.377%
9	6.100	691	702	715	rVB	20633	62721	5.33%	0.681%
10	6.240	715	725	735	rBV2	11065	34638	2.94%	0.376%
11	6.539	763	774	783	rBV5	14009	42398	3.60%	0.461%
12	6.691	788	799	806	rVV3	13892	43158	3.66%	0.469%
13	6.795	806	816	825	rVV2	49716	155850	13.23%	1.693%
14	6.868	825	828	837	rVB3	6288	14850	1.26%	0.161%
15	7.527	928	936	946	rVB2	37973	97025	8.24%	1.054%
16	7.716	957	967	972	rBV	72452	187387	15.91%	2.036%
17	7.789	972	979	986	rVV2	146961	366998	31.16%	3.987%
18	7.868	986	992	1003	rVB3	47781	124182	10.54%	1.349%
19	7.996	1003	1013	1019	rBV2	25322	59190	5.03%	0.643%
20	8.142	1028	1037	1048	rBV2	76480	195404	16.59%	2.123%
21	8.325	1048	1067	1077	rVV3	64686	278039	23.61%	3.020%
22	8.417	1077	1082	1092	rVV3	13521	35405	3.01%	0.385%
23	8.545	1092	1103	1110	rVB6	13500	35705	3.03%	0.388%
24	8.685	1118	1126	1138	rBV	222815	496490	42.16%	5.393%
25	8.783	1138	1142	1148	rVV4	6840	14383	1.22%	0.156%
26	8.850	1148	1153	1158	rVV2	7571	13564	1.15%	0.147%
27	8.923	1159	1165	1168	rVV2	13161	26366	2.24%	0.286%
28	8.965	1168	1172	1185	rVB	19120	40518	3.44%	0.440%
29	9.148	1194	1202	1203	rBV3	22856	39162	3.33%	0.425%
30	9.185	1203	1208	1213	rVV5	46444	99599	8.46%	1.082%
31	9.240	1213	1217	1224	rVB3	12520	24541	2.08%	0.267%
32	9.362	1233	1237	1245	rVB3	10133	19696	1.67%	0.214%
33	9.459	1245	1253	1259	rBV6	7582	18311	1.55%	0.199%
34	9.612	1272	1278	1287	rVB	51864	100629	8.54%	1.093%
35	9.709	1287	1294	1296	rBV2	33438	63291	5.37%	0.688%
36	9.746	1296	1300	1308	rVV3	68335	155330	13.19%	1.687%

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37	9.819	1308	1312	1315	rVV4	9775	17959	1.52%	0.195%
38	9.849	1315	1317	1325	rVB5	8733	18777	1.59%	0.204%
39	9.959	1325	1335	1345	rBV8	13737	50899	4.32%	0.553%
40	10.057	1345	1351	1355	rBV2	20416	39277	3.33%	0.427%
41	10.112	1355	1360	1364	rVV2	17510	28800	2.45%	0.313%
42	10.179	1364	1371	1378	rVV	288515	516197	43.83%	5.607%
43	10.240	1378	1381	1385	rVV	14909	24231	2.06%	0.263%
44	10.294	1385	1390	1394	rVV3	9087	17064	1.45%	0.185%
45	10.374	1394	1403	1409	rVB7	14897	37834	3.21%	0.411%
46	10.599	1434	1440	1450	rVB3	28480	68982	5.86%	0.749%
47	10.697	1450	1456	1464	rVB9	5607	13134	1.12%	0.143%
48	10.794	1464	1472	1476	rBV8	4604	11852	1.01%	0.129%
49	10.874	1476	1485	1487	rBV5	6283	14575	1.24%	0.158%
50	11.288	1547	1553	1562	rVB9	8661	23850	2.03%	0.259%
51	11.374	1562	1567	1576	rBV6	9136	20945	1.78%	0.228%
52	11.489	1579	1586	1592	rVV	302872	487747	41.41%	5.298%
53	11.587	1594	1602	1611	rVV	208922	339261	28.81%	3.685%
54	11.703	1613	1621	1631	rVB	53556	104504	8.87%	1.135%
55	12.026	1664	1674	1682	rVB4	8293	21771	1.85%	0.236%
56	12.325	1717	1723	1729	rBV2	32008	53274	4.52%	0.579%
57	12.483	1741	1749	1758	rVB3	639105	1177732	100.00%	12.793%
58	12.666	1770	1779	1785	rVB	42810	69803	5.93%	0.758%
59	12.733	1785	1790	1792	rVV2	12325	18746	1.59%	0.204%
60	12.764	1792	1795	1799	rVV	19088	29163	2.48%	0.317%
61	12.806	1799	1802	1807	rVB2	14758	22054	1.87%	0.240%
62	12.971	1823	1829	1837	rVB2	29534	48271	4.10%	0.524%
63	13.117	1847	1853	1859	rVB	29784	46505	3.95%	0.505%
64	13.245	1866	1874	1879	rBV3	13972	32348	2.75%	0.351%
65	13.422	1897	1903	1913	rBV	304426	501478	42.58%	5.447%
66	13.587	1924	1930	1933	rBV	98093	159954	13.58%	1.738%
67	13.623	1933	1936	1941	rVV	281836	419450	35.62%	4.556%
68	13.672	1941	1944	1952	rVB3	42498	77608	6.59%	0.843%
69	13.751	1952	1957	1962	rBV3	17448	27536	2.34%	0.299%
70	13.818	1962	1968	1974	rVB2	13228	21830	1.85%	0.237%
71	13.891	1974	1980	1986	rBV3	19449	40767	3.46%	0.443%
72	13.965	1986	1992	1998	rVB3	215385	343599	29.17%	3.732%
73	14.026	1998	2002	2006	rBV	21288	33110	2.81%	0.360%
74	14.080	2006	2011	2017	rVB	96149	152531	12.95%	1.657%
75	14.294	2037	2046	2050	rBV	67262	106421	9.04%	1.156%
76	14.373	2056	2059	2062	rBV2	8957	12573	1.07%	0.137%

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77	14.562	2085	2090	2095	rBV	61675	94008	7.98%	1.021%
78	14.684	2102	2110	2115	rBV2	110999	206746	17.55%	2.246%
79	14.739	2115	2119	2125	rVB3	10582	19653	1.67%	0.213%
80	14.830	2129	2134	2139	rVB2	15563	23029	1.96%	0.250%
81	14.940	2146	2152	2156	rBV3	18645	31216	2.65%	0.339%
82	15.050	2162	2170	2178	rVB3	18893	45005	3.82%	0.489%
83	15.233	2190	2200	2207	rVB	63466	108465	9.21%	1.178%
84	15.342	2213	2218	2222	rBV4	7318	13008	1.10%	0.141%
85	15.391	2222	2226	2241	rVB4	7163	16420	1.39%	0.178%
86	15.525	2241	2248	2255	rBV2	10005	19512	1.66%	0.212%
87	15.696	2267	2276	2287	rBV7	5771	18623	1.58%	0.202%
88	15.885	2300	2307	2313	rVB6	4595	11982	1.02%	0.130%
89	16.287	2367	2373	2380	rVB2	21972	44222	3.75%	0.480%
90	16.488	2398	2406	2413	rBV3	14821	30932	2.63%	0.336%

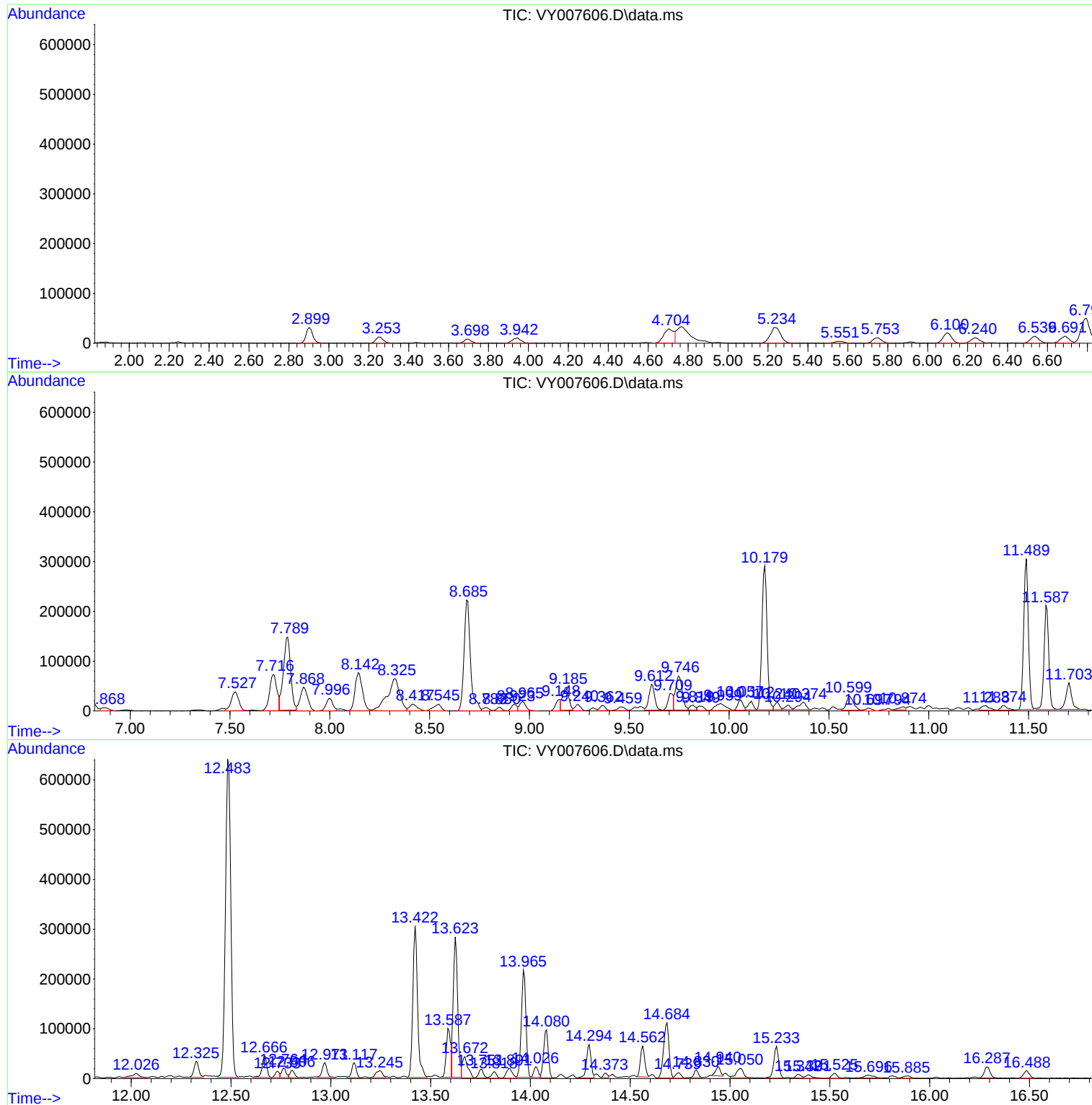
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y022322S.M
Quant Title : SW846 8260

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



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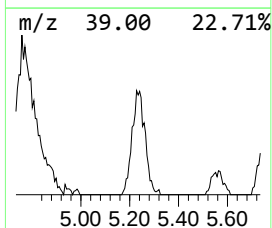
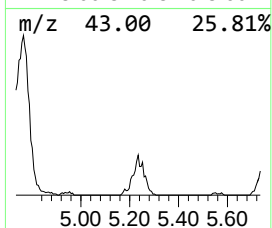
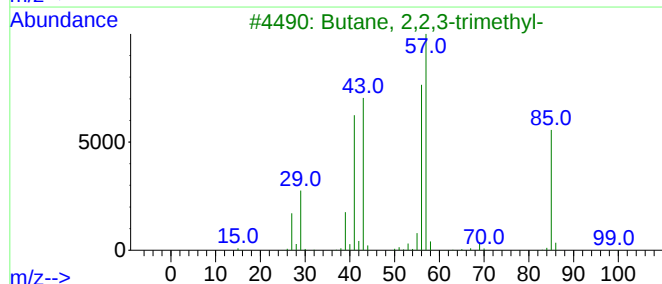
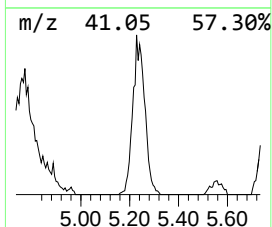
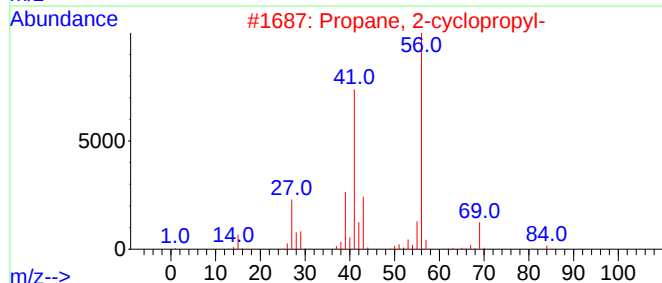
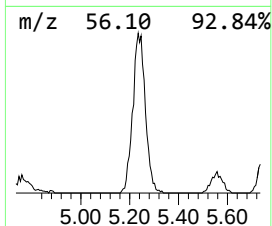
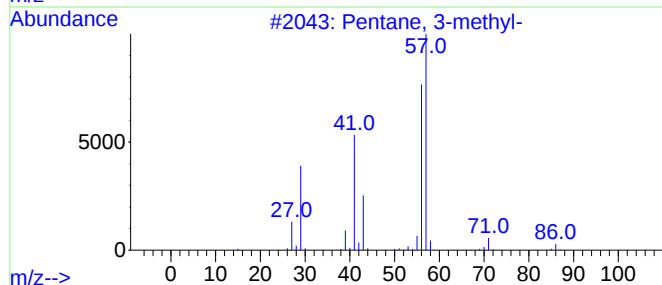
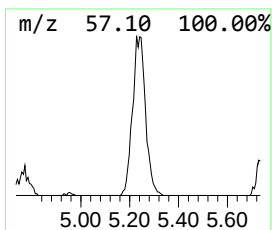
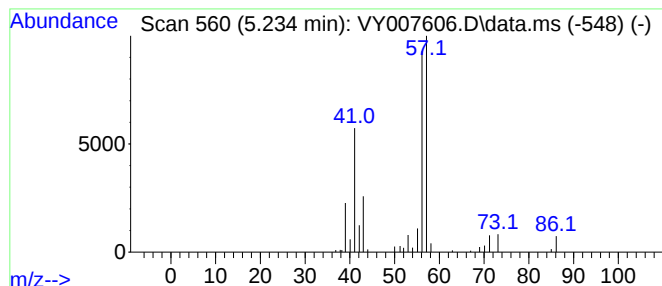
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y022322S.M
Quant Title : SW846 8260

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

Peak Number 1 Pentane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.234	15.82 ug/l	116143	Pentafluorobenzene	7.789

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-methyl-	86	C6H14	000096-14-0	72
2			Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	45
3			Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	40
4			Cyclobutanecarbonitrile, 3,3-dim...	109	C7H11N	053783-86-1	40
5			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	39



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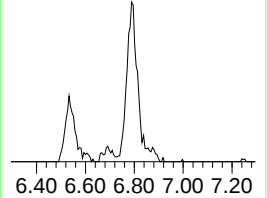
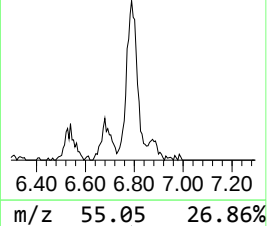
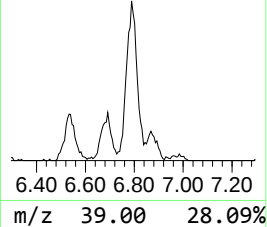
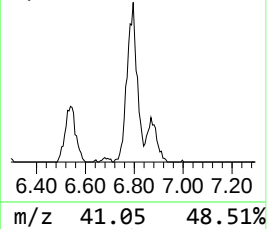
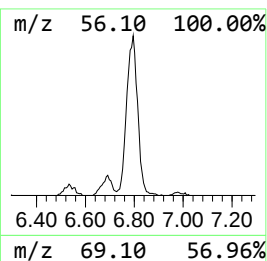
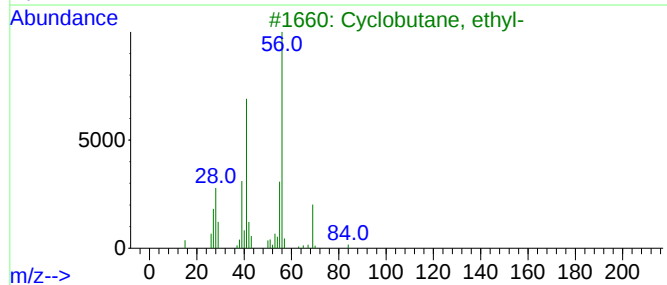
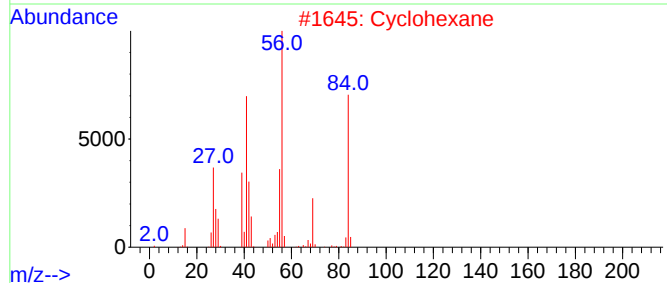
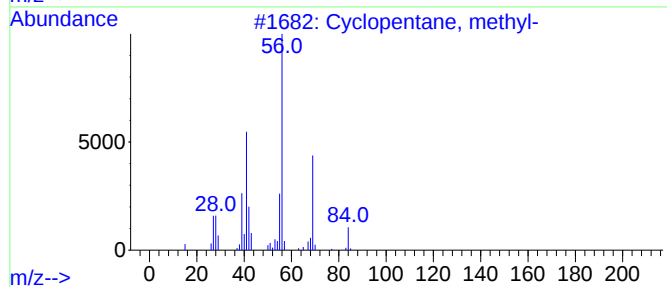
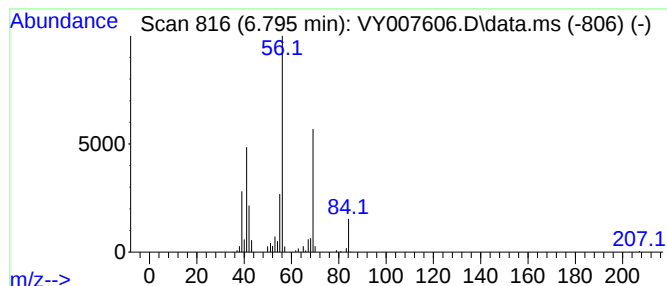
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y022322S.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Cyclopentane, methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.795	21.23 ug/l	155850	Pentafluorobenzene	7.789

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2			Cyclohexane	84	C6H12	000110-82-7	86
3			Cyclobutane, ethyl-	84	C6H12	004806-61-5	59
4			1-Hexene, 2-methyl-	98	C7H14	006094-02-6	50
5			1-Heptene, 2-methyl-	112	C8H16	015870-10-7	50



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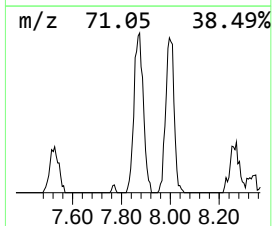
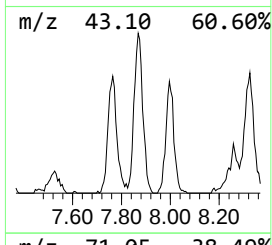
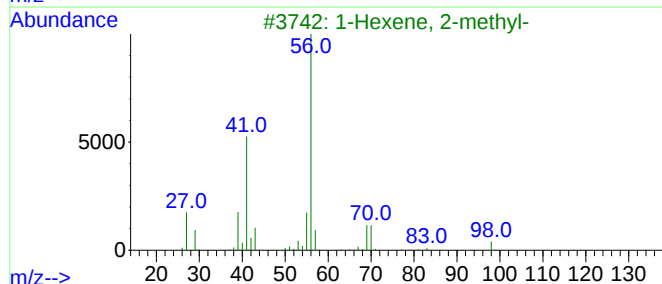
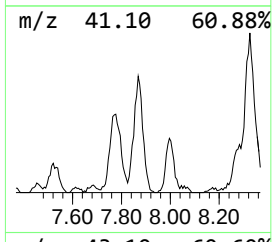
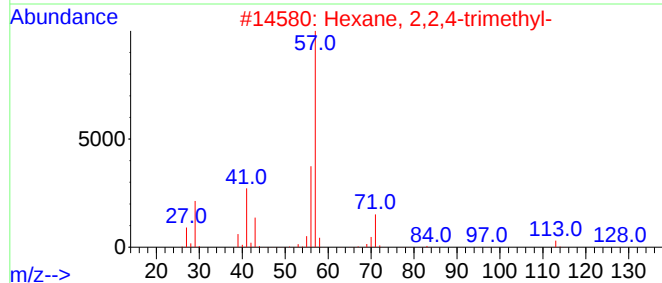
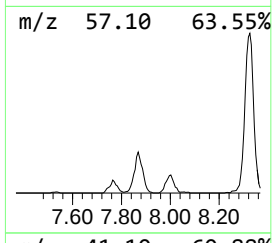
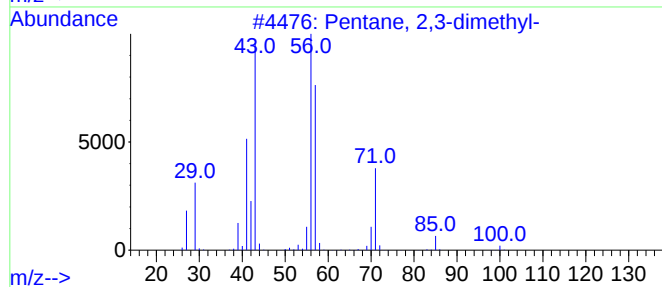
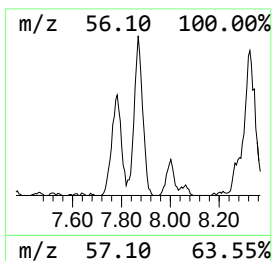
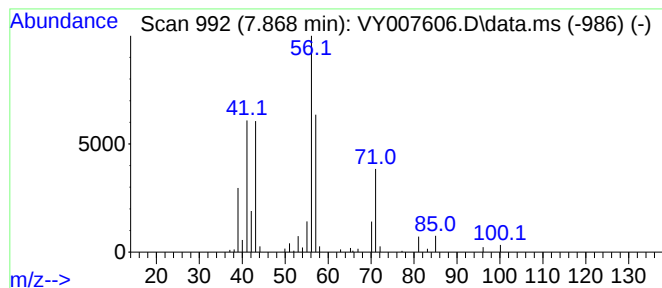
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Quant Title : SW846 8260

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

Peak Number 3 Pentane, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.868	16.92 ug/l	124182	Pentafluorobenzene	7.789

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	83
2			Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	56
3			1-Hexene, 2-methyl-	98	C7H14	006094-02-6	47
4			Heptane	100	C7H16	000142-82-5	37
5			Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	37



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D15-6

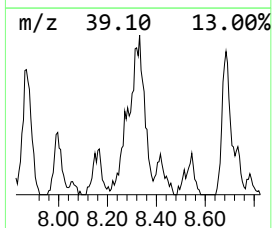
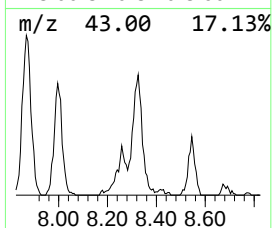
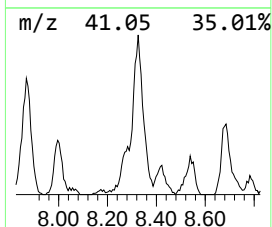
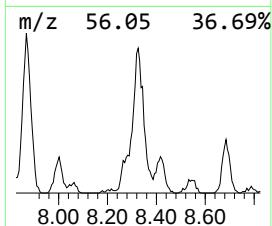
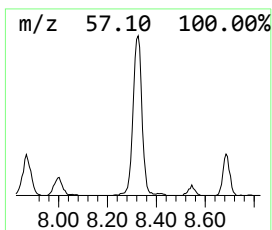
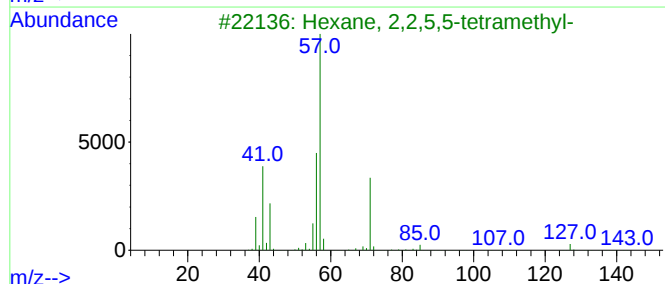
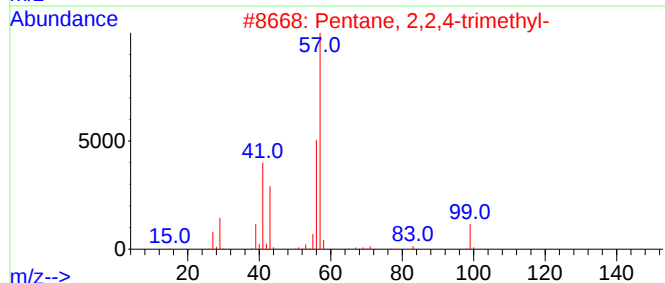
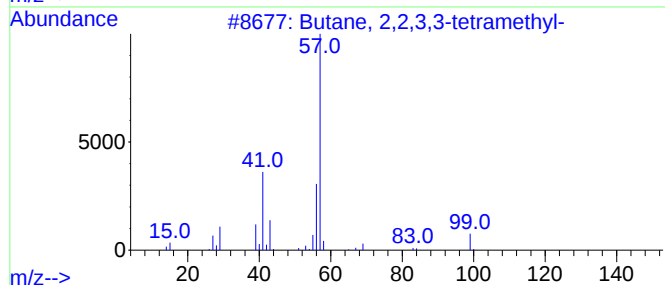
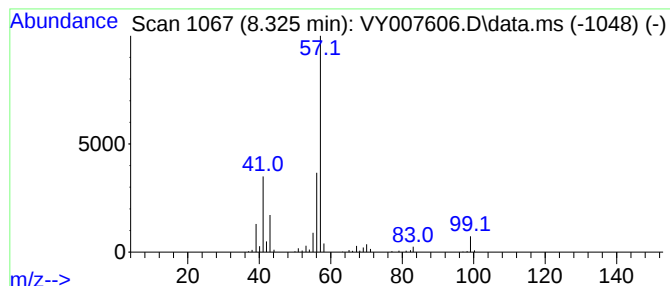
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Quant Title : SW846 8260

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

Peak Number 4 Butane, 2,2,3,3-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.325	28.00 ug/l	278039	1,4-Difluorobenzene	8.685

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
2			Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	64
3			Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	59
4			Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	53
5			Heptane, 2,2,3,5-tetramethyl-	156	C11H24	061868-42-6	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY022322\
Data File : VY007606.D
Acq On : 23 Feb 2022 18:29
Operator : SY/MD
Sample : N1584-11
Misc : 5.12g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
D15-6

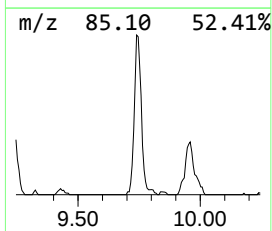
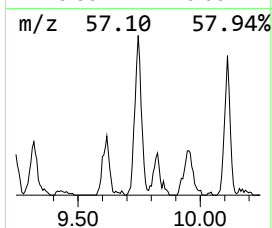
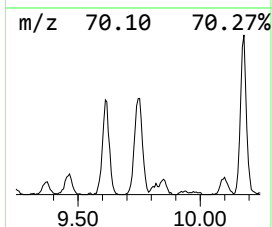
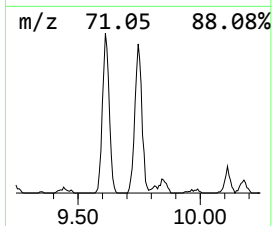
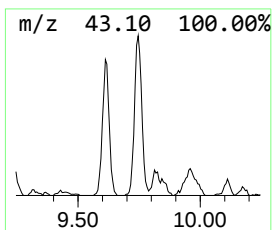
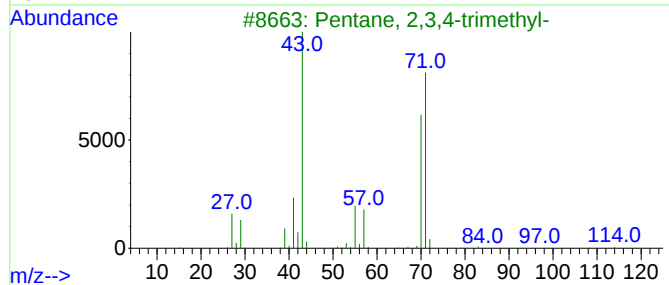
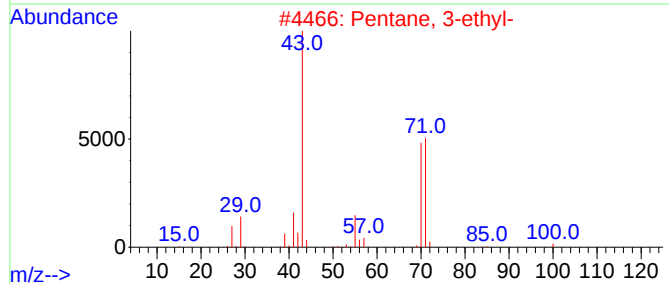
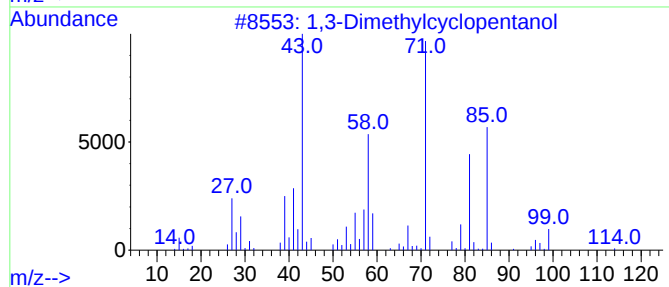
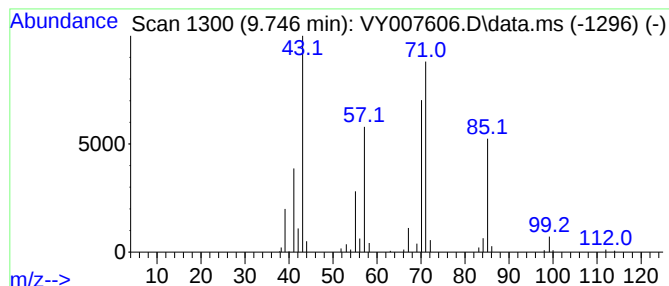
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Quant Title : SW846 8260

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

Peak Number 5 1,3-Dimethylcyclopentanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.746	15.64 ug/l	155330	1,4-Difluorobenzene	8.685

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dimethylcyclopentanol	114	C7H14O	019550-46-0	64
2			Pentane, 3-ethyl-	100	C7H16	000617-78-7	62
3			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	58
4			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	53
5			Heptane, 2,5,5-trimethyl-	142	C10H22	001189-99-7	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY022322\
Data File : VY007606.D
Acq On : 23 Feb 2022 18:29
Operator : SY/MD
Sample : N1584-11
Misc : 5.12g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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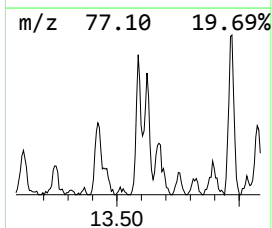
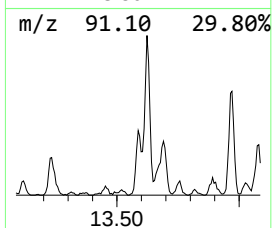
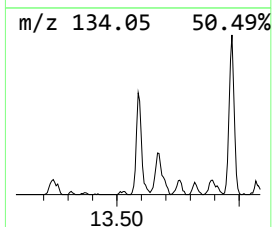
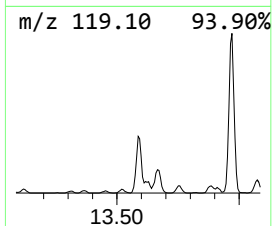
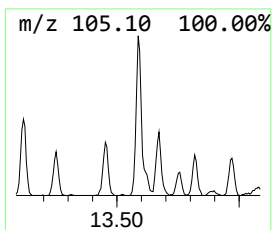
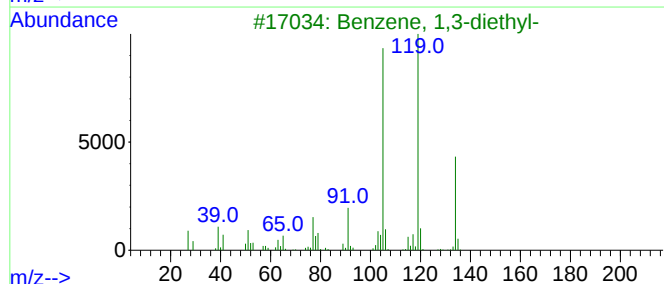
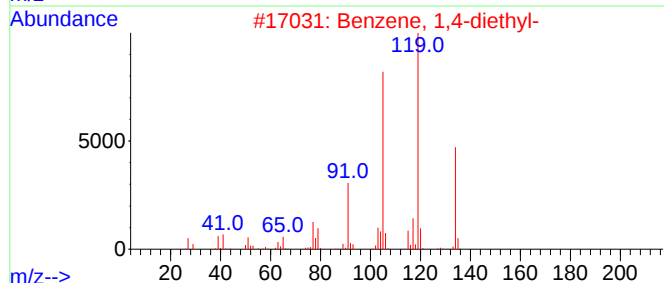
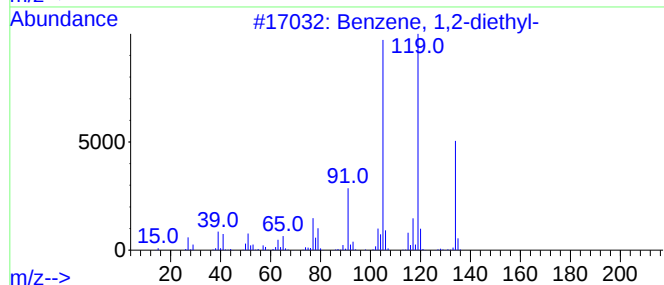
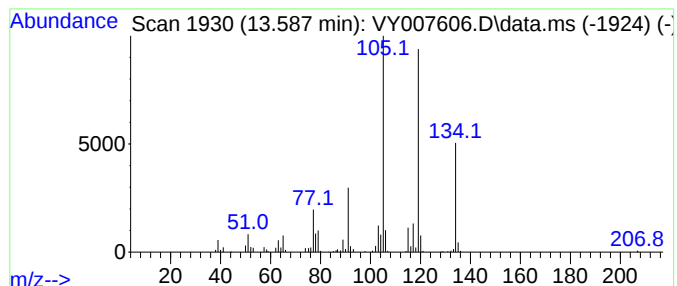
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.587	15.95 ug/l	159954	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
2			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	91
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	80
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY022322\
Data File : VY007606.D
Acq On : 23 Feb 2022 18:29
Operator : SY/MD
Sample : N1584-11
Misc : 5.12g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
D15-6

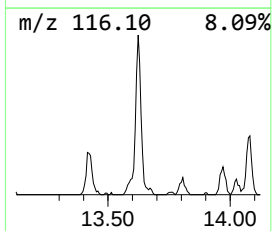
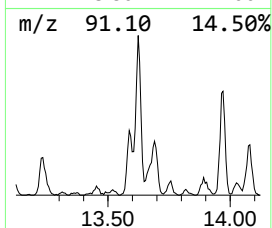
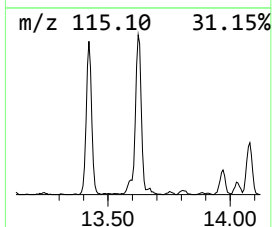
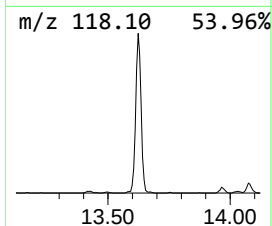
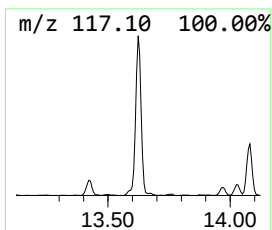
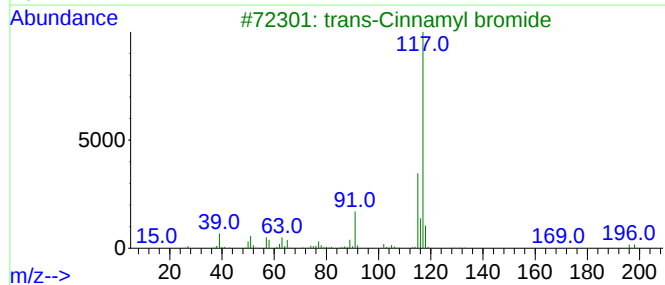
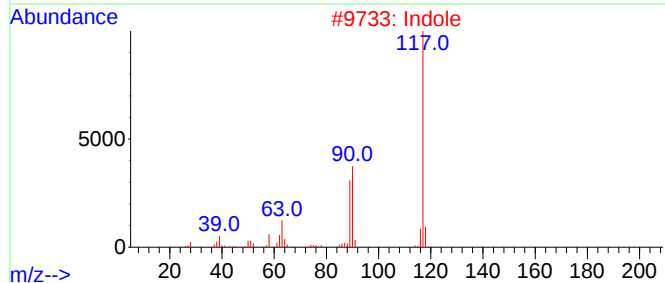
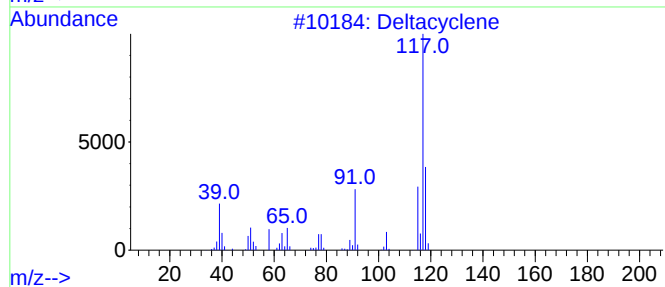
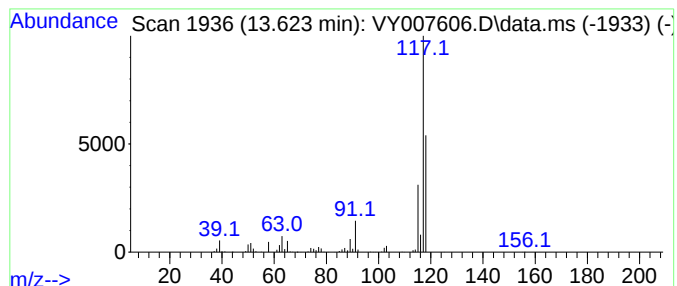
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Quant Title : SW846 8260

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

Peak Number 7 Deltacyclene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.623	41.82 ug/l	419450	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Deltacyclene	118	C9H10	007785-10-6	59
2			Indole	117	C8H7N	000120-72-9	46
3			trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	45
4			Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	42
5			Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	40



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY022322\
Data File : VY007606.D
Acq On : 23 Feb 2022 18:29
Operator : SY/MD
Sample : N1584-11
Misc : 5.12g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
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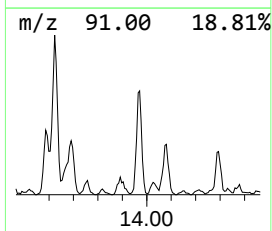
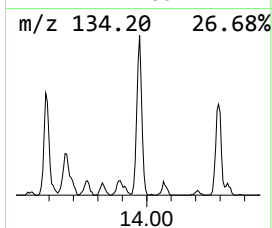
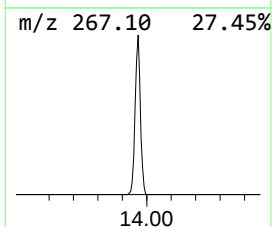
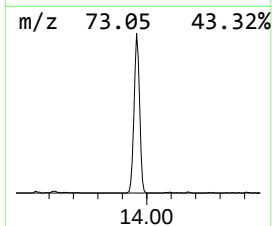
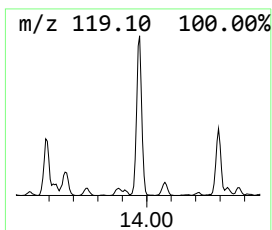
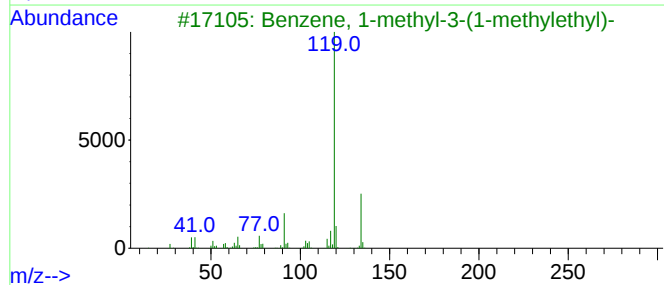
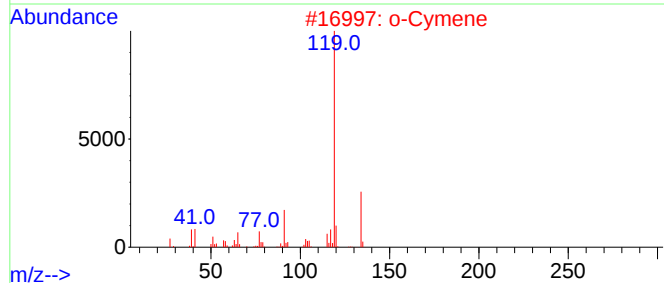
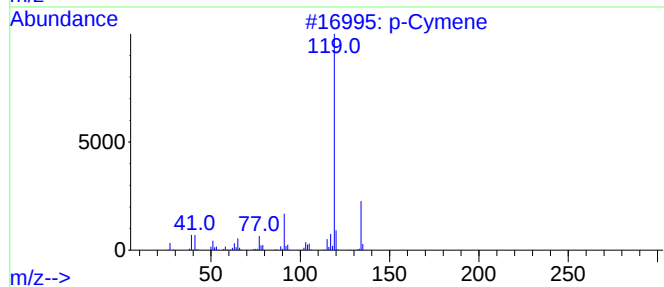
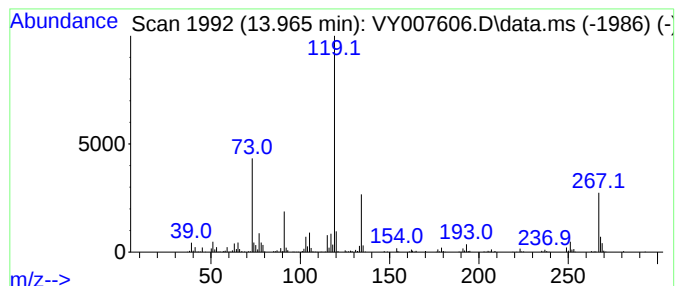
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Quant Title : SW846 8260

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

Peak Number 8 p-Cymene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.964	34.26 ug/l	343599	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	93
2			o-Cymene	134	C10H14	000527-84-4	93
3			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	92
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	92
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY022322\
Data File : VY007606.D
Acq On : 23 Feb 2022 18:29
Operator : SY/MD
Sample : N1584-11
Misc : 5.12g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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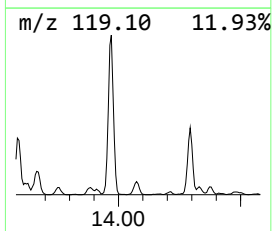
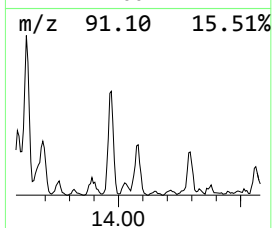
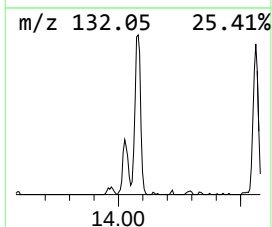
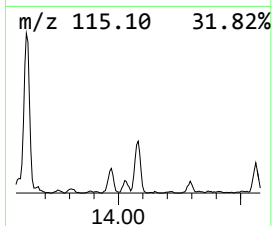
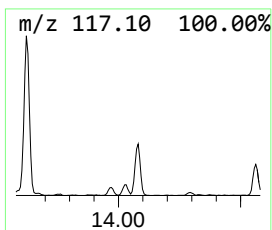
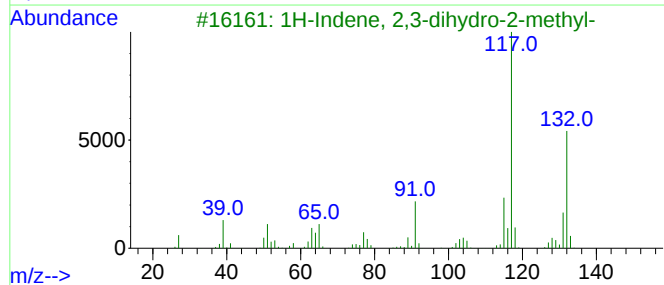
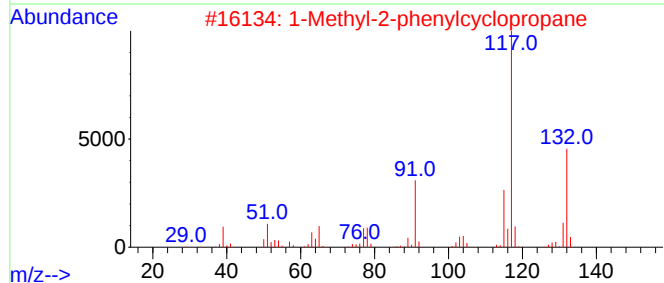
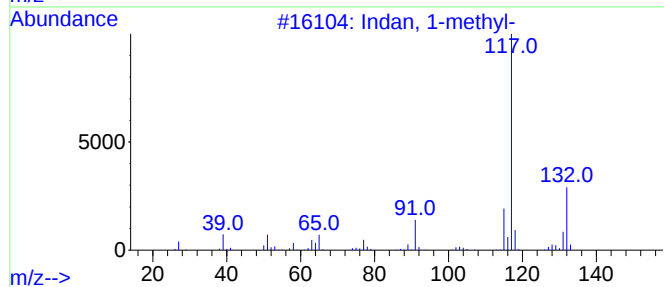
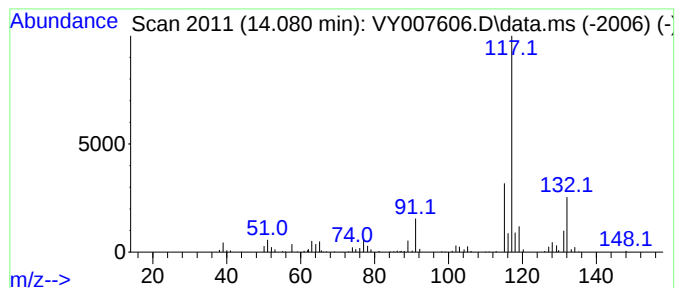
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Quant Title : SW846 8260

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

Peak Number 9 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.080	15.21 ug/l	152531	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	87
2			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	83
3			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	83
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	83
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY022322\
Data File : VY007606.D
Acq On : 23 Feb 2022 18:29
Operator : SY/MD
Sample : N1584-11
Misc : 5.12g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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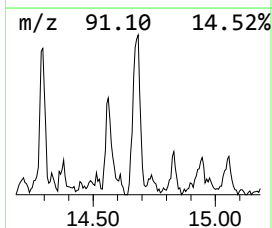
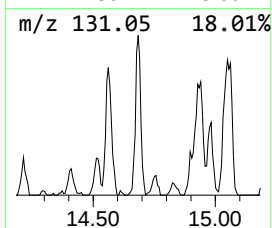
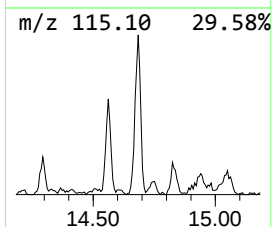
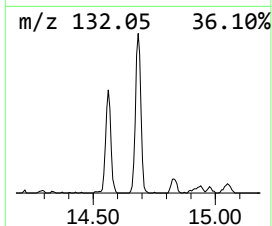
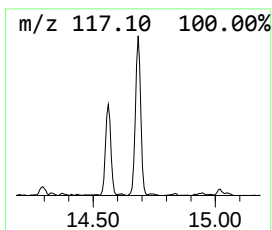
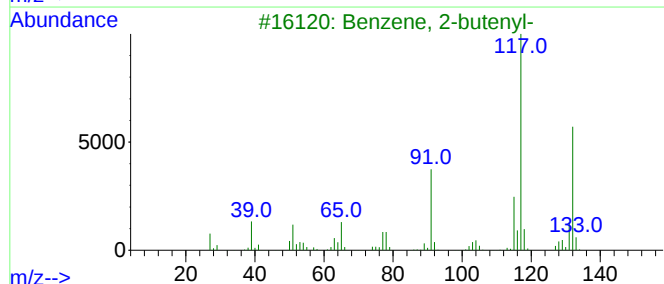
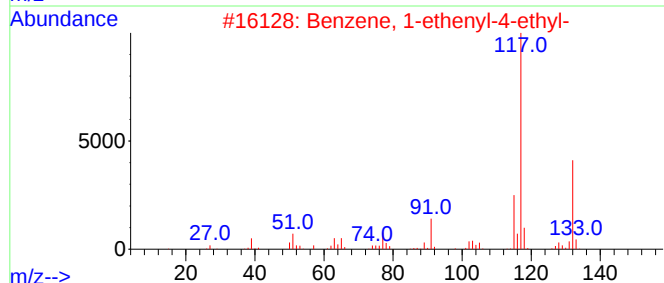
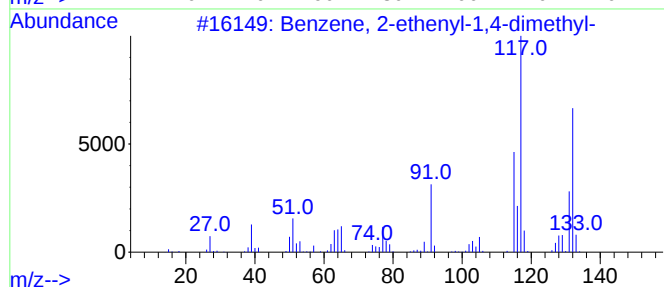
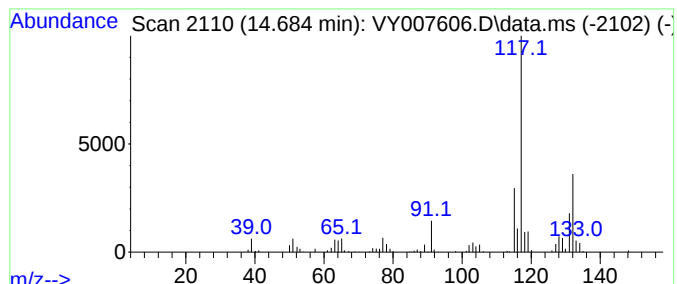
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	20.61 ug/l	206746	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	90
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
5			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY022322\
Data File : VY007606.D
Acq On : 23 Feb 2022 18:29
Operator : SY/MD
Sample : N1584-11
Misc : 5.12g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
D15-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y022322S.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
Pentane, 3-methyl-	5.234	15.8	ug/l	116143	1	7.789	366998	50.0
Cyclopentane, m...	6.795	21.2	ug/l	155850	1	7.789	366998	50.0
Pentane, 2,3-di...	7.868	16.9	ug/l	124182	1	7.789	366998	50.0
Butane, 2,2,3,3...	8.325	28.0	ug/l	278039	2	8.685	496490	50.0
1,3-Dimethylcyc...	9.746	15.6	ug/l	155330	2	8.685	496490	50.0
Benzene, 1,2-di...	13.587	15.9	ug/l	159954	4	13.422	501478	50.0
Deltacyclene	13.623	41.8	ug/l	419450	4	13.422	501478	50.0
p-Cymene	13.964	34.3	ug/l	343599	4	13.422	501478	50.0
Indan, 1-methyl-	14.080	15.2	ug/l	152531	4	13.422	501478	50.0
Benzene, 2-ethe...	14.684	20.6	ug/l	206746	4	13.422	501478	50.0