

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY032922\
 Data File : VY008098.D
 Acq On : 29 Mar 2022 19:33
 Operator : SY/MD
 Sample : N2106-01
 Misc : 5.06g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 OILY-DEBRIS-DRUM-1

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\82Y032522S.M
 Title : SW846 8260

Signal : TIC: VY008098.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.863	3	7	12	rBV2	5166	7098	1.01%	0.100%
2	2.223	60	66	81	rBV	37995	65022	9.24%	0.916%
3	2.388	88	93	102	rVB	17307	30466	4.33%	0.429%
4	3.381	247	256	264	rBV	4757	13218	1.88%	0.186%
5	3.826	322	329	338	rBV	3380	8626	1.23%	0.122%
6	3.948	338	349	365	rBV	125140	344222	48.93%	4.851%
7	4.210	382	392	404	rVB2	4848	14665	2.08%	0.207%
8	4.412	416	425	435	rBV2	6449	20475	2.91%	0.289%
9	4.722	465	476	494	rBV6	7860	35991	5.12%	0.507%
10	4.954	498	514	540	rBV2	77613	291453	41.43%	4.107%
11	5.161	540	548	556	rBV3	4398	15269	2.17%	0.215%
12	5.582	607	617	630	rBV3	5819	19143	2.72%	0.270%
13	5.753	636	645	655	rVB5	5783	17200	2.44%	0.242%
14	6.716	790	803	811	rBV3	11703	40393	5.74%	0.569%
15	6.984	838	847	859	rBV	48362	132756	18.87%	1.871%
16	7.326	894	903	913	rBV5	2858	9798	1.39%	0.138%
17	7.722	958	968	973	rBV	99297	233286	33.16%	3.288%
18	7.789	973	979	989	rVV	167329	385897	54.85%	5.438%
19	8.002	1006	1014	1023	rVB4	4450	10596	1.51%	0.149%
20	8.143	1027	1037	1048	rBV2	99434	243008	34.54%	3.425%
21	8.246	1048	1054	1061	rVV2	8385	22291	3.17%	0.314%
22	8.325	1061	1067	1073	rVV2	8705	24331	3.46%	0.343%
23	8.393	1073	1078	1090	rVB3	9604	23713	3.37%	0.334%
24	8.551	1096	1104	1114	rBV2	26313	57431	8.16%	0.809%
25	8.691	1119	1127	1136	rVV	251099	497537	70.72%	7.011%
26	8.886	1150	1159	1166	rBV2	3876	9676	1.38%	0.136%
27	9.136	1192	1200	1205	rBV	45859	85311	12.13%	1.202%
28	9.179	1205	1207	1214	rVV3	6943	12755	1.81%	0.180%
29	9.283	1214	1224	1233	rVB4	5470	19134	2.72%	0.270%
30	9.466	1244	1254	1257	rBV6	4067	10113	1.44%	0.143%
31	9.520	1257	1263	1269	rVV	11375	24217	3.44%	0.341%
32	9.618	1273	1279	1285	rBV3	5984	11770	1.67%	0.166%
33	9.746	1295	1300	1307	rBV	7117	15260	2.17%	0.215%
34	9.825	1307	1313	1324	rVB6	4148	10750	1.53%	0.151%
35	9.959	1327	1335	1347	rVB5	5872	17480	2.48%	0.246%
36	10.069	1347	1353	1357	rBV2	9987	18119	2.58%	0.255%

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37	10.179	1363	1371	1377	rBV	406394	703510	100.00%	9.914%
38	10.246	1377	1382	1388	rVB	163226	264017	37.53%	3.721%
39	10.368	1395	1402	1409	rVB4	13068	26233	3.73%	0.370%
40	10.496	1409	1423	1431	rBV5	13133	33973	4.83%	0.479%
41	10.606	1433	1441	1445	rBV7	3529	9043	1.29%	0.127%
42	10.721	1450	1460	1468	rVB	26477	51337	7.30%	0.723%
43	10.831	1471	1478	1485	rBV	15845	28186	4.01%	0.397%
44	10.935	1488	1495	1503	rVV3	8330	16797	2.39%	0.237%
45	11.276	1545	1551	1556	rBV4	3603	7143	1.02%	0.101%
46	11.490	1579	1586	1597	rVV	365619	587868	83.56%	8.284%
47	11.593	1597	1603	1609	rVV	47114	78929	11.22%	1.112%
48	11.697	1609	1620	1632	rVV	158488	324367	46.11%	4.571%
49	11.807	1632	1638	1644	rVB4	8411	13804	1.96%	0.195%
50	11.965	1656	1664	1668	rBV3	4755	8620	1.23%	0.121%
51	12.032	1669	1675	1682	rVB	65250	104106	14.80%	1.467%
52	12.105	1682	1687	1695	rVB2	9965	16640	2.37%	0.234%
53	12.209	1698	1704	1708	rBV4	8117	14350	2.04%	0.202%
54	12.331	1718	1724	1730	rBV	4885	7741	1.10%	0.109%
55	12.416	1730	1738	1742	rBV4	4808	10986	1.56%	0.155%
56	12.483	1743	1749	1759	rVB	256997	406038	57.72%	5.722%
57	12.672	1775	1780	1785	rVB	15254	23038	3.27%	0.325%
58	12.733	1785	1790	1794	rBV	50139	84840	12.06%	1.196%
59	12.770	1794	1796	1799	rVV	28603	36814	5.23%	0.519%
60	12.806	1799	1802	1809	rVB3	46728	73123	10.39%	1.030%
61	12.934	1817	1823	1824	rBV3	4916	7526	1.07%	0.106%
62	12.971	1824	1829	1835	rVV	17999	33919	4.82%	0.478%
63	13.062	1838	1844	1848	rVB6	4832	9904	1.41%	0.140%
64	13.123	1848	1854	1859	rBV	85437	130700	18.58%	1.842%
65	13.282	1871	1880	1881	rBV6	6332	13885	1.97%	0.196%
66	13.428	1897	1904	1913	rBV2	243415	414604	58.93%	5.843%
67	13.623	1930	1936	1940	rBV2	21519	36767	5.23%	0.518%
68	13.666	1940	1943	1952	rVB3	11435	23612	3.36%	0.333%
69	13.751	1952	1957	1962	rBV3	29974	54472	7.74%	0.768%
70	13.916	1982	1984	1988	rVB4	7575	8899	1.26%	0.125%
71	13.971	1989	1993	1998	rVB2	19230	31580	4.49%	0.445%
72	14.081	2006	2011	2015	rBV3	18836	33280	4.73%	0.469%
73	14.129	2015	2019	2026	rBV6	10742	24676	3.51%	0.348%
74	14.209	2027	2032	2036	rBV7	10165	18844	2.68%	0.266%
75	14.257	2036	2040	2043	rBV4	29085	48126	6.84%	0.678%
76	14.391	2057	2062	2064	rBV4	18656	32831	4.67%	0.463%

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77	14.422	2064	2067	2071	rVV2	22557	33112	4.71%	0.467%
78	14.477	2071	2076	2081	rVV8	17337	33832	4.81%	0.477%
79	14.617	2095	2099	2103	rVB4	22961	31633	4.50%	0.446%
80	14.684	2104	2110	2115	rBV5	37862	86749	12.33%	1.222%
81	14.733	2115	2118	2125	rVB4	37885	63539	9.03%	0.895%
82	14.800	2125	2129	2130	rBV3	7271	10641	1.51%	0.150%
83	14.843	2131	2136	2139	rBV3	50784	94700	13.46%	1.335%
84	14.946	2149	2153	2157	rVB3	33840	52103	7.41%	0.734%
85	15.446	2232	2235	2240	rBV7	22603	36150	5.14%	0.509%

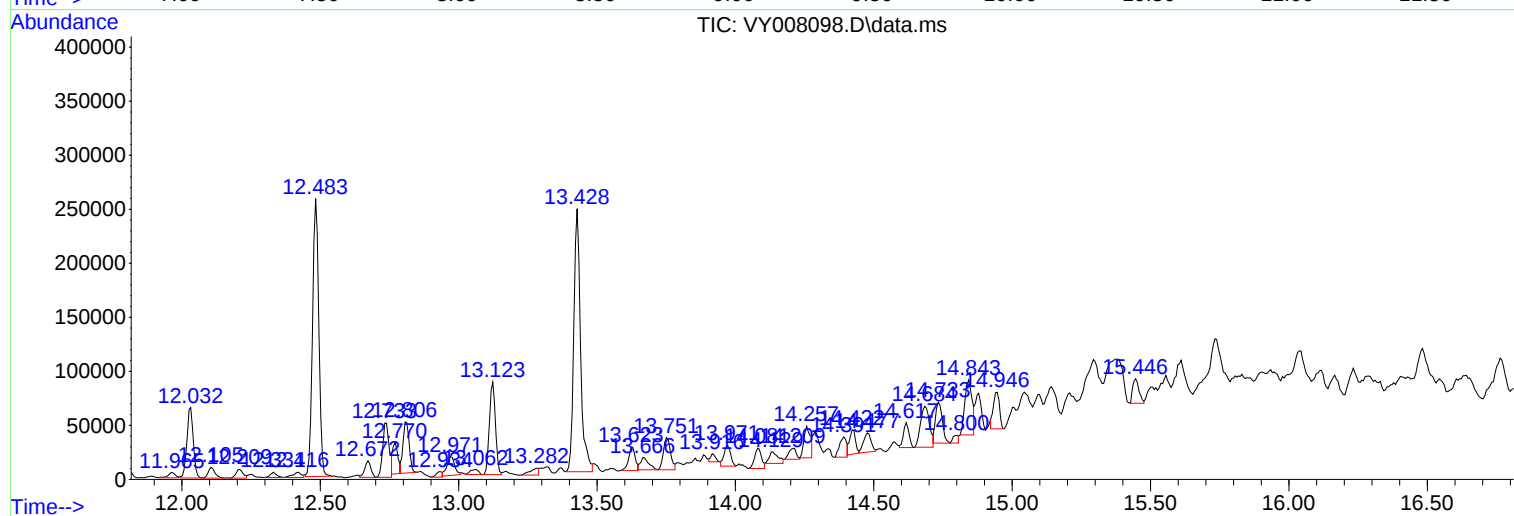
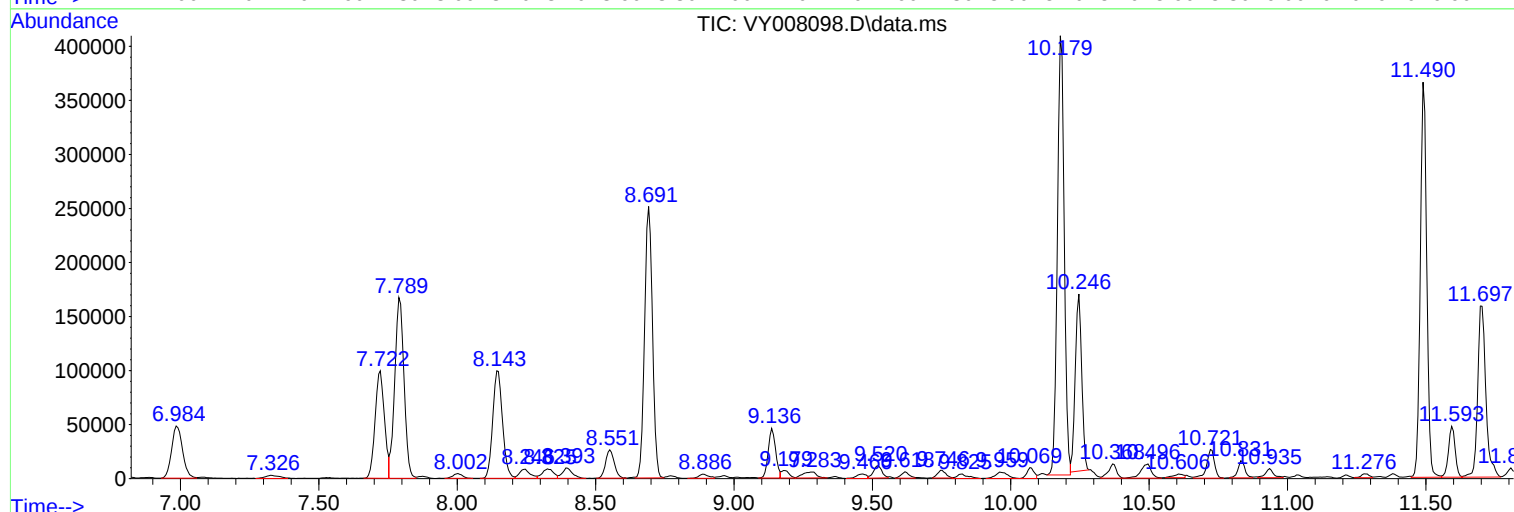
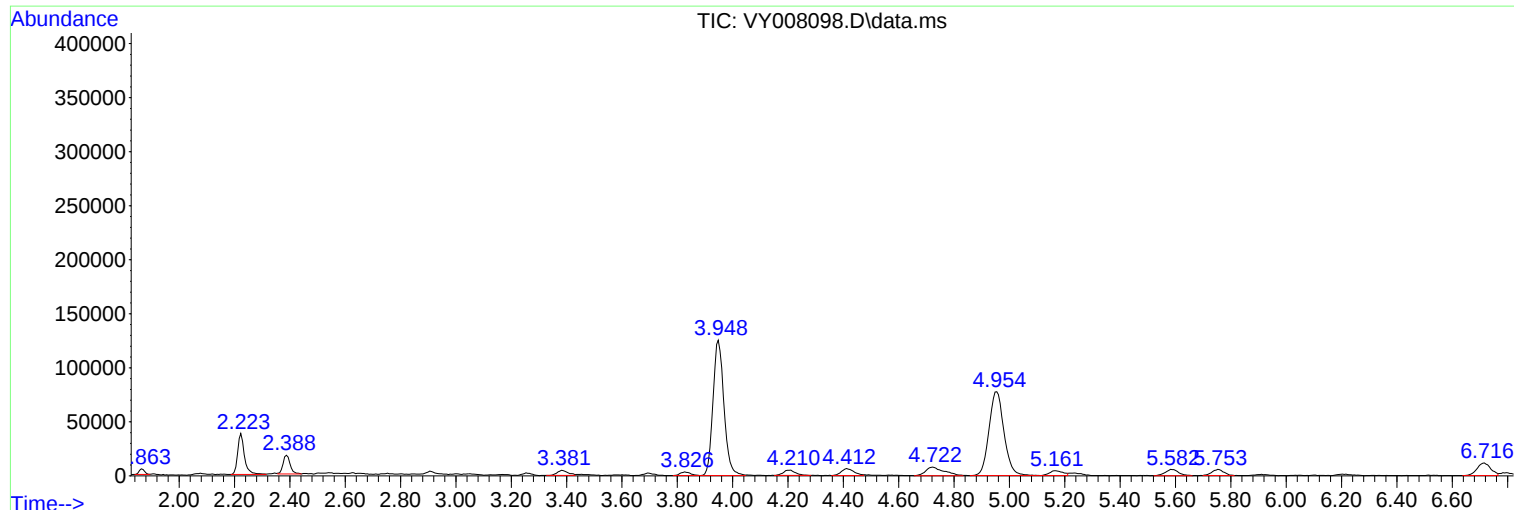
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032522S.M
Quant Title : SW846 8260

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



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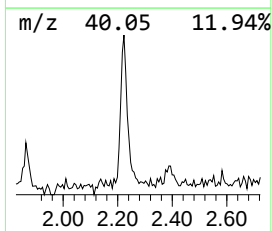
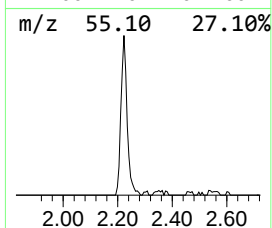
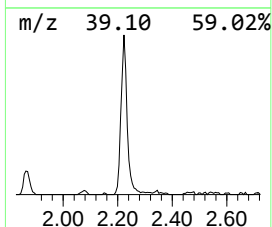
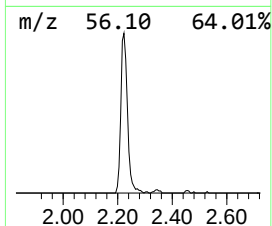
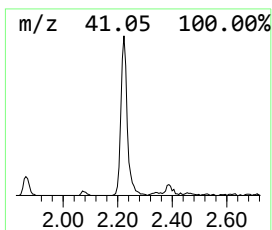
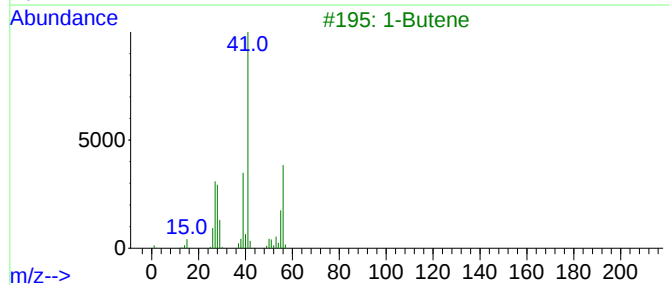
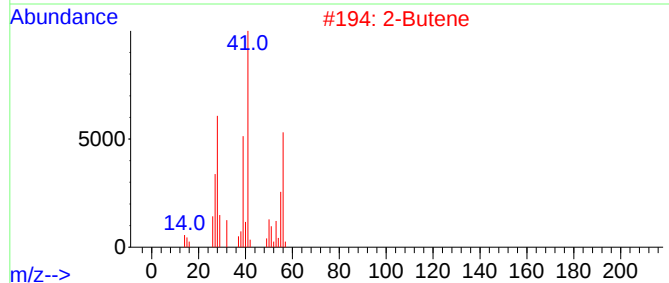
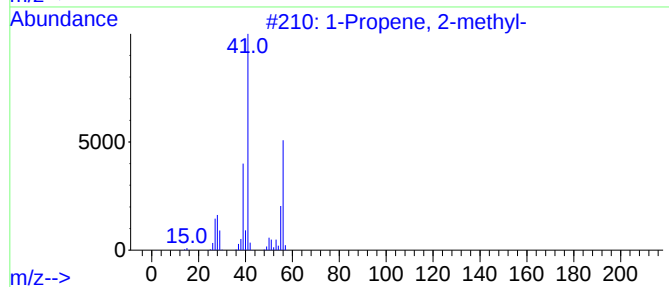
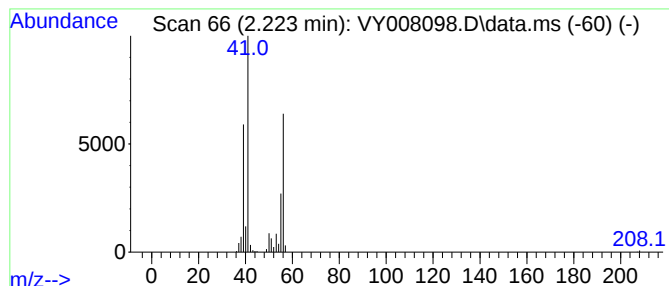
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032522S.M
Quant Title : SW846 8260

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

Peak Number 1 1-Propene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.223	8.42 ug/l	65022	Pentafluorobenzene	7.789

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Propene, 2-methyl-	56	C4H8	000115-11-7	87
2		2-Butene	56	C4H8	000107-01-7	83
3		1-Butene	56	C4H8	000106-98-9	52
4		Cyclobutane	56	C4H8	000287-23-0	52
5		2-Butene, (Z)-	56	C4H8	000590-18-1	52



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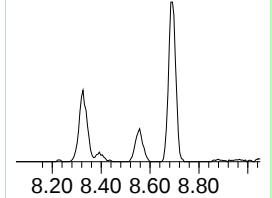
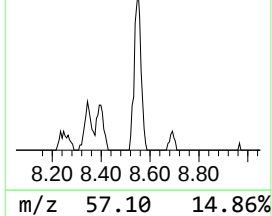
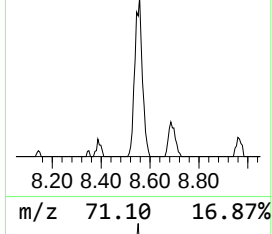
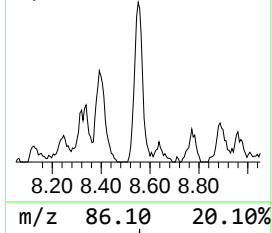
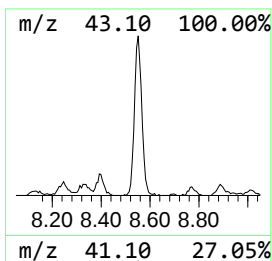
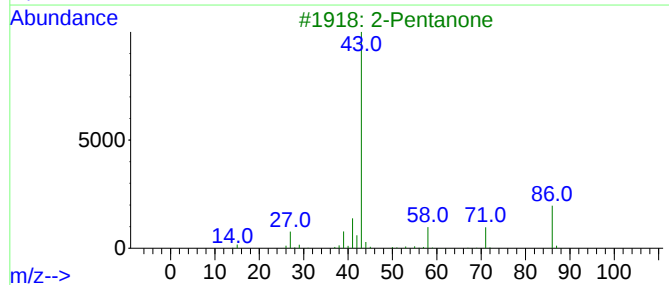
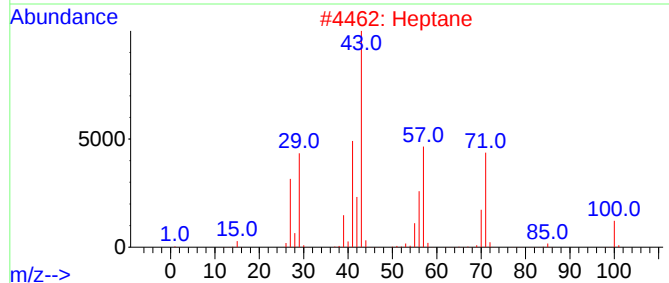
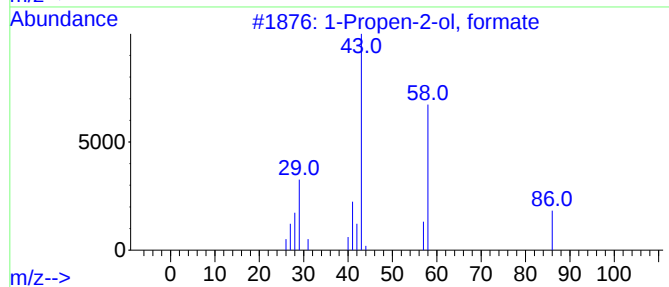
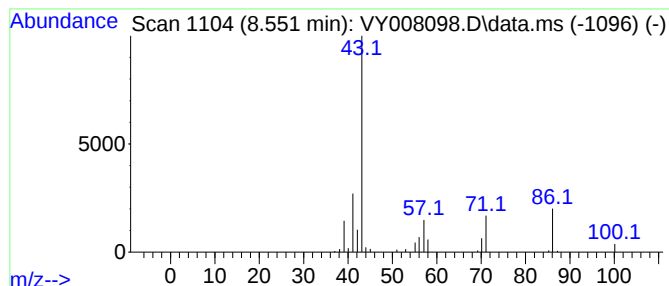
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032522S.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown8.551 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.551	5.77 ug/l	57431	1,4-Difluorobenzene	8.691

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Propen-2-ol, formate	86	C4H6O2	032978-00-0	47
2		Heptane	100	C7H16	000142-82-5	43
3		2-Pentanone	86	C5H10O	000107-87-9	38
4		Oxalic acid, monoamide, n-propyl...	215	C11H21N03	1000309-64-4	38
5		2,3-Butanedione	86	C4H6O2	000431-03-8	32



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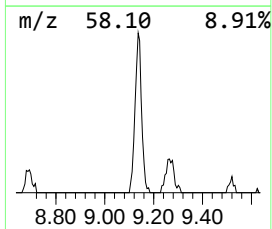
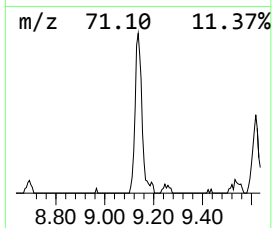
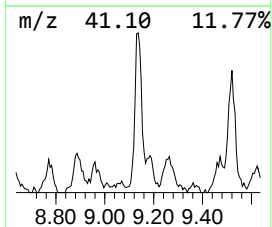
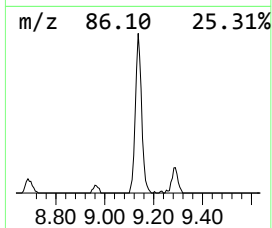
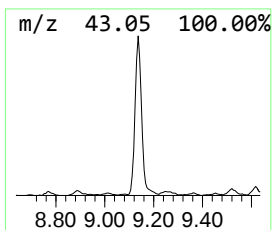
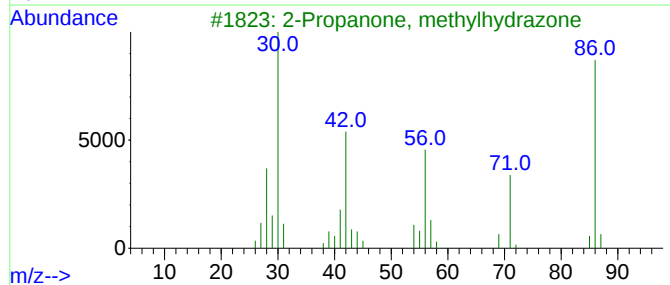
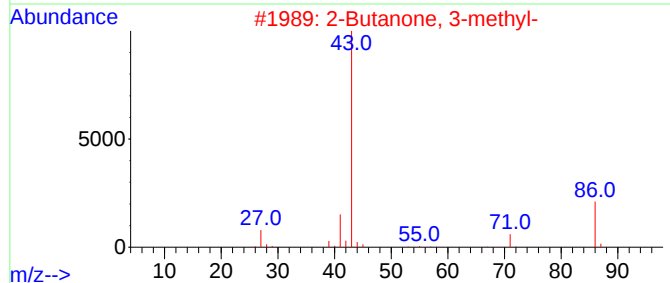
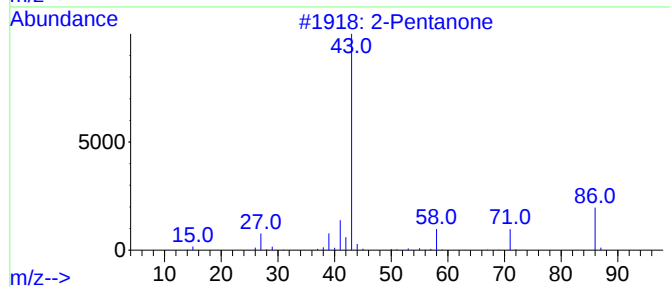
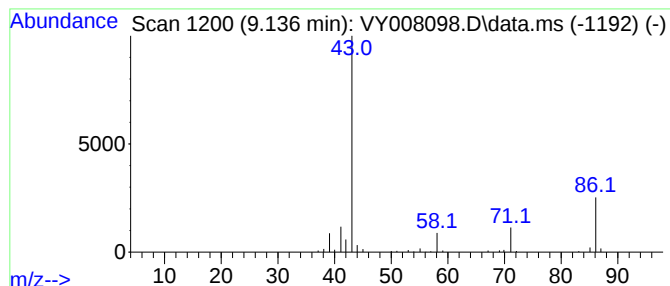
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032522S.M
Quant Title : SW846 8260

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

Peak Number 3 2-Pentanone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.136	8.57 ug/l	85311	1,4-Difluorobenzene	8.691

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone	86	C5H10O	000107-87-9	86
2			2-Butanone, 3-methyl-	86	C5H10O	000563-80-4	64
3			2-Propanone, methylhydrazone	86	C4H10N2	005771-02-8	23
4			Propane, 2-(ethenyloxy)-	86	C5H10O	000926-65-8	9
5			2,3-Butanedione	86	C4H6O2	000431-03-8	9



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ClientSampleId :
OILY-DEBRIS-DRUM-1

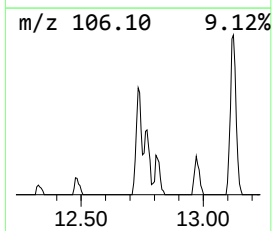
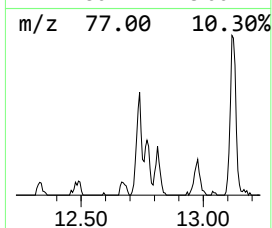
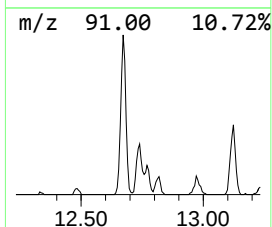
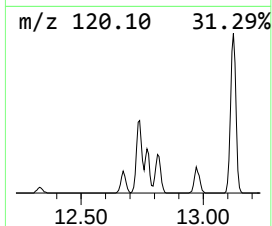
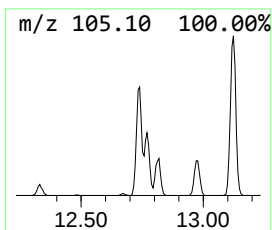
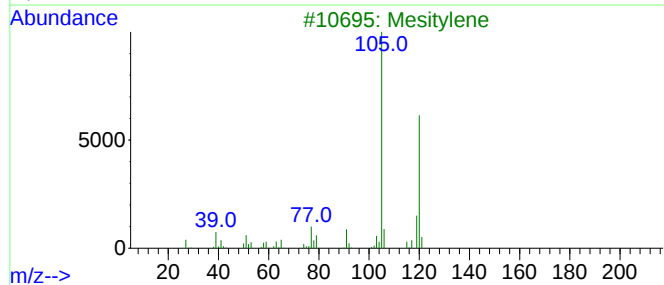
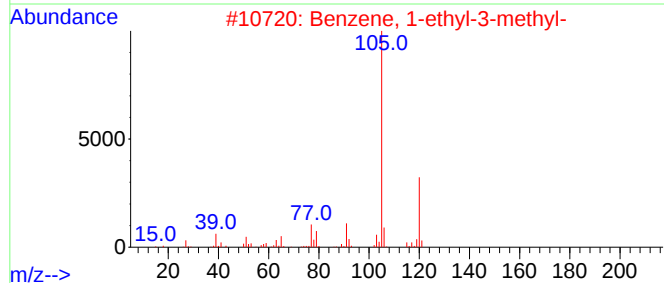
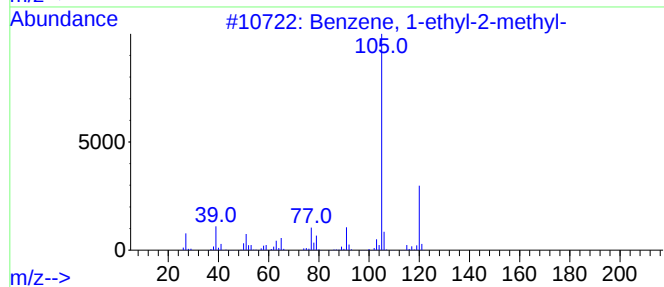
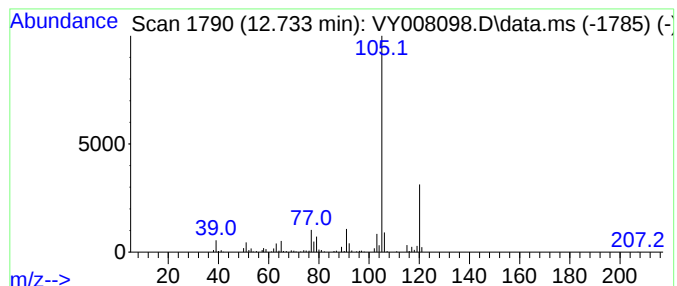
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.733	10.23 ug/l	84840	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY032922\
Data File : VY008098.D
Acq On : 29 Mar 2022 19:33
Operator : SY/MD
Sample : N2106-01
Misc : 5.06g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OILY-DEBRIS-DRUM-1

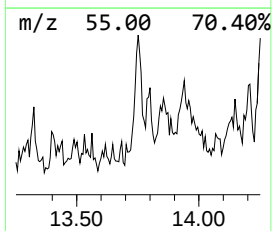
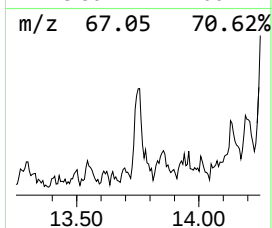
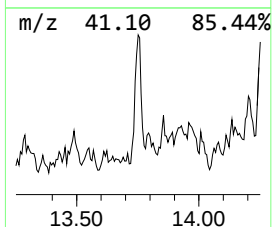
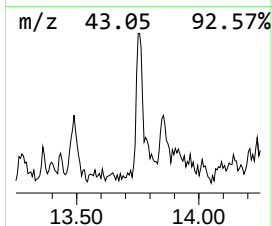
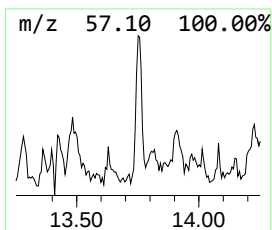
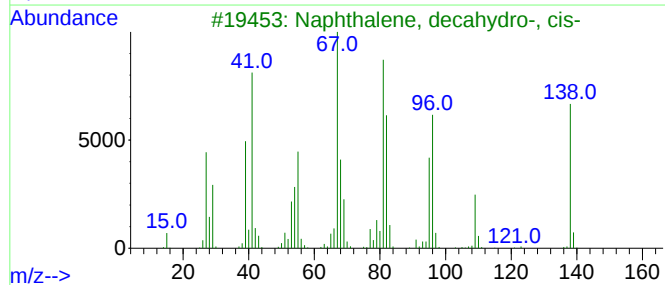
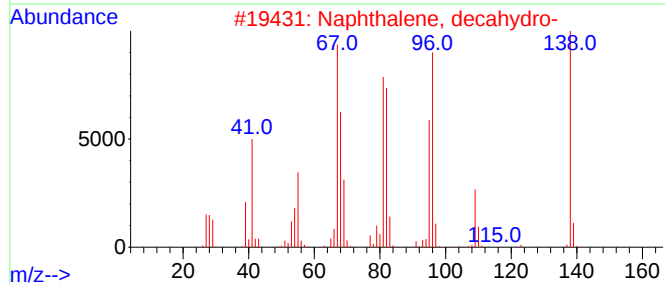
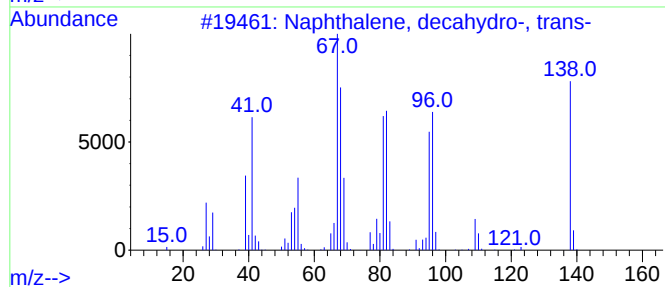
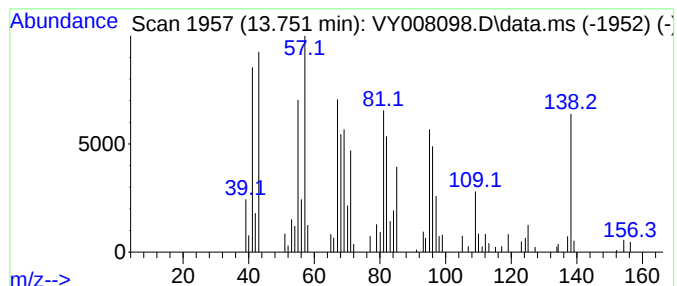
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Quant Title : SW846 8260

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

Peak Number 5 Naphthalene, decahydro-, tr... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.751	6.57 ug/l	54472	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	70
2			Naphthalene, decahydro-	138	C10H18	000091-17-8	55
3			Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	50
4			2-Decen-1-ol, (E)-	156	C10H20O	018409-18-2	49
5			Citronellyl butyrate	226	C14H26O2	000141-16-2	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY032922\
Data File : VY008098.D
Acq On : 29 Mar 2022 19:33
Operator : SY/MD
Sample : N2106-01
Misc : 5.06g/5.0mL/MSVOA_Y/SoIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OILY-DEBRIS-DRUM-1

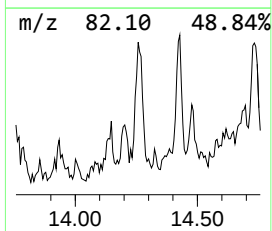
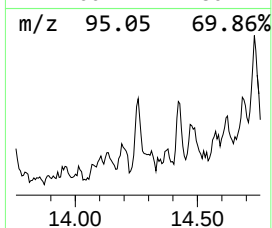
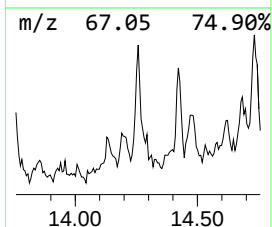
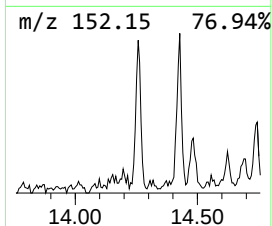
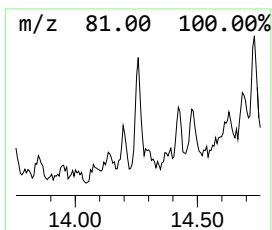
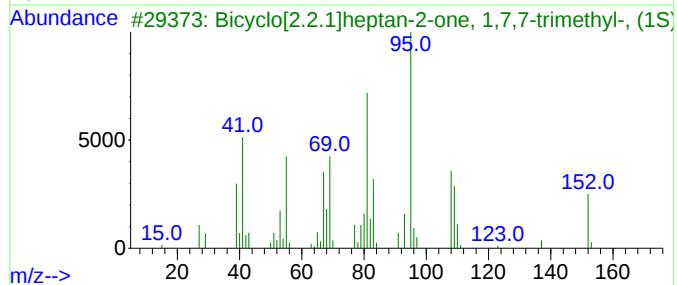
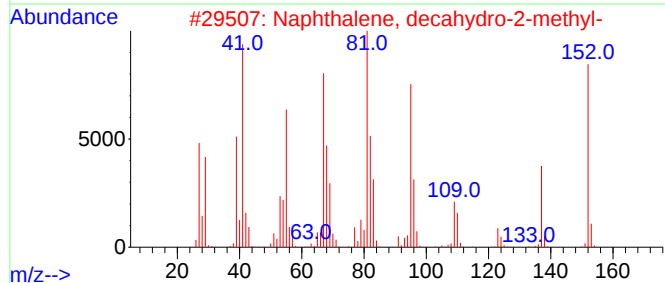
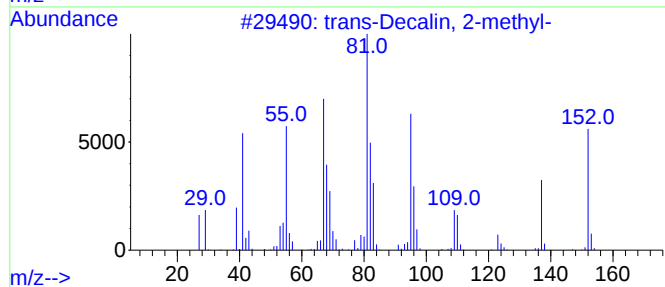
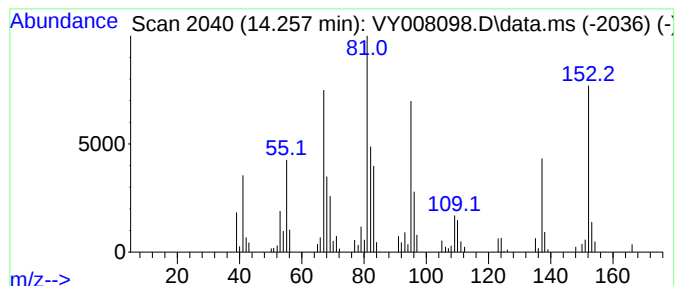
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Quant Title : SW846 8260

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

Peak Number 6 trans-Decalin, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.257	5.80 ug/l	48126	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	96
2			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	93
3			Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	76
4			Cyclohexanone, 5-methyl-2-(1-met...	152	C10H16O	015932-80-6	76
5			Camphor	152	C10H16O	000076-22-2	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY032922\
Data File : VY008098.D
Acq On : 29 Mar 2022 19:33
Operator : SY/MD
Sample : N2106-01
Misc : 5.06g/5.0mL/MSVOA_Y/soil
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OILY-DEBRIS-DRUM-1

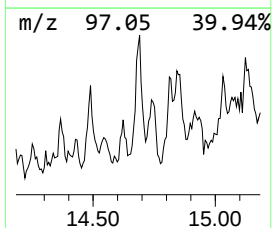
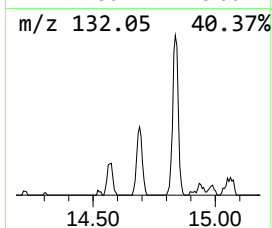
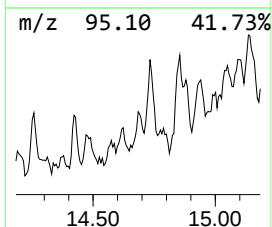
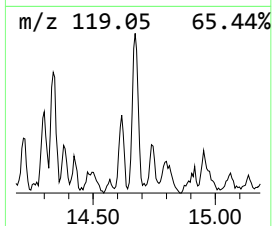
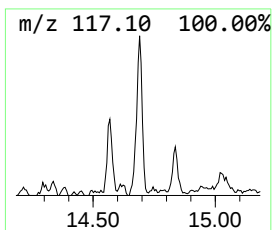
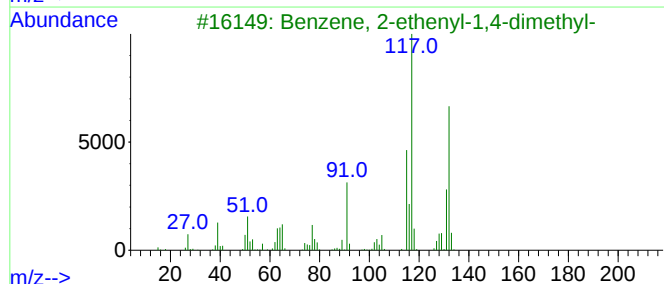
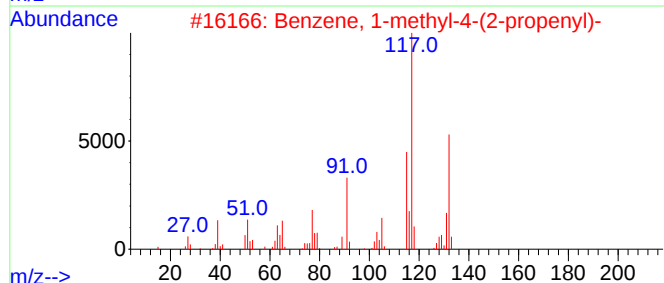
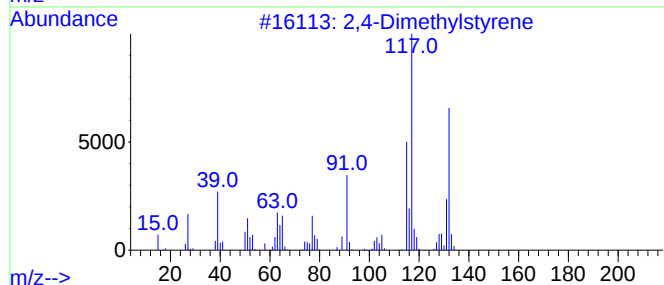
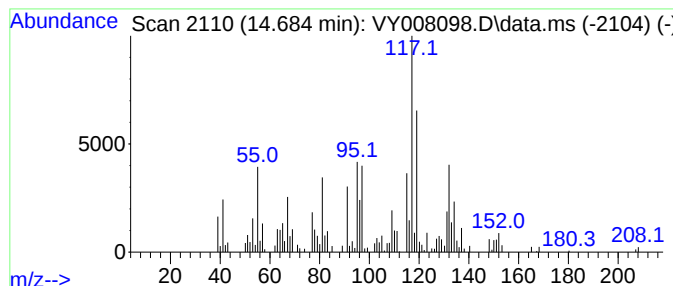
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Quant Title : SW846 8260

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

Peak Number 7 2,4-Dimethylstyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	10.46 ug/l	86749	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4-Dimethylstyrene	132	C10H12	002234-20-0	58
2			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	56
3			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	56
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	55
5			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY032922\
Data File : VY008098.D
Acq On : 29 Mar 2022 19:33
Operator : SY/MD
Sample : N2106-01
Misc : 5.06g/5.0mL/MSVOA_Y/soil
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OILY-DEBRIS-DRUM-1

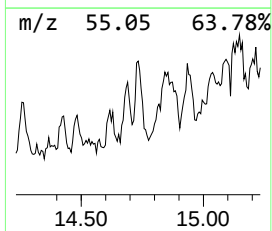
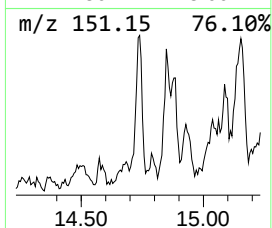
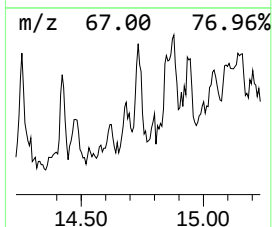
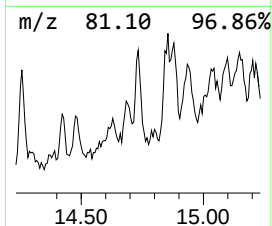
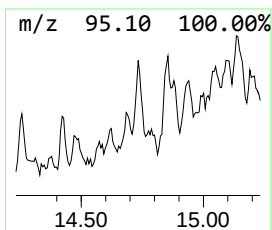
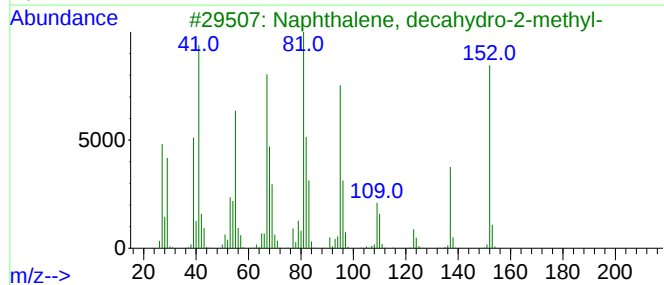
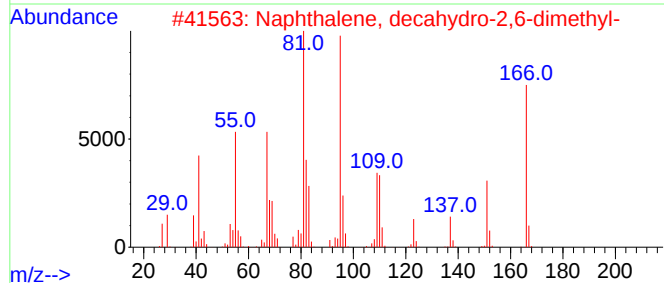
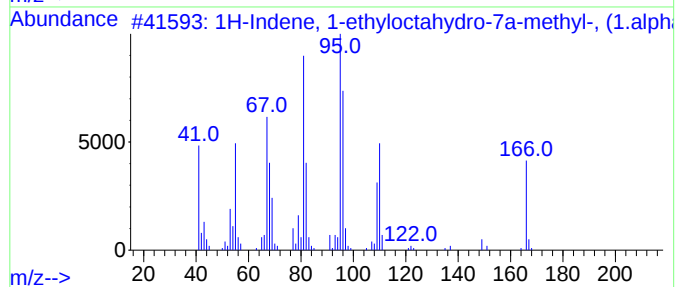
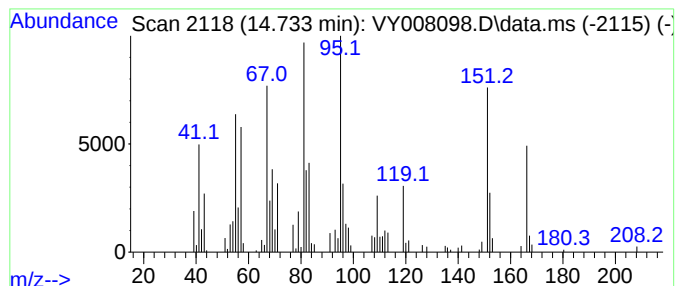
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Quant Title : SW846 8260

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

Peak Number 8 Naphthalene, decahydro-2,6-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.733	7.66 ug/l	63539	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 1-ethyloctahydro-7a-m...	166	C12H22	056324-71-1	78
2			Naphthalene, decahydro-2,6-dimet...	166	C12H22	001618-22-0	70
3			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	62
4			cis,trans-1,6-Dimethylspiro[4.5]...	166	C12H22	1000111-72-3	53
5			cis-Decalin, 2-syn-methyl-	152	C11H20	1000155-85-6	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY032922\
Data File : VY008098.D
Acq On : 29 Mar 2022 19:33
Operator : SY/MD
Sample : N2106-01
Misc : 5.06g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OILY-DEBRIS-DRUM-1

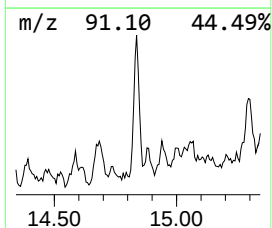
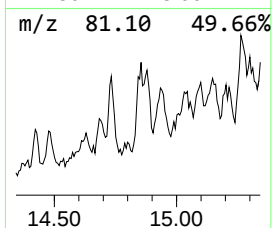
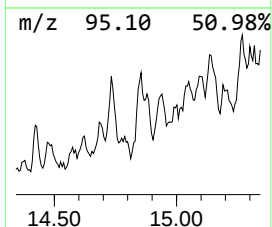
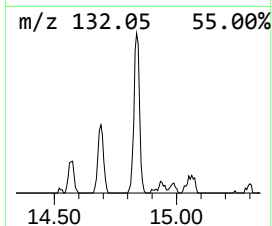
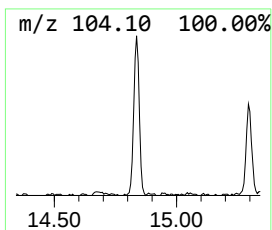
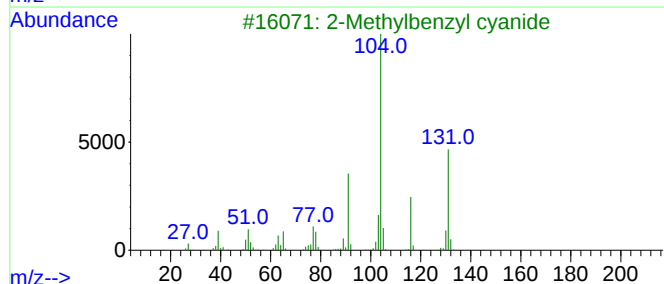
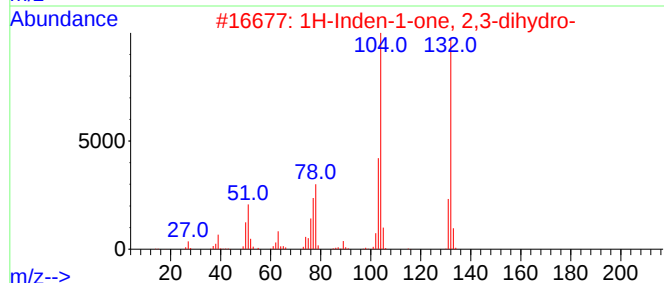
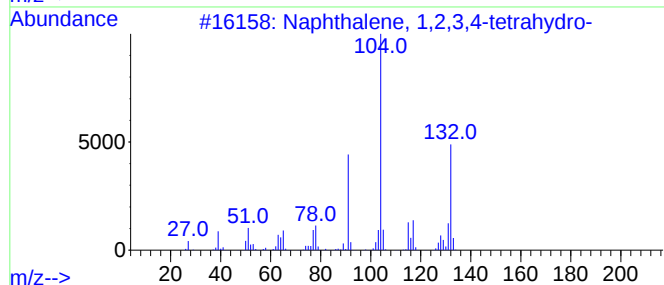
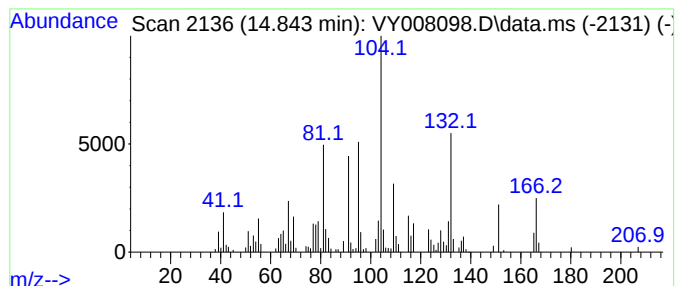
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.843	11.42 ug/l	94700	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2			1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	38
3			2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	38
4			2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	35
5			1H-Indol-3-amine	132	C8H8N2	007250-19-3	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY032922\
Data File : VY008098.D
Acq On : 29 Mar 2022 19:33
Operator : SY/MD
Sample : N2106-01
Misc : 5.06g/5.0mL/MSVOA_Y/soil
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OILY-DEBRIS-DRUM-1

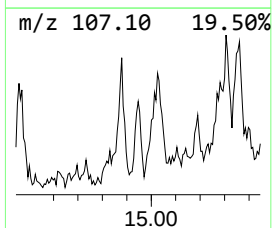
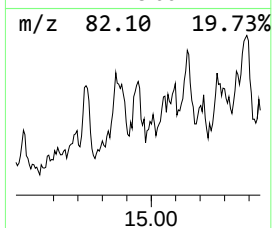
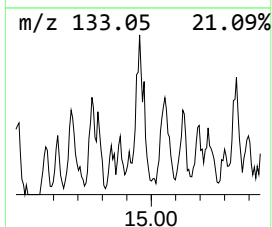
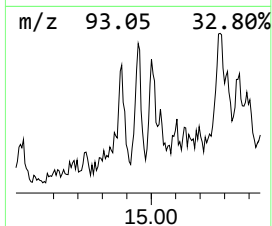
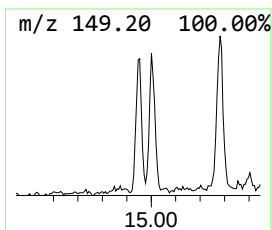
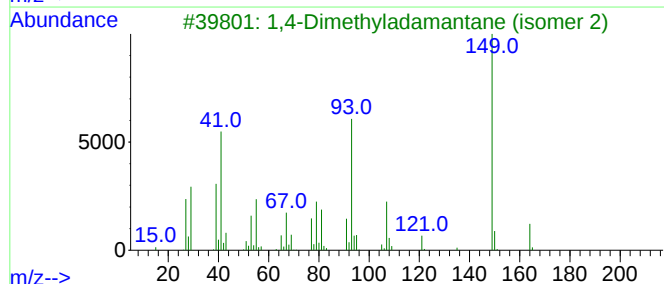
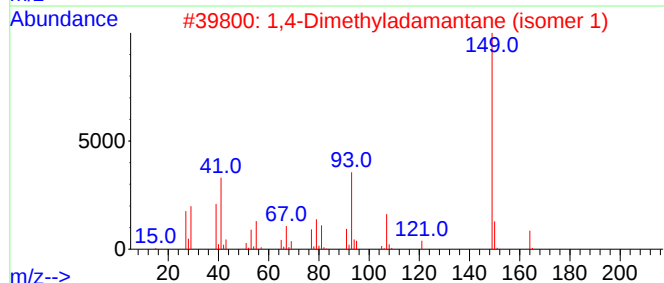
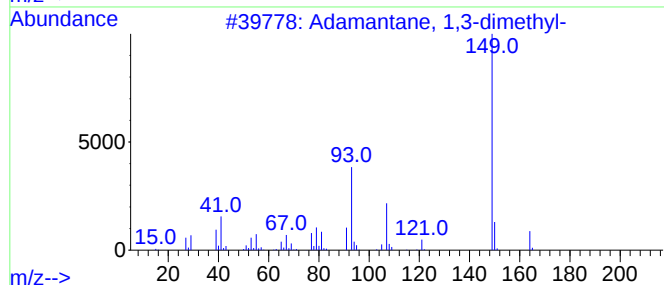
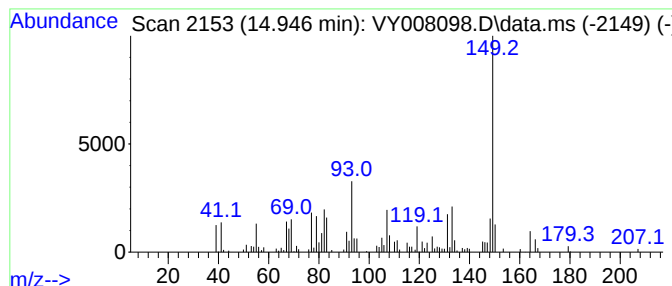
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032522S.M
Quant Title : SW846 8260

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

Peak Number 10 Adamantane, 1,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.946	6.28 ug/l	52103	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Adamantane, 1,3-dimethyl-	164	C12H20	000702-79-4	64
2			1,4-Dimethyladamantane (isomer 1)	164	C12H20	1000460-67-0	55
3			1,4-Dimethyladamantane (isomer 2)	164	C12H20	1000460-67-1	49
4			Pyridine, pentamethyl-	149	C10H15N	003748-83-2	46
5			Benzothiazole, 2-methyl-	149	C8H7NS	000120-75-2	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY032922\
Data File : VY008098.D
Acq On : 29 Mar 2022 19:33
Operator : SY/MD
Sample : N2106-01
Misc : 5.06g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
OILY-DEBRIS-DRUM-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032522S.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
1-Propene, 2-me...	2.223	8.4	ug/l	65022	1	7.789	385897	50.0
unknown8.551	8.551	5.8	ug/l	57431	2	8.691	497537	50.0
2-Pentanone	9.136	8.6	ug/l	85311	2	8.691	497537	50.0
Benzene, 1-ethy...	12.733	10.2	ug/l	84840	4	13.428	414604	50.0
Naphthalene, de...	13.751	6.6	ug/l	54472	4	13.428	414604	50.0
trans-Decalin, ...	14.257	5.8	ug/l	48126	4	13.428	414604	50.0
2,4-Dimethylsty...	14.684	10.5	ug/l	86749	4	13.428	414604	50.0
Naphthalene, de...	14.733	7.7	ug/l	63539	4	13.428	414604	50.0
Naphthalene, 1,...	14.843	11.4	ug/l	94700	4	13.428	414604	50.0
Adamantane, 1,3...	14.946	6.3	ug/l	52103	4	13.428	414604	50.0