

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY060321\
 Data File : VY004945.D
 Acq On : 03 Jun 2021 18:32
 Operator : SY/MD
 Sample : M2589-15
 Misc : 12.82G/5ML/MSVOA_Y/SOIL
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 18-B11(2-4)

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\82Y060221S.M

Title : SW846 8260

Signal : TIC: VY004945.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.229	63	67	76	rVB3	12789	29789	1.84%	0.248%
2	2.912	173	179	185	rVB2	8887	17372	1.07%	0.145%
3	3.960	340	351	359	rBV	21908	63122	3.90%	0.526%
4	4.204	381	391	409	rBV	76650	221737	13.69%	1.848%
5	7.728	959	969	974	rBV	152391	356207	21.99%	2.968%
6	7.795	974	980	990	rVV	265355	613565	37.88%	5.113%
7	8.008	1006	1015	1021	rBV4	9092	22225	1.37%	0.185%
8	8.149	1028	1038	1052	rBV2	194167	477017	29.45%	3.975%
9	8.331	1063	1068	1077	rVV3	8970	28364	1.75%	0.236%
10	8.551	1096	1104	1113	rBV3	9003	20458	1.26%	0.170%
11	8.697	1119	1128	1141	rBV	435533	844992	52.17%	7.041%
12	9.191	1196	1209	1215	rBV3	25710	68011	4.20%	0.567%
13	9.618	1271	1279	1287	rBV3	9826	19436	1.20%	0.162%
14	9.758	1290	1302	1307	rBV4	10869	25760	1.59%	0.215%
15	9.825	1307	1313	1316	rVV	10661	18680	1.15%	0.156%
16	9.959	1326	1335	1347	rVB4	11898	35452	2.19%	0.295%
17	10.185	1364	1372	1378	rBV	708107	1223548	75.55%	10.195%
18	10.246	1378	1382	1388	rVB	39154	67511	4.17%	0.563%
19	10.374	1395	1403	1411	rVB5	17623	40452	2.50%	0.337%
20	10.526	1423	1428	1433	rVB	10868	17795	1.10%	0.148%
21	10.618	1433	1443	1450	rBV9	9122	26672	1.65%	0.222%
22	10.715	1454	1459	1465	rVB3	9982	17918	1.11%	0.149%
23	11.038	1508	1512	1516	rVB3	13485	19668	1.21%	0.164%
24	11.093	1516	1521	1526	rVB	16664	27668	1.71%	0.231%
25	11.282	1544	1552	1563	rVB4	19584	54071	3.34%	0.451%
26	11.380	1563	1568	1573	rBV	13797	22012	1.36%	0.183%
27	11.496	1579	1587	1597	rBV	611632	1014894	62.66%	8.457%
28	11.593	1597	1603	1612	rVV	33547	67089	4.14%	0.559%
29	11.703	1612	1621	1631	rVV3	89213	226169	13.96%	1.885%
30	12.032	1662	1675	1683	rBV2	68768	157318	9.71%	1.311%
31	12.123	1685	1690	1694	rVB3	16862	26976	1.67%	0.225%
32	12.203	1698	1703	1706	rVV3	15653	27932	1.72%	0.233%
33	12.239	1706	1709	1714	rVB3	19264	35275	2.18%	0.294%
34	12.337	1719	1725	1730	rBV2	95506	160942	9.94%	1.341%
35	12.489	1744	1750	1761	rVB3	922391	1619628	100.00%	13.496%
36	12.581	1761	1765	1769	rVB7	11554	19097	1.18%	0.159%

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Integrator: RTE

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Filtering: 5

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Max Peaks: 100

Stop Thrs : 0

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37	12.648	1772	1776	1784	rVB3	25243	57136	3.53%	0.476%
38	12.733	1784	1790	1794	rBV2	40673	72297	4.46%	0.602%
39	12.806	1794	1802	1808	rVB4	124755	238700	14.74%	1.989%
40	12.971	1822	1829	1837	rBV3	41429	103153	6.37%	0.860%
41	13.038	1837	1840	1848	rVB2	33421	52993	3.27%	0.442%
42	13.123	1848	1854	1859	rBV	80393	124336	7.68%	1.036%
43	13.251	1867	1875	1878	rBV5	11063	24431	1.51%	0.204%
44	13.337	1883	1889	1892	rBV4	25957	61195	3.78%	0.510%
45	13.428	1897	1904	1913	rVV	600655	987021	60.94%	8.224%
46	13.629	1932	1937	1940	rVV2	44238	73662	4.55%	0.614%
47	13.666	1940	1943	1951	rVB2	34524	63838	3.94%	0.532%
48	13.751	1952	1957	1962	rBV	137614	199473	12.32%	1.662%
49	13.916	1981	1984	1987	rVV4	22325	30365	1.87%	0.253%
50	13.971	1988	1993	1998	rVB2	59456	95924	5.92%	0.799%
51	14.019	1998	2001	2006	rVB4	12567	18783	1.16%	0.157%
52	14.080	2006	2011	2016	rBV2	34975	51883	3.20%	0.432%
53	14.215	2028	2033	2035	rBV5	19121	35069	2.17%	0.292%
54	14.269	2036	2042	2044	rVV6	27804	71188	4.40%	0.593%
55	14.312	2046	2049	2057	rVV2	48735	112114	6.92%	0.934%
56	14.385	2057	2061	2064	rVV3	26030	40090	2.48%	0.334%
57	14.422	2064	2067	2072	rVB3	21863	31593	1.95%	0.263%
58	14.568	2087	2091	2094	rBV2	13750	23655	1.46%	0.197%
59	14.611	2094	2098	2103	rVB2	94244	125690	7.76%	1.047%
60	14.684	2103	2110	2114	rBV5	44549	102439	6.32%	0.854%
61	14.733	2114	2118	2125	rVB2	44822	86294	5.33%	0.719%
62	14.946	2148	2153	2161	rBV5	35934	62007	3.83%	0.517%
63	15.056	2166	2171	2175	rBV4	21746	41866	2.58%	0.349%
64	15.135	2180	2184	2190	rVB7	25762	51453	3.18%	0.429%
65	15.239	2192	2201	2206	rBV	277117	515622	31.84%	4.296%
66	15.342	2213	2218	2222	rBV7	11599	23235	1.43%	0.194%
67	15.434	2229	2233	2242	rVB2	57861	97388	6.01%	0.811%
68	15.531	2242	2249	2255	rBV5	16040	41581	2.57%	0.346%
69	15.605	2257	2261	2266	rVB6	13430	22210	1.37%	0.185%
70	15.690	2268	2275	2278	rVV6	11233	27847	1.72%	0.232%
71	15.727	2278	2281	2291	rVB4	14715	28625	1.77%	0.239%
72	15.891	2302	2308	2314	rBV6	11544	29543	1.82%	0.246%
73	16.031	2324	2331	2336	rBV7	13579	30818	1.90%	0.257%
74	16.159	2347	2352	2356	rBV3	11324	18327	1.13%	0.153%
75	16.233	2358	2364	2368	rVV5	12473	29062	1.79%	0.242%
76	16.293	2368	2374	2380	rVV	88045	174785	10.79%	1.456%

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18-B11(2-4)

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Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

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77	16.361	2380	2385	2392	rVB	25397	46048	2.84%	0.384%
78	16.489	2399	2406	2421	rVB	60359	142467	8.80%	1.187%

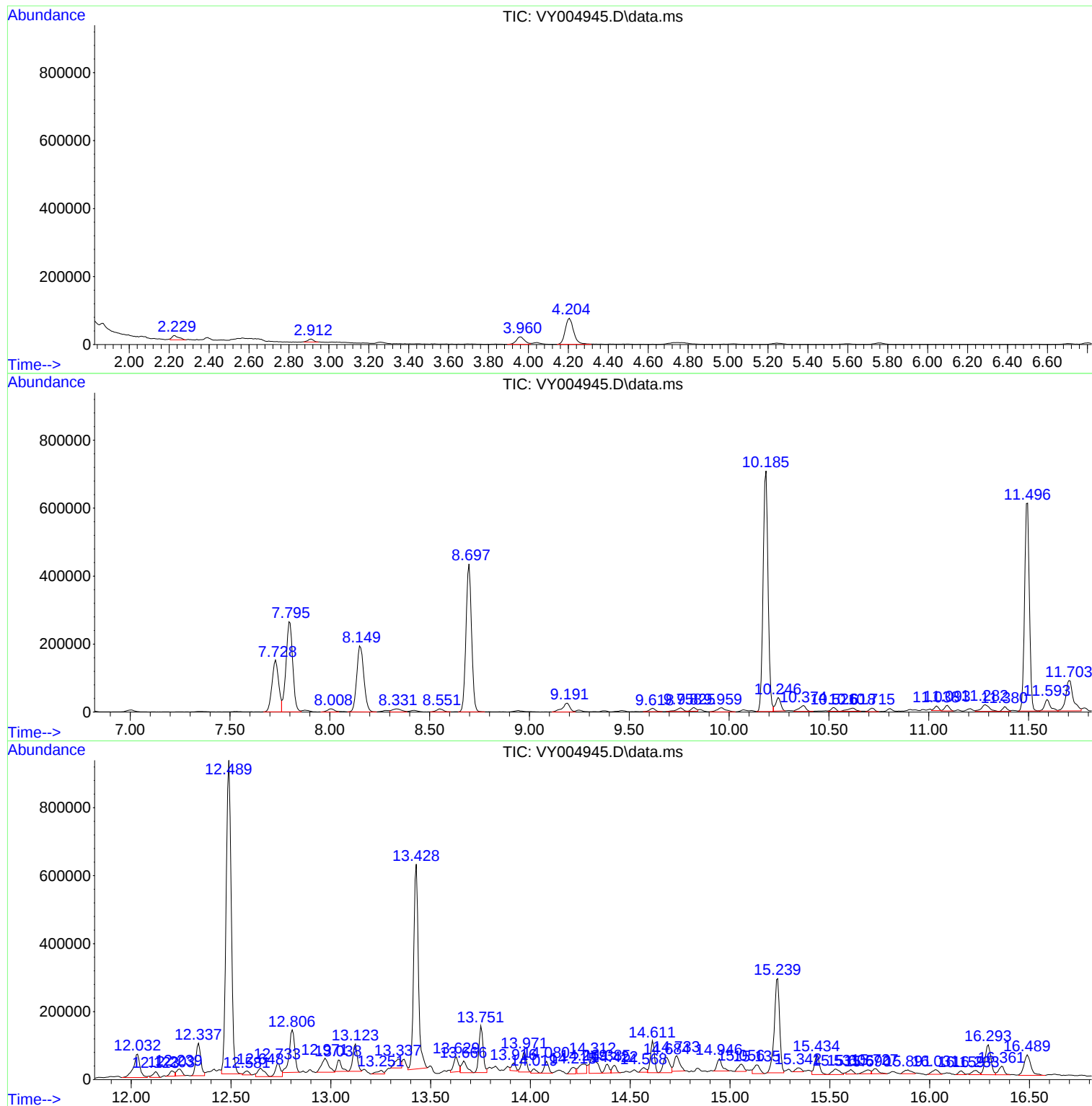
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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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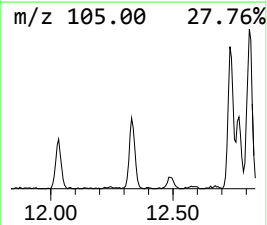
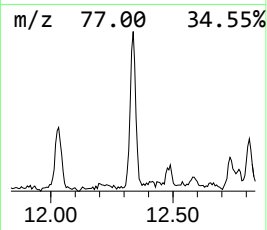
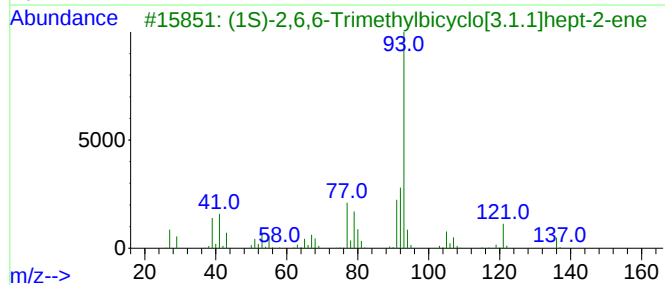
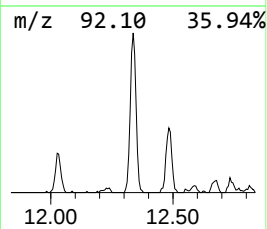
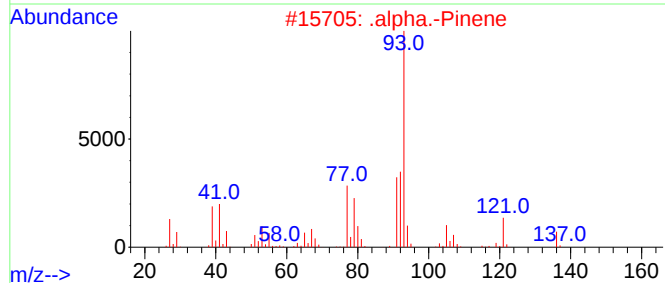
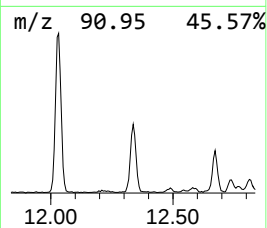
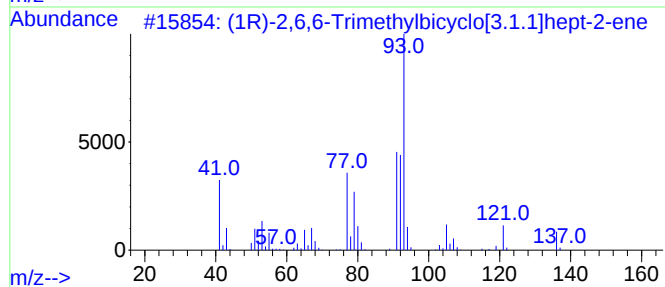
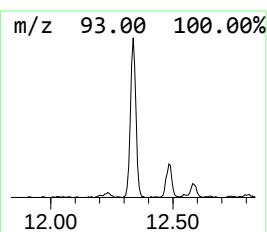
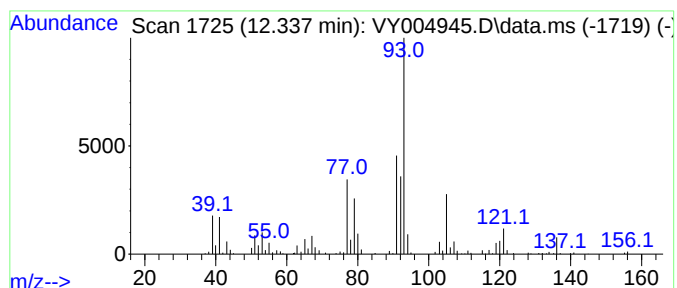
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.337	7.93 ug/l	160942	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1R)-2,6,6-Trimethylbicyclo[3.1...	136	C10H16	007785-70-8	96		
2	.alpha.-Pinene	136	C10H16	000080-56-8	93		
3	(1S)-2,6,6-Trimethylbicyclo[3.1...	136	C10H16	007785-26-4	90		
4	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	90		
5	.gamma.-Terpinene	136	C10H16	000099-85-4	87		



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Instrument :
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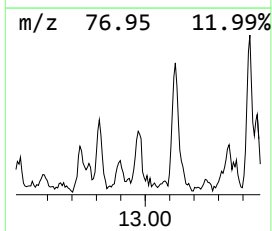
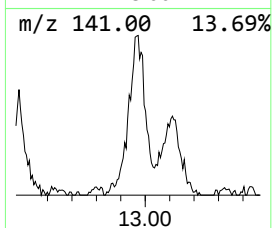
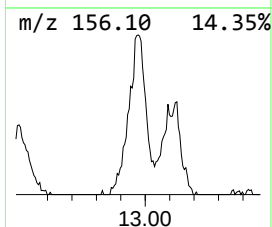
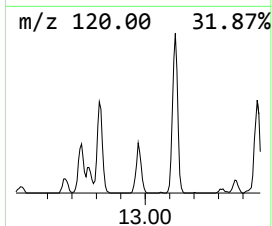
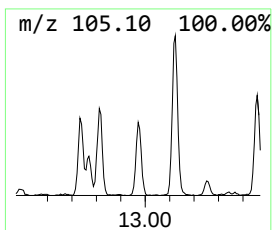
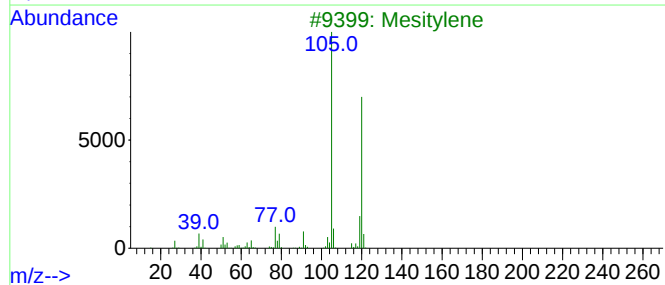
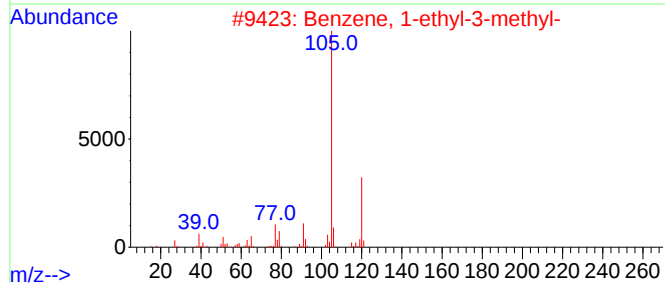
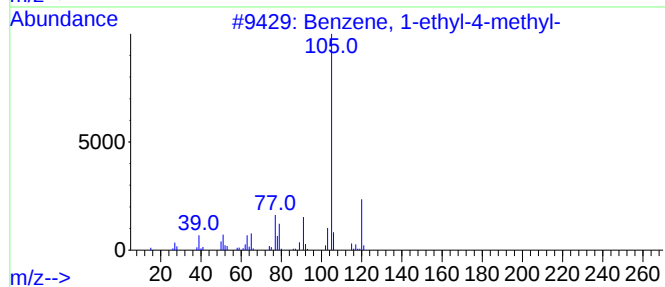
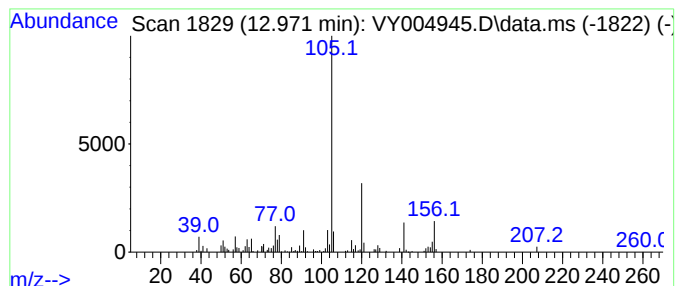
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1-ethyl-4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.971	5.23 ug/l	103153	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	76
3			Mesitylene	120	C9H12	000108-67-8	76
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	76
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	76



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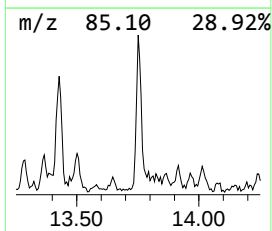
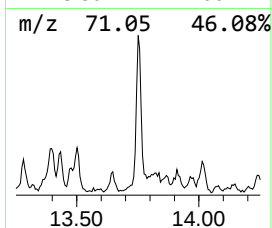
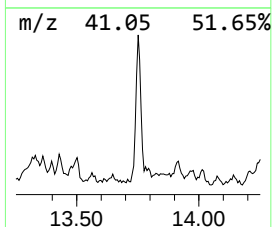
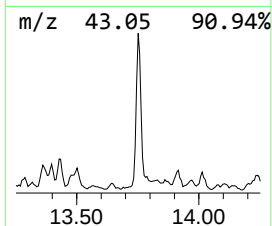
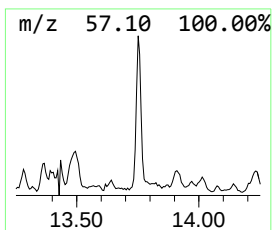
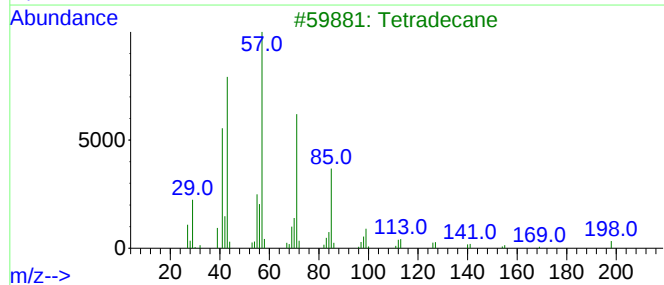
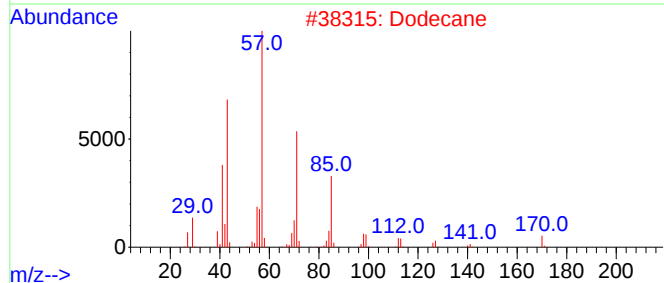
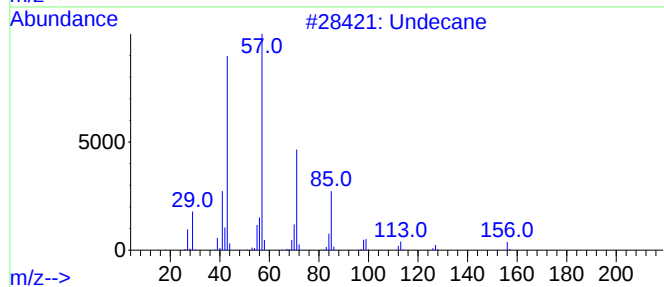
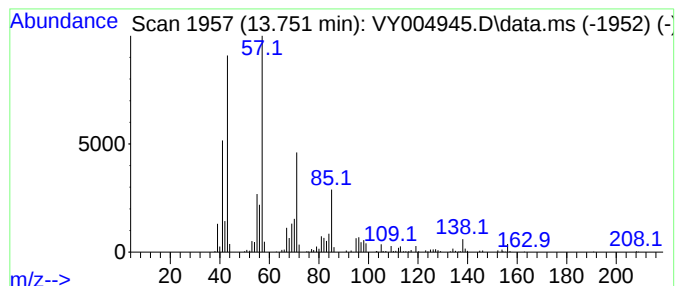
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Undecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.751	10.10 ug/l	199473	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	87
2			Dodecane	170	C12H26	000112-40-3	72
3			Tetradecane	198	C14H30	000629-59-4	64
4			Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	59
5			Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	53



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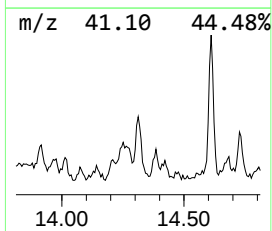
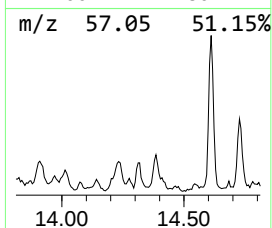
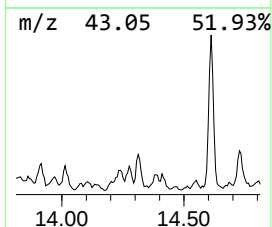
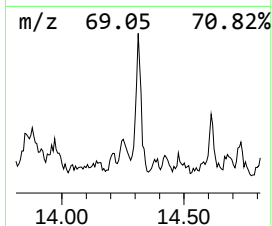
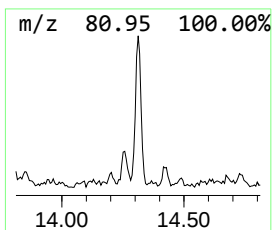
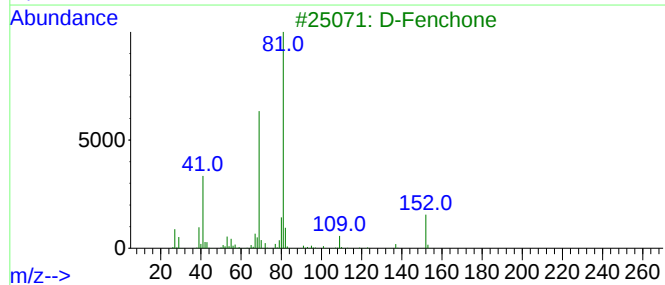
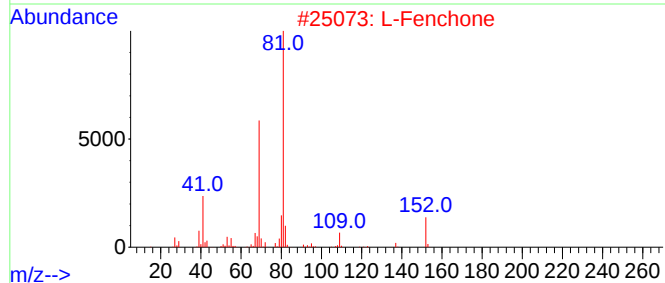
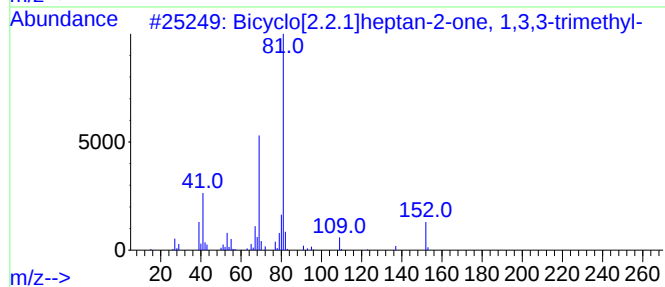
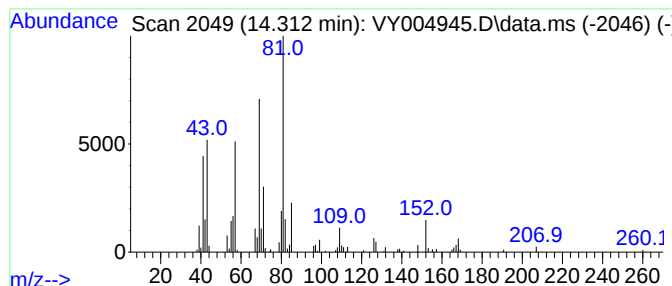
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.312	5.68 ug/l	112114	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	58
2			L-Fenchone	152	C10H16O	007787-20-4	52
3			D-Fenchone	152	C10H16O	004695-62-9	52
4			Furan, 2-hexyl-	152	C10H16O	003777-70-6	38
5			Undecane, 2-methyl-	170	C12H26	007045-71-8	14



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY060321\
Data File : VY004945.D
Acq On : 03 Jun 2021 18:32
Operator : SY/MD
Sample : M2589-15
Misc : 12.82G/5ML/MSVOA_Y/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
18-B11(2-4)

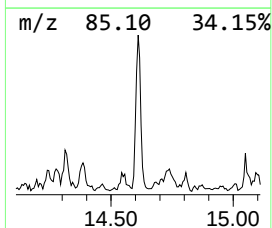
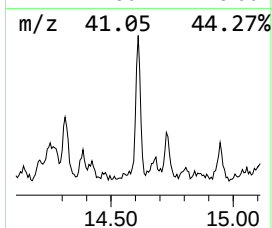
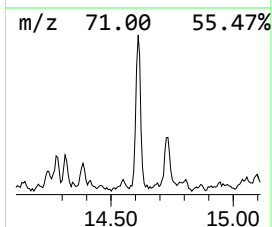
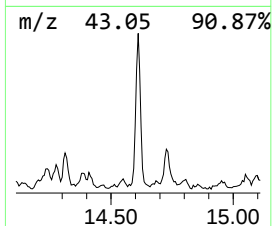
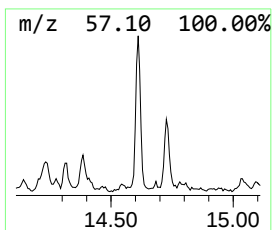
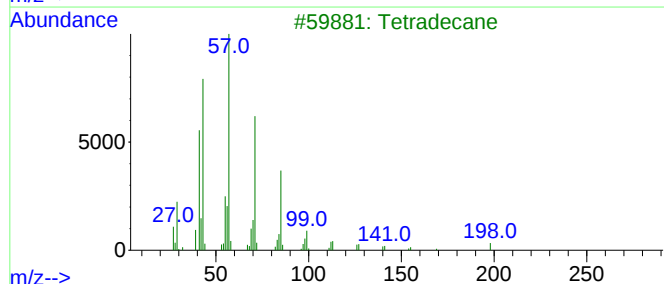
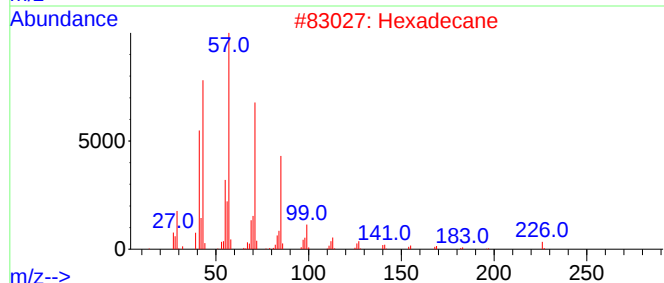
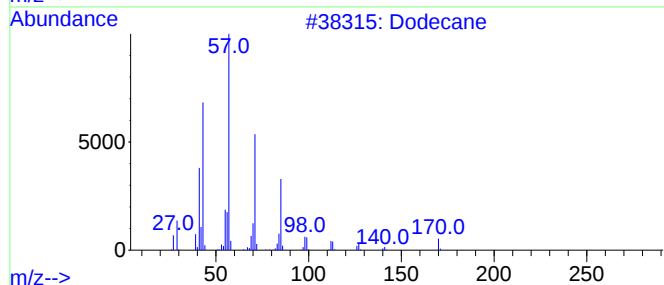
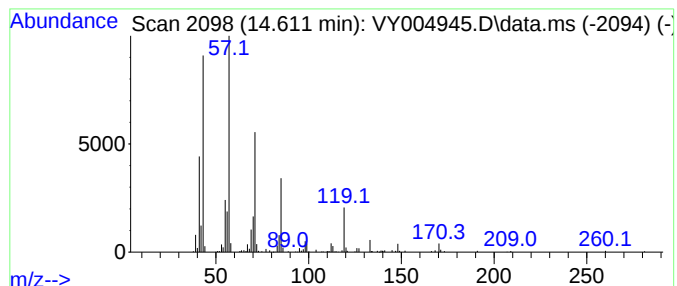
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Dodecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.611	6.37 ug/l	125690	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	76
2			Hexadecane	226	C16H34	000544-76-3	58
3			Tetradecane	198	C14H30	000629-59-4	58
4			Tridecane	184	C13H28	000629-50-5	58
5			Octadecane	254	C18H38	000593-45-3	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY060321\
Data File : VY004945.D
Acq On : 03 Jun 2021 18:32
Operator : SY/MD
Sample : M2589-15
Misc : 12.82G/5ML/MSVOA_Y/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
18-B11(2-4)

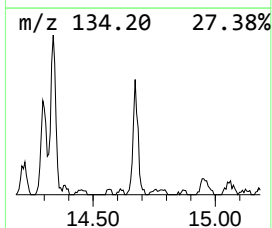
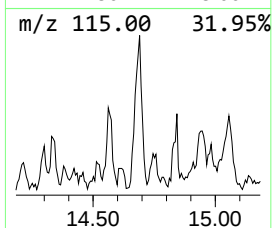
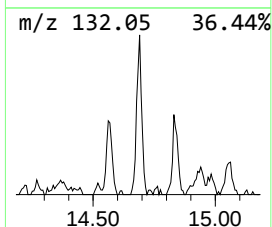
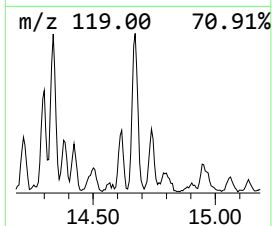
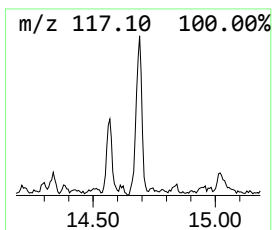
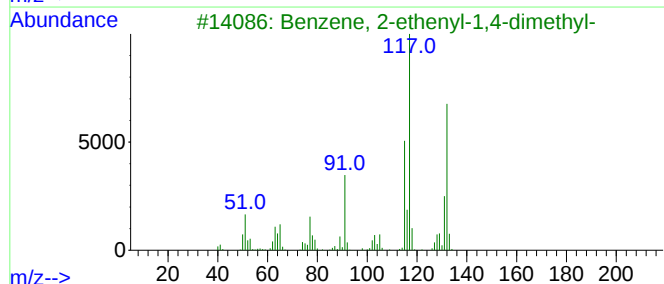
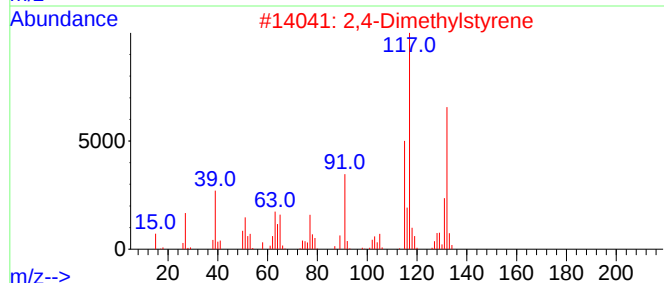
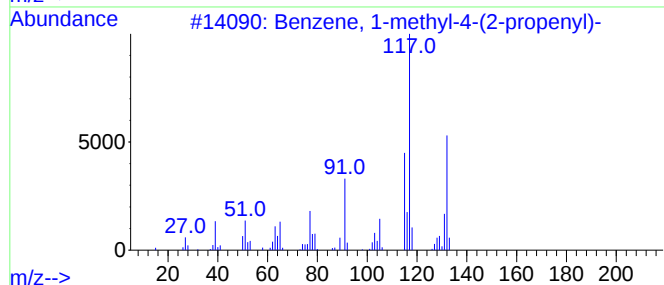
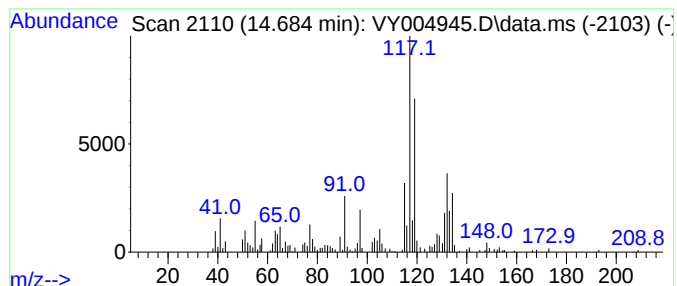
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-methyl-4-(2-prop... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	5.19 ug/l	102439	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	92
2			2,4-Dimethylstyrene	132	C10H12	002234-20-0	83
3			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	80
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	60
5			1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY060321\
Data File : VY004945.D
Acq On : 03 Jun 2021 18:32
Operator : SY/MD
Sample : M2589-15
Misc : 12.82G/5ML/MSVOA_Y/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
18-B11(2-4)

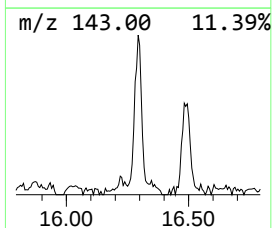
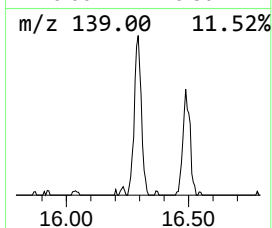
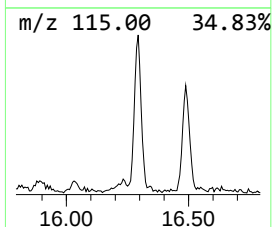
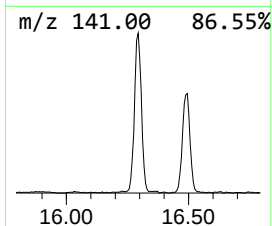
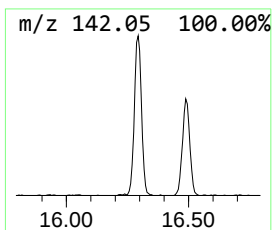
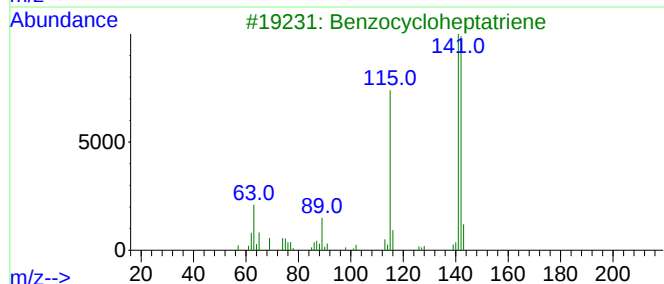
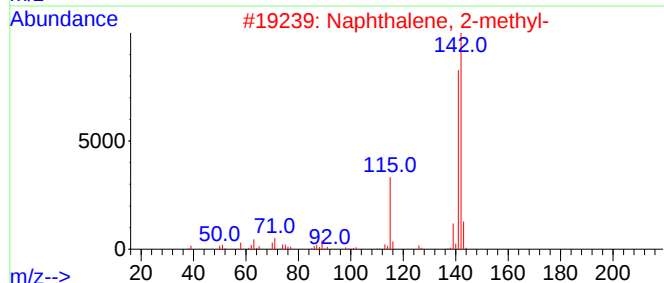
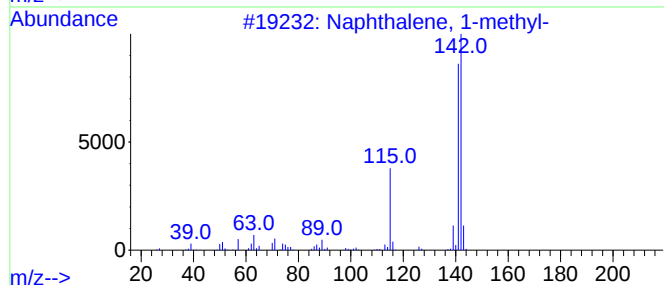
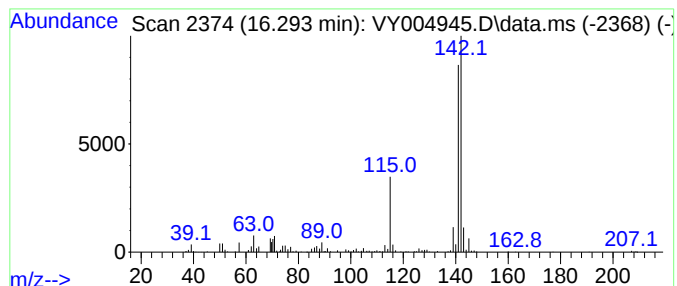
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.294	8.85 ug/l	174785	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
5			Naphthalene, 1-[(2-propenyloxy)m...	198	C14H14O	074685-39-5	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY060321\
Data File : VY004945.D
Acq On : 03 Jun 2021 18:32
Operator : SY/MD
Sample : M2589-15
Misc : 12.82G/5ML/MSVOA_Y/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
18-B11(2-4)

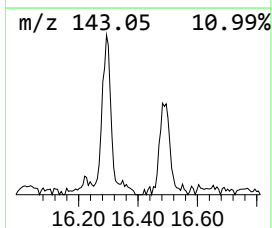
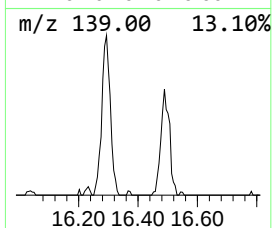
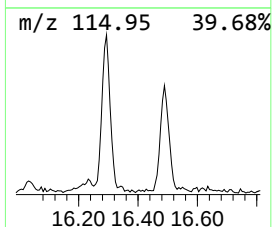
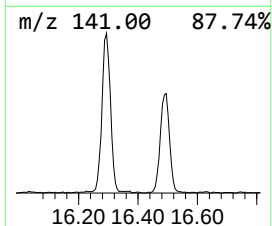
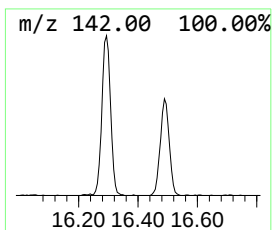
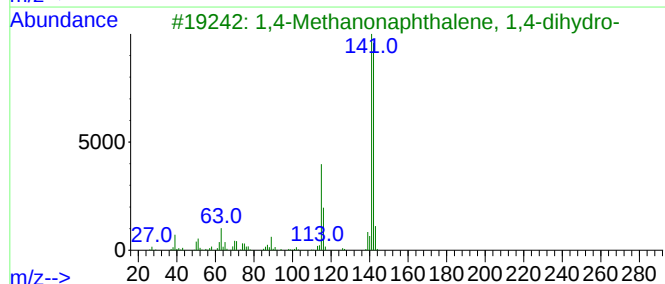
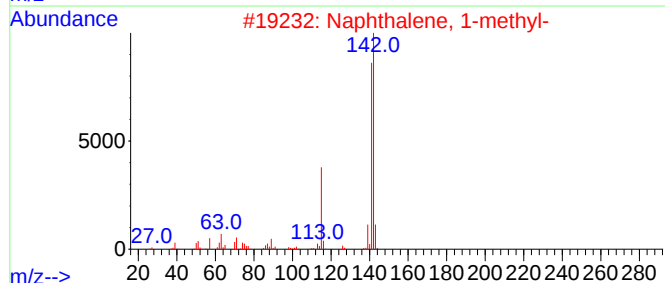
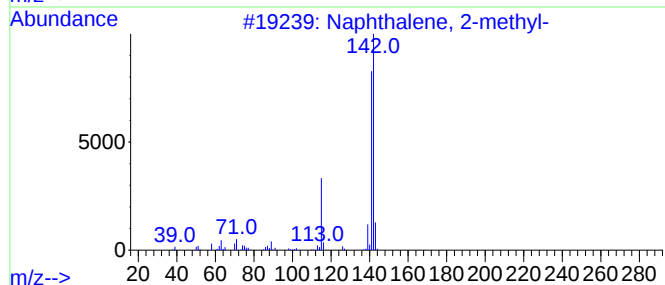
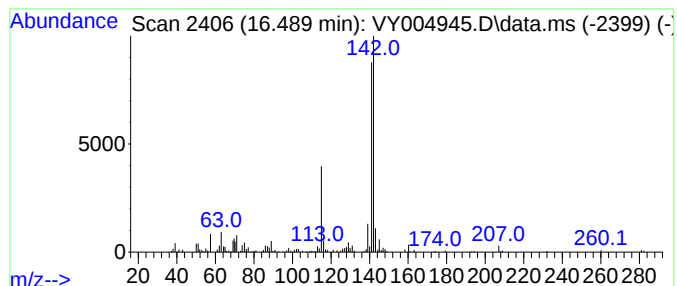
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1,4-Methanonaphthalene, 1,4... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.489	7.22 ug/l	142467	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
4			Benzocycloheptatriene	142	C11H10	000264-09-5	90
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY060321\
Data File : VY004945.D
Acq On : 03 Jun 2021 18:32
Operator : SY/MD
Sample : M2589-15
Misc : 12.82G/5ML/MSVOA_Y/SOIL
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
18-B11(2-4)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y060221S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(1R)-2,6,6-Trim...	12.337	7.9	ug/l	160942	3	11.496	1014890	50.0
Benzene, 1-ethy...	12.971	5.2	ug/l	103153	4	13.428	987021	50.0
Undecane	13.751	10.1	ug/l	199473	4	13.428	987021	50.0
Bicyclo[2.2.1]h...	14.312	5.7	ug/l	112114	4	13.428	987021	50.0
Dodecane	14.611	6.4	ug/l	125690	4	13.428	987021	50.0
Benzene, 1-meth...	14.684	5.2	ug/l	102439	4	13.428	987021	50.0
Naphthalene, 1-...	16.294	8.8	ug/l	174785	4	13.428	987021	50.0
1,4-Methanonaph...	16.489	7.2	ug/l	142467	4	13.428	987021	50.0