

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070121\
 Data File : VY005313.D
 Acq On : 01 Jul 2021 14:52
 Operator : SY/MD
 Sample : M2896-05
 Misc : 9.26G/5ML/MSVOA_Y/SOIL
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 FS-03

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\82Y062221S.M
 Title : SW846 8260

Signal : TIC: VY005313.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.869	3	8	15	rVB	18525	27477	1.34%	0.078%
2	3.960	341	351	364	rBV	76847	214098	10.46%	0.608%
3	4.722	467	476	495	rBV4	10395	35403	1.73%	0.101%
4	7.728	958	969	974	rBV	148462	344524	16.83%	0.978%
5	7.795	974	980	990	rVV	239338	547667	26.76%	1.555%
6	8.149	1029	1038	1049	rVB	151037	326608	15.96%	0.927%
7	8.551	1094	1104	1117	rVB	66204	133941	6.54%	0.380%
8	8.697	1117	1128	1140	rBV	386477	762409	37.25%	2.165%
9	9.356	1226	1236	1246	rVB3	15142	38612	1.89%	0.110%
10	9.807	1299	1310	1317	rBV4	51894	127607	6.24%	0.362%
11	9.959	1326	1335	1338	rBV4	8232	24316	1.19%	0.069%
12	10.185	1364	1372	1379	rBV	639250	1097553	53.63%	3.117%
13	10.246	1379	1382	1386	rVV2	13572	25072	1.23%	0.071%
14	10.295	1386	1390	1394	rVV4	13036	21392	1.05%	0.061%
15	10.374	1394	1403	1409	rVV	389873	629914	30.78%	1.789%
16	10.435	1409	1413	1417	rBV	20557	29850	1.46%	0.085%
17	10.605	1436	1441	1452	rVB2	42342	94826	4.63%	0.269%
18	10.697	1452	1456	1463	rVB3	11041	22731	1.11%	0.065%
19	10.801	1463	1473	1478	rBV5	16845	52623	2.57%	0.149%
20	10.886	1478	1487	1491	rBV3	21762	49035	2.40%	0.139%
21	10.965	1495	1500	1503	rBV	26705	43195	2.11%	0.123%
22	11.002	1503	1506	1510	rVV	35405	56536	2.76%	0.161%
23	11.038	1510	1512	1517	rVB2	20940	24920	1.22%	0.071%
24	11.160	1524	1532	1536	rBV4	51758	101885	4.98%	0.289%
25	11.209	1536	1540	1547	rVB2	43827	78616	3.84%	0.223%
26	11.294	1547	1554	1564	rBV8	39469	121481	5.94%	0.345%
27	11.386	1564	1569	1571	rBV	18214	33319	1.63%	0.095%
28	11.496	1577	1587	1595	rVB	618505	1031686	50.41%	2.930%
29	11.599	1598	1604	1608	rBV2	70070	140827	6.88%	0.400%
30	11.715	1611	1623	1637	rVB2	743911	1381746	67.52%	3.924%
31	11.831	1637	1642	1647	rVB3	34109	58483	2.86%	0.166%
32	11.892	1647	1652	1657	rBV2	95784	160806	7.86%	0.457%
33	11.947	1657	1661	1666	rVB	41932	66167	3.23%	0.188%
34	12.032	1666	1675	1683	rBV4	93190	251433	12.29%	0.714%
35	12.105	1683	1687	1694	rVB3	59863	104341	5.10%	0.296%
36	12.197	1697	1702	1704	rBV3	85710	128960	6.30%	0.366%

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37	12.239	1704	1709	1717	rVV	594680	960235	46.92%	2.727%
38	12.319	1717	1722	1728	rVV3	166413	278826	13.62%	0.792%
39	12.379	1728	1732	1740	rVB	86758	148373	7.25%	0.421%
40	12.489	1743	1750	1759	rVB2	1128018	2046534	100.00%	5.812%
41	12.568	1759	1763	1767	rBV4	39531	65003	3.18%	0.185%
42	12.629	1770	1773	1776	rVB	28863	33045	1.61%	0.094%
43	12.672	1776	1780	1786	rVB	160141	229189	11.20%	0.651%
44	12.733	1786	1790	1793	rBV	220462	346086	16.91%	0.983%
45	12.770	1793	1796	1797	rVV	168711	225583	11.02%	0.641%
46	12.806	1797	1802	1809	rVB3	745452	1256950	61.42%	3.569%
47	12.977	1821	1830	1837	rVB	209593	432430	21.13%	1.228%
48	13.044	1837	1841	1844	rBV4	47432	83770	4.09%	0.238%
49	13.123	1849	1854	1860	rVB	389379	619823	30.29%	1.760%
50	13.263	1868	1877	1881	rBV6	46112	158479	7.74%	0.450%
51	13.318	1881	1886	1891	rBV3	117442	215102	10.51%	0.611%
52	13.373	1891	1895	1897	rBV	68333	98488	4.81%	0.280%
53	13.428	1899	1904	1908	rBV	527066	758535	37.06%	2.154%
54	13.501	1913	1916	1918	rBV	21882	30614	1.50%	0.087%
55	13.593	1926	1931	1933	rBV3	132597	220378	10.77%	0.626%
56	13.629	1933	1937	1941	rVV3	390048	774845	37.86%	2.200%
57	13.672	1941	1944	1946	rVV3	326271	509002	24.87%	1.445%
58	13.696	1946	1948	1954	rVV2	337306	564431	27.58%	1.603%
59	13.751	1954	1957	1963	rVB	428734	594834	29.07%	1.689%
60	13.824	1963	1969	1974	rBV	207629	330765	16.16%	0.939%
61	13.885	1975	1979	1981	rVV	120333	170153	8.31%	0.483%
62	13.916	1981	1984	1989	rVB	202170	274644	13.42%	0.780%
63	13.971	1989	1993	1998	rVB	304134	448721	21.93%	1.274%
64	14.032	1999	2003	2006	rBV	57940	76582	3.74%	0.217%
65	14.080	2006	2011	2016	rBV	284512	453394	22.15%	1.288%
66	14.166	2023	2025	2029	rVB2	39048	42305	2.07%	0.120%
67	14.215	2029	2033	2037	rVB2	69778	95239	4.65%	0.270%
68	14.269	2037	2042	2045	rBV	126042	215080	10.51%	0.611%
69	14.336	2050	2053	2056	rBV	98292	132028	6.45%	0.375%
70	14.385	2056	2061	2065	rVB2	292539	471069	23.02%	1.338%
71	14.428	2065	2068	2074	rVB5	75993	141892	6.93%	0.403%
72	14.483	2074	2077	2079	rBV3	23013	32022	1.56%	0.091%
73	14.519	2080	2083	2086	rVB	54317	65414	3.20%	0.186%
74	14.593	2086	2095	2096	rBV2	359292	794494	38.82%	2.256%
75	14.690	2103	2111	2116	rBV3	340917	936936	45.78%	2.661%
76	14.751	2116	2121	2130	rVB3	233945	435637	21.29%	1.237%

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77	14.836	2130	2135	2140	rVB	159814	224787	10.98%	0.638%
78	14.903	2142	2146	2148	rBV2	72753	97385	4.76%	0.277%
79	14.946	2148	2153	2157	rBV3	201477	324074	15.84%	0.920%
80	15.056	2166	2171	2178	rVB2	124456	281438	13.75%	0.799%
81	15.117	2178	2181	2184	rBV3	21732	27137	1.33%	0.077%
82	15.196	2184	2194	2209	rBV3	466847	1854138	90.60%	5.265%
83	15.318	2209	2214	2224	rVV	803268	1566809	76.56%	4.449%
84	15.403	2224	2228	2231	rVV	227218	364513	17.81%	1.035%
85	15.440	2231	2234	2242	rVB4	244235	532853	26.04%	1.513%
86	15.531	2242	2249	2256	rBV4	131575	352190	17.21%	1.000%
87	15.611	2257	2262	2266	rBV3	70031	137472	6.72%	0.390%
88	15.665	2267	2271	2276	rBV3	74564	155892	7.62%	0.443%
89	15.763	2283	2287	2292	rVB	121100	206885	10.11%	0.587%
90	15.861	2292	2303	2314	rVB5	233782	1027536	50.21%	2.918%
91	15.964	2315	2320	2322	rBV2	59152	102562	5.01%	0.291%
92	16.001	2322	2326	2329	rBV	54679	90189	4.41%	0.256%
93	16.135	2342	2348	2354	rVV	197048	409558	20.01%	1.163%
94	16.202	2354	2359	2366	rVV	293861	701436	34.27%	1.992%
95	16.287	2366	2373	2380	rVB2	391053	948287	46.34%	2.693%
96	16.360	2380	2385	2390	rVB	90135	164996	8.06%	0.469%
97	16.434	2390	2397	2401	rBV2	223976	507450	24.80%	1.441%
98	16.489	2401	2406	2416	rVB4	284206	689589	33.70%	1.958%
99	16.696	2434	2440	2446	rVB	101982	225144	11.00%	0.639%
100	16.763	2446	2451	2455	rBV3	30380	65594	3.21%	0.186%

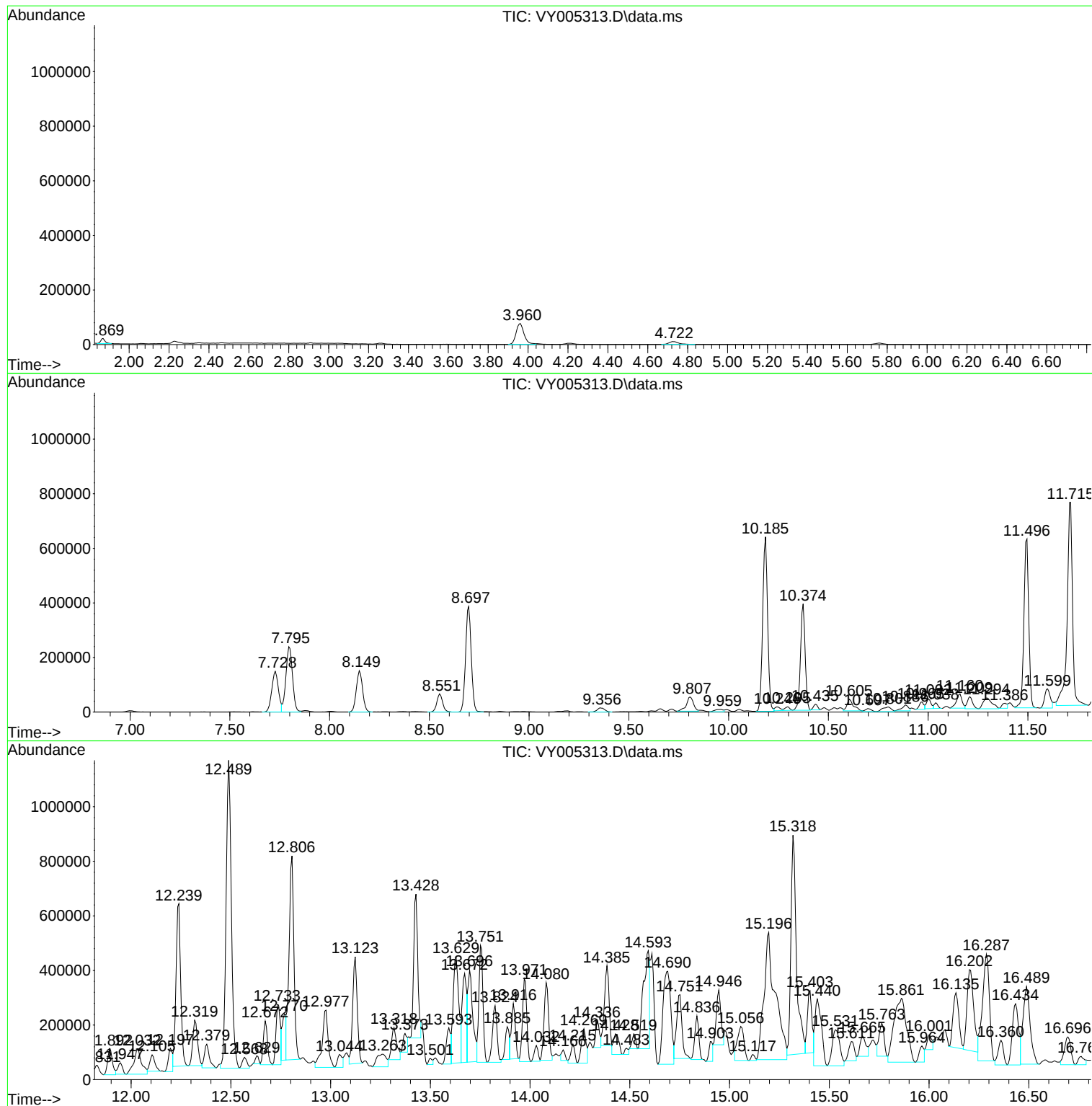
Sum of corrected areas: 35214903

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Instrument :
MSVOA_Y
ClientSampled :
FS-03

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062221S.M
Quant Title : SW846 8260

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



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Instrument :
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ClientSampleId :
FS-03

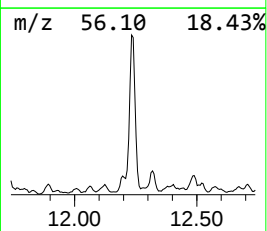
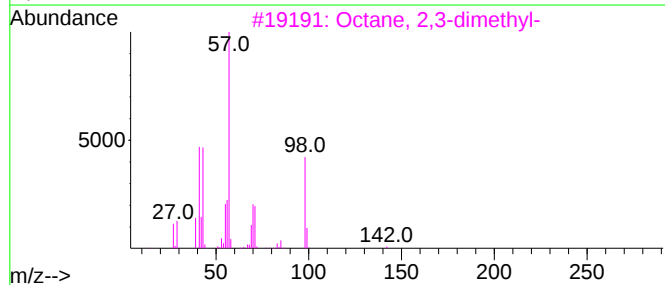
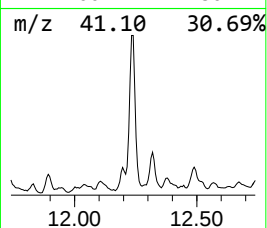
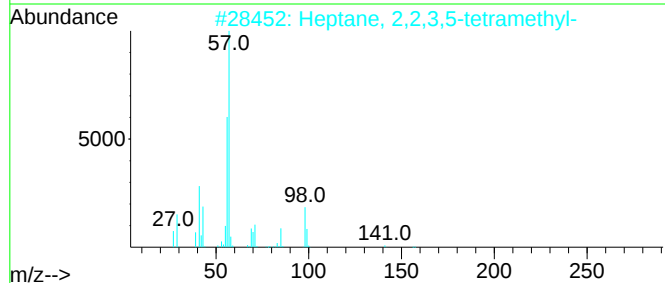
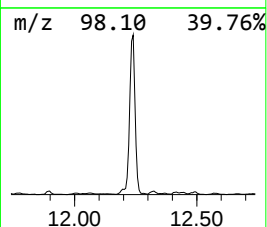
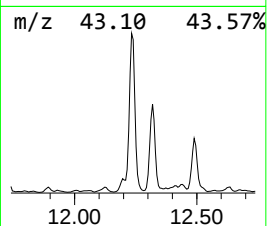
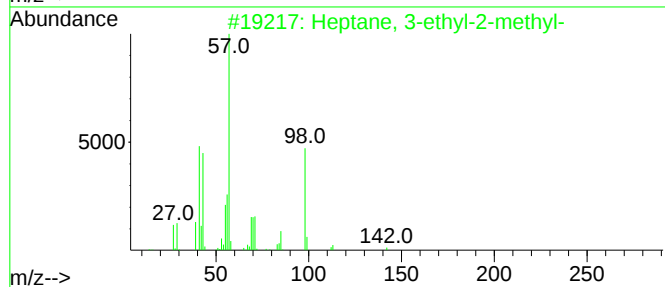
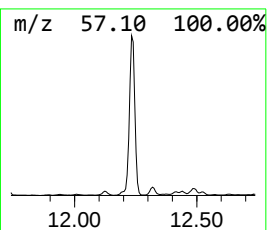
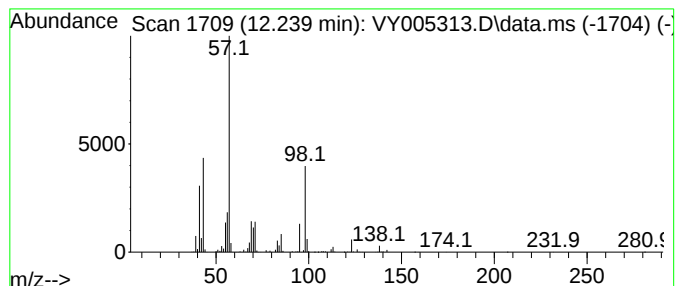
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062221S.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane, 3-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.239	46.54 ug/l	960235	Chlorobenzene-d5	11.496

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	81
2			Heptane, 2,2,3,5-tetramethyl-	156	C11H24	061868-42-6	53
3			Octane, 2,3-dimethyl-	142	C10H22	007146-60-3	50
4			Hexane, 3-ethyl-2,5-dimethyl-	142	C10H22	052897-04-8	47
5			Heptane, 4-propyl-	142	C10H22	003178-29-8	38



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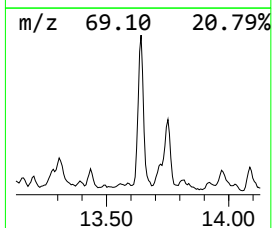
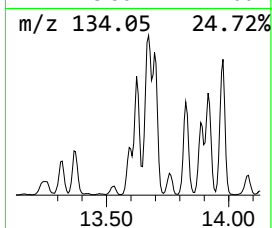
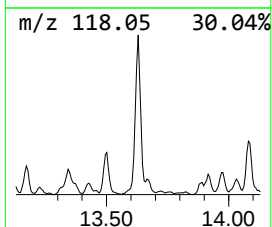
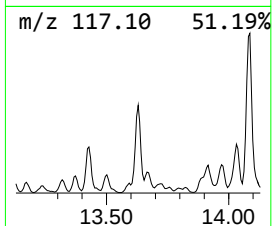
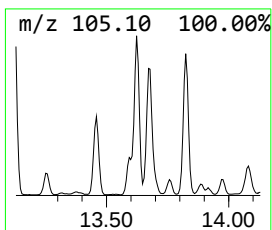
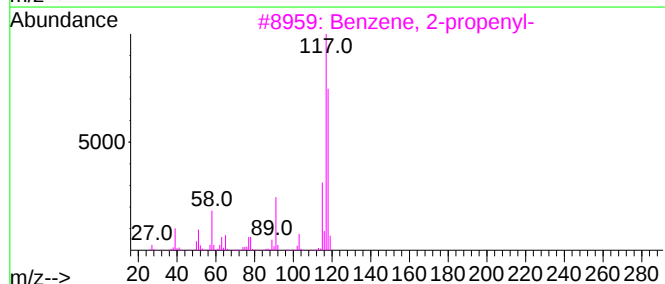
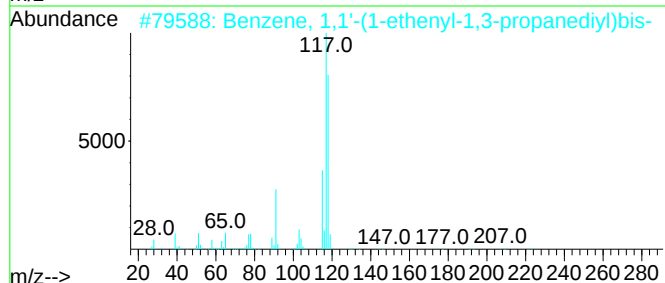
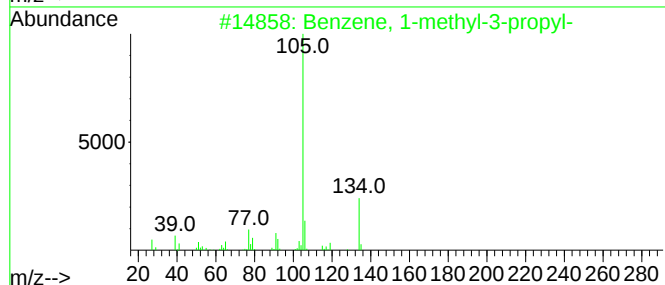
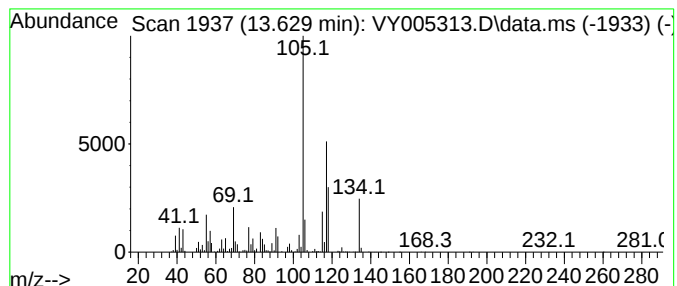
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown13.629 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.629	51.08 ug/l	774845	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	46
2			Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	43
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	43
4			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	43
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	43



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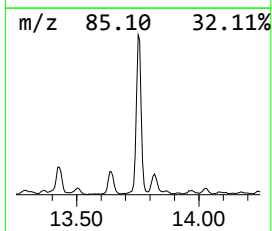
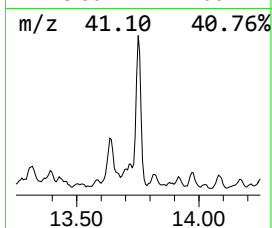
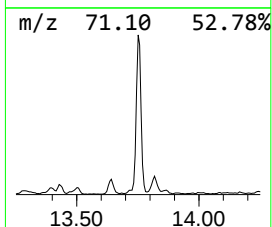
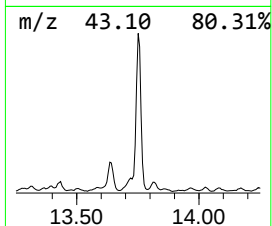
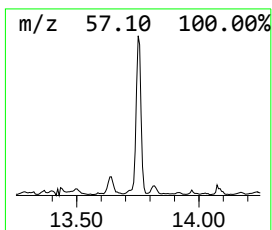
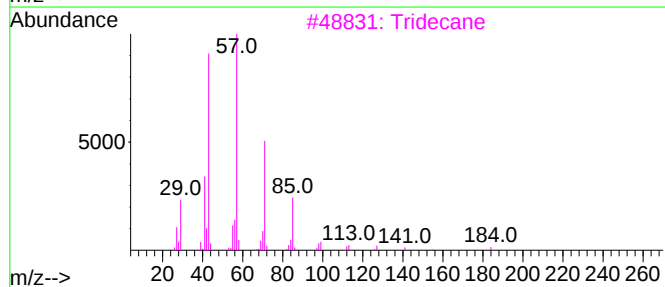
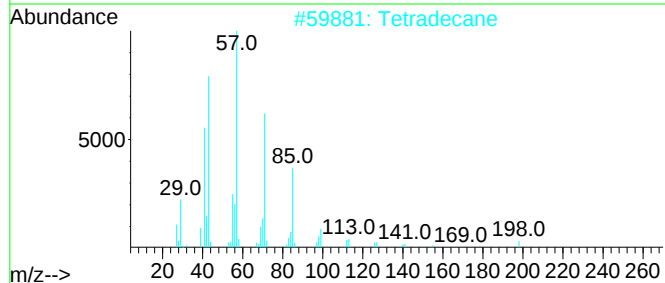
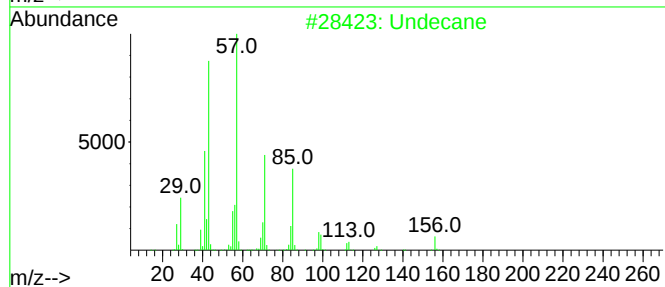
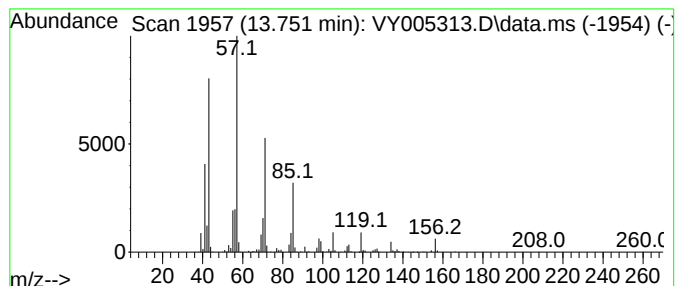
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Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 3 Undecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.751	39.21 ug/l	594834	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	91
2	Tetradecane			198	C14H30	000629-59-4	64
3	Tridecane			184	C13H28	000629-50-5	59
4	Hexadecane			226	C16H34	000544-76-3	58
5	Decane			142	C10H22	000124-18-5	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070121\
Data File : VY005313.D
Acq On : 01 Jul 2021 14:52
Operator : SY/MD
Sample : M2896-05
Misc : 9.26G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FS-03

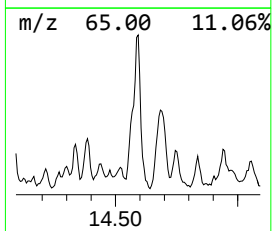
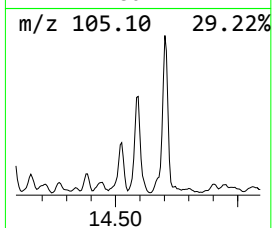
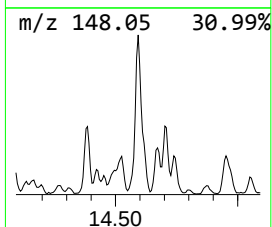
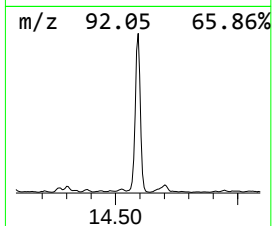
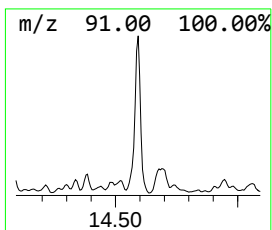
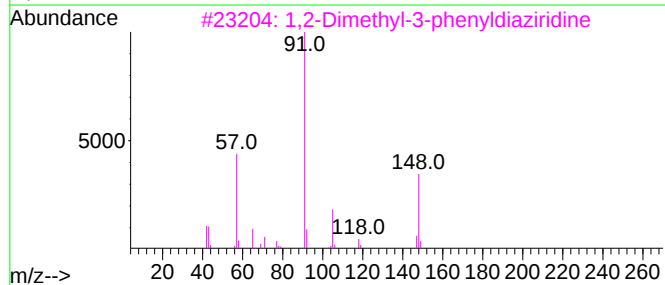
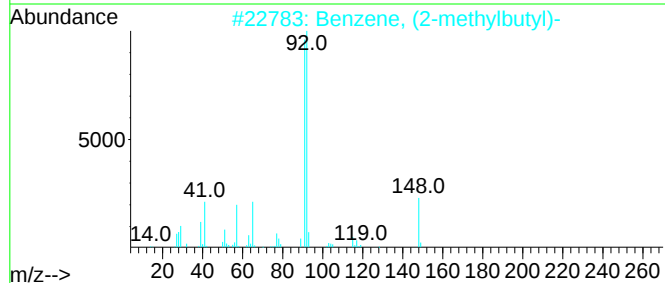
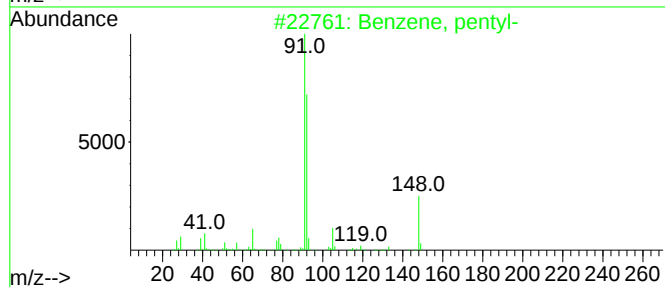
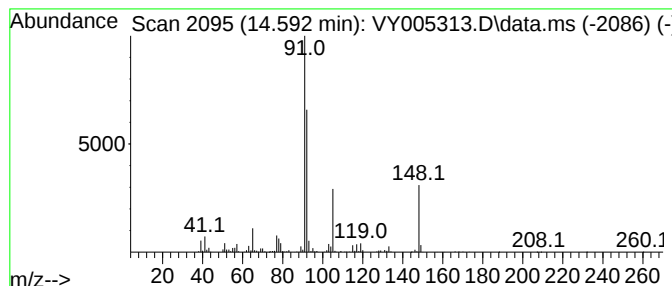
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, pentyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.592	52.37 ug/l	794494	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, pentyl-	148	C11H16	000538-68-1	87
2			Benzene, (2-methylbutyl)-	148	C11H16	003968-85-2	53
3			1,2-Dimethyl-3-phenyldiaziridine	148	C9H12N2	051456-63-4	49
4			Benzene, hexyl-	162	C12H18	001077-16-3	43
5			Benzene, (3-methylbutyl)-	148	C11H16	002049-94-7	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070121\
Data File : VY005313.D
Acq On : 01 Jul 2021 14:52
Operator : SY/MD
Sample : M2896-05
Misc : 9.26G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FS-03

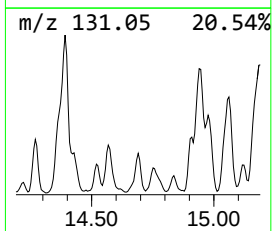
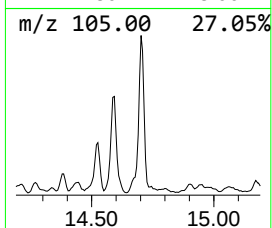
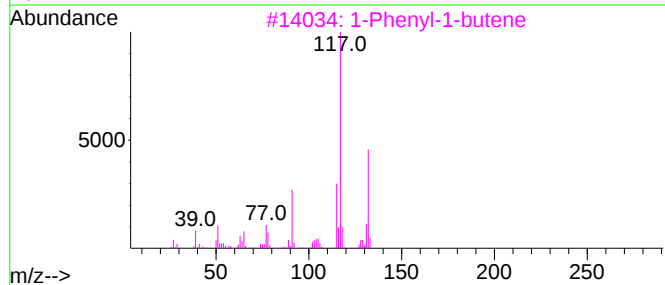
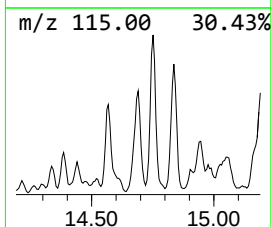
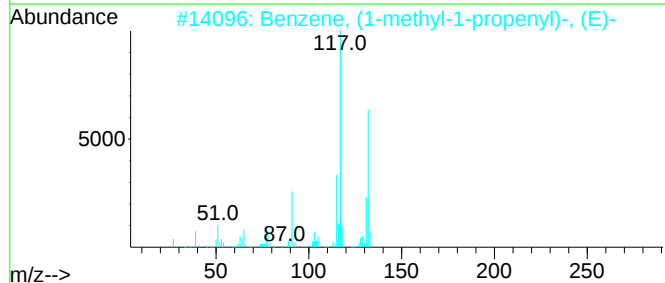
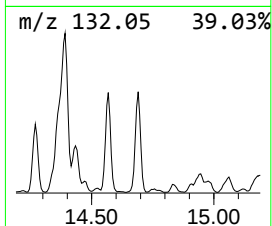
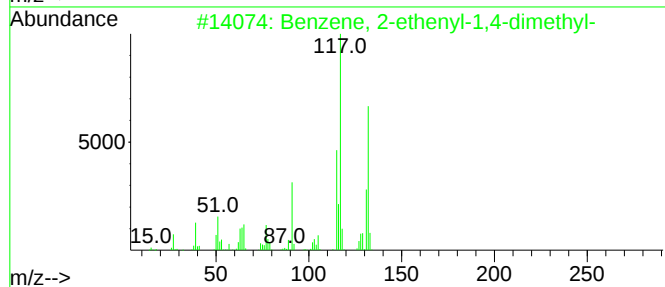
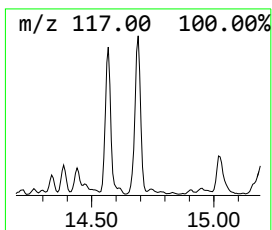
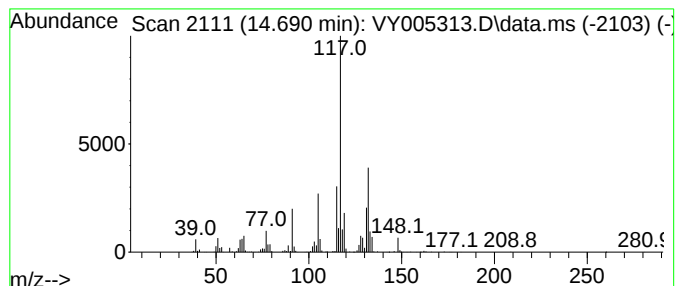
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.690	61.76 ug/l	936936	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
2			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	76
4			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	76
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070121\
Data File : VY005313.D
Acq On : 01 Jul 2021 14:52
Operator : SY/MD
Sample : M2896-05
Misc : 9.26G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FS-03

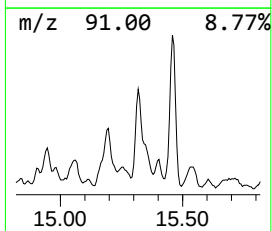
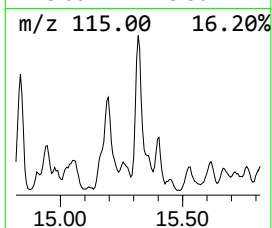
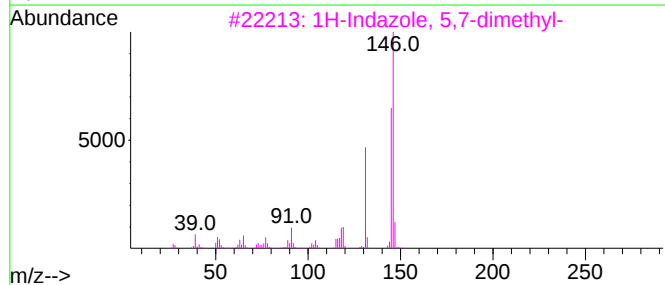
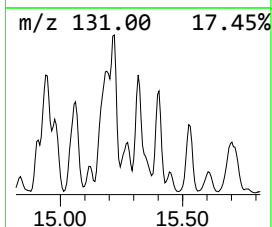
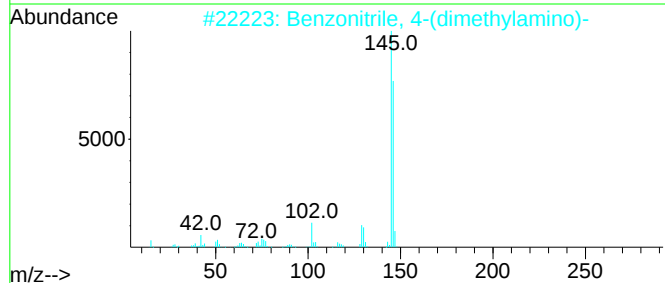
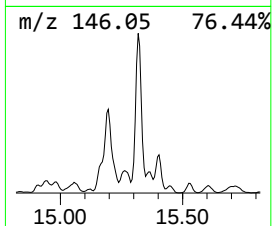
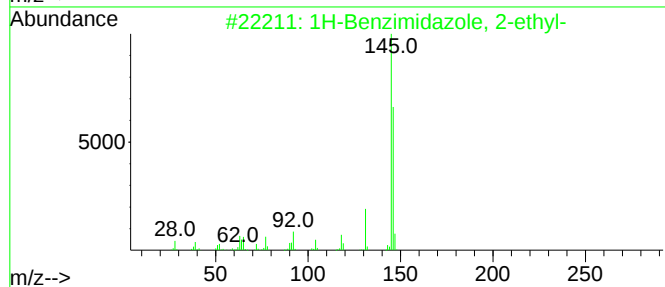
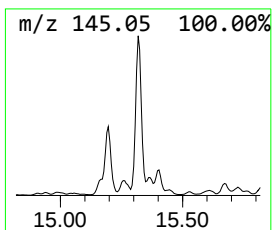
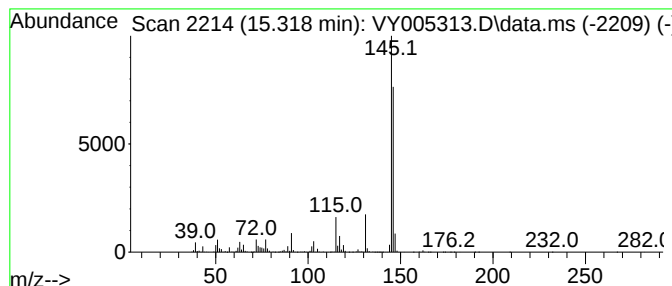
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Quant Title : SW846 8260

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

Peak Number 6 1H-Benzimidazole, 2-ethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.318	103.28 ug/l	1566810	1,4-Dichlorobenzene-d4	13.428

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Benzimidazole, 2-ethyl-	146	C9H10N2	001848-84-6	80
2		Benzonitrile, 4-(dimethylamino)-	146	C9H10N2	001197-19-9	64
3		1H-Indazole, 5,7-dimethyl-	146	C9H10N2	043067-41-0	58
4		1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	50
5		1,2-Dimethylbenzimidazole	146	C9H10N2	002876-08-6	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070121\
Data File : VY005313.D
Acq On : 01 Jul 2021 14:52
Operator : SY/MD
Sample : M2896-05
Misc : 9.26G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FS-03

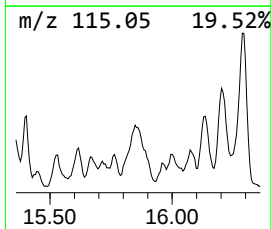
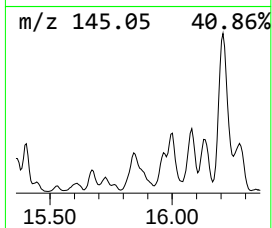
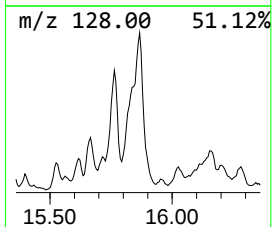
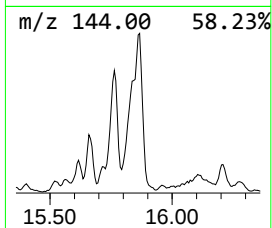
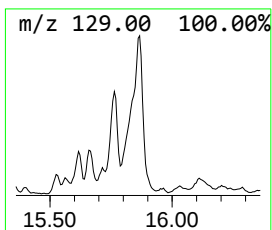
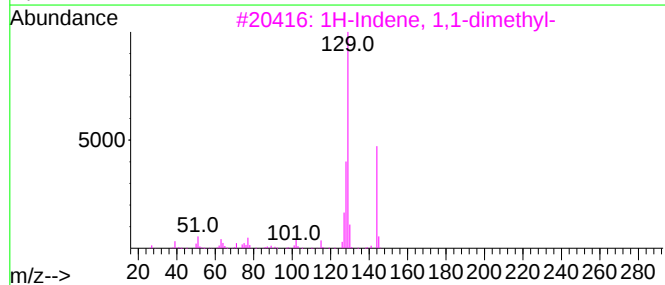
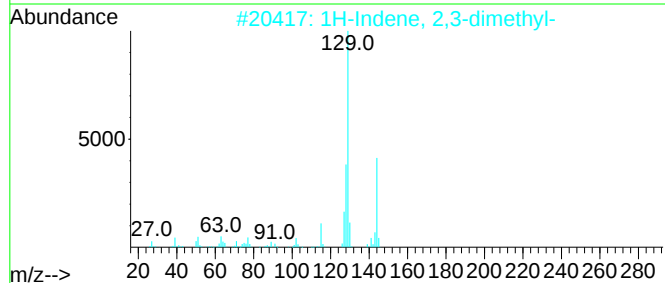
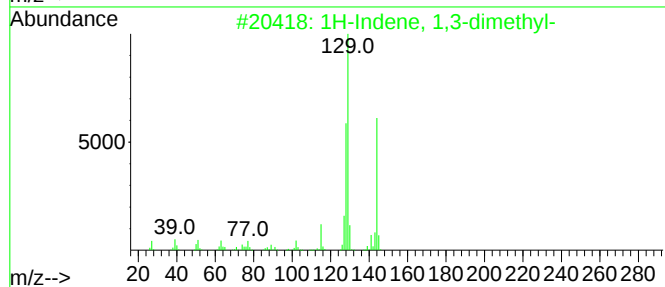
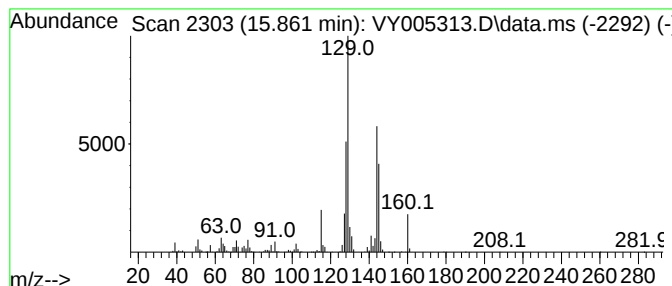
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Quant Title : SW846 8260

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

Peak Number 7 1H-Indene, 1,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.861	67.73 ug/l	1027540	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 1,3-dimethyl-	144	C11H12	002177-48-2	93
2			1H-Indene, 2,3-dimethyl-	144	C11H12	004773-82-4	76
3			1H-Indene, 1,1-dimethyl-	144	C11H12	018636-55-0	70
4			Naphthalene, 1,2-dihydro-3-methyl-	144	C11H12	002717-44-4	68
5			1H-Indene, 4,7-dimethyl-	144	C11H12	006974-97-6	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070121\
Data File : VY005313.D
Acq On : 01 Jul 2021 14:52
Operator : SY/MD
Sample : M2896-05
Misc : 9.26G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FS-03

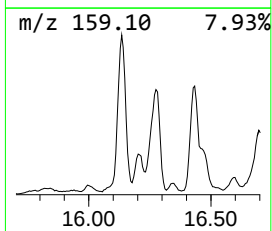
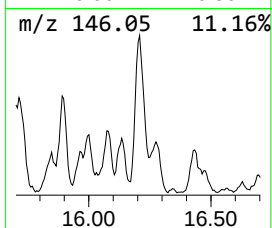
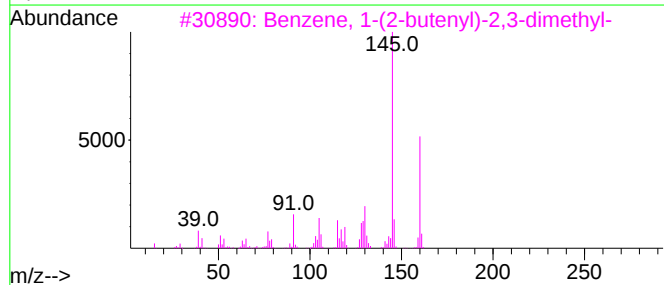
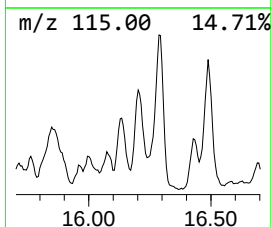
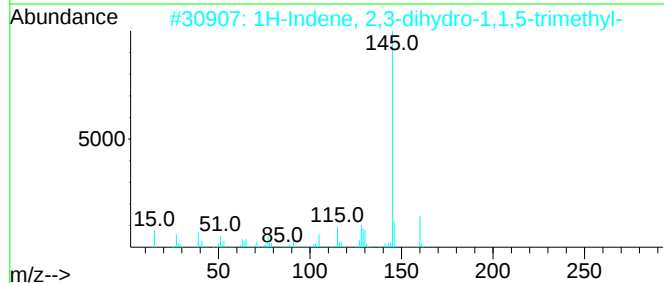
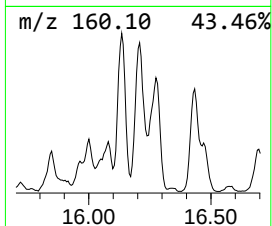
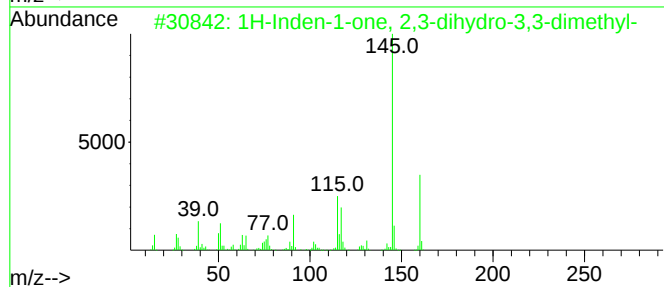
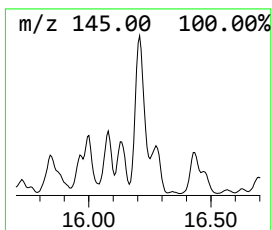
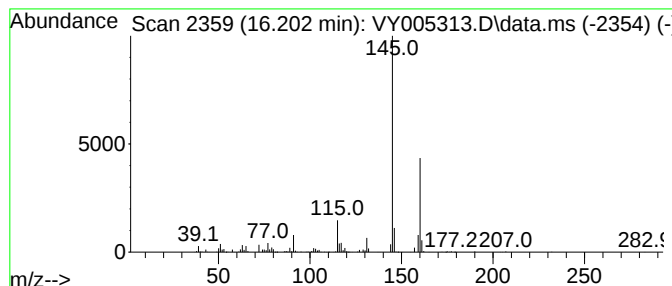
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Quant Title : SW846 8260

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

Peak Number 8 1H-Inden-1-one, 2,3-dihydro... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.202	46.24 ug/l	701436	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Inden-1-one, 2,3-dihydro-3,3-...	160	C11H12O	026465-81-6	90
2			1H-Indene, 2,3-dihydro-1,1,5-tri...	160	C12H16	040650-41-7	86
3			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	86
4			1H-Benzimidazole, 2-(1-methyleth...	160	C10H12N2	005851-43-4	80
5			Silane, methyltripropynyl-	160	C10H12Si	1000158-62-1	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070121\
Data File : VY005313.D
Acq On : 01 Jul 2021 14:52
Operator : SY/MD
Sample : M2896-05
Misc : 9.26G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FS-03

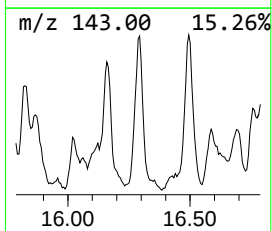
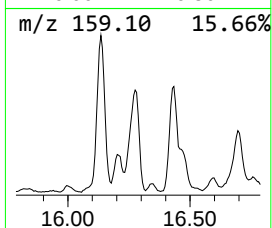
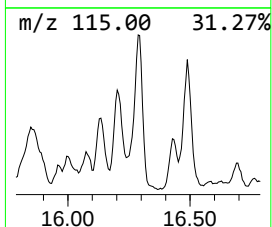
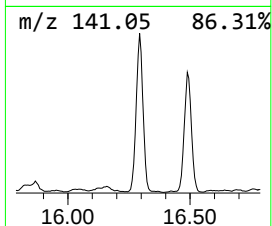
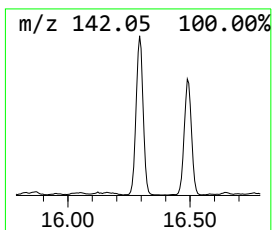
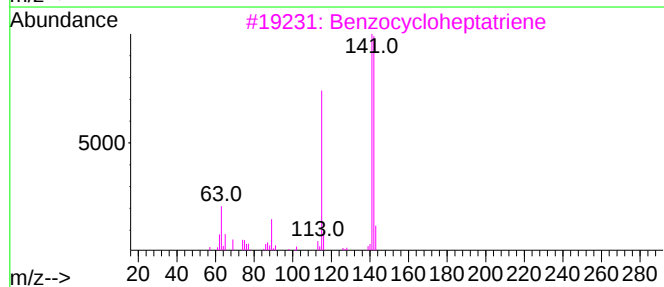
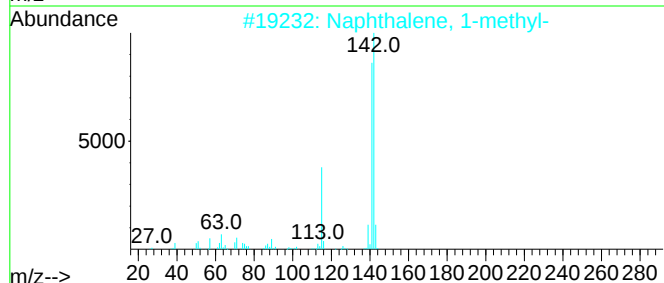
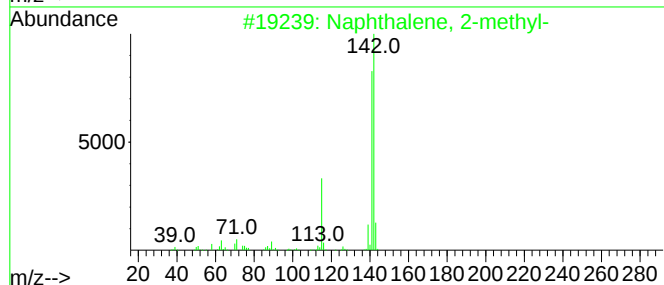
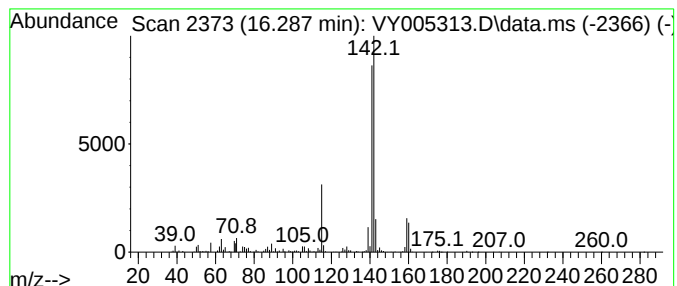
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Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.287	62.51 ug/l	948287	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
3			Benzocycloheptatriene	142	C11H10	000264-09-5	87
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	83
5			Naphthalene, 1-(2-hydroxypropyl)	186	C13H14O	027653-13-0	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070121\
Data File : VY005313.D
Acq On : 01 Jul 2021 14:52
Operator : SY/MD
Sample : M2896-05
Misc : 9.26G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FS-03

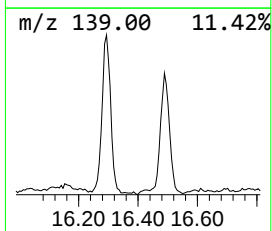
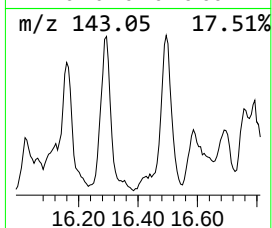
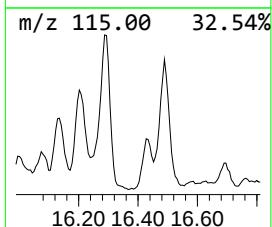
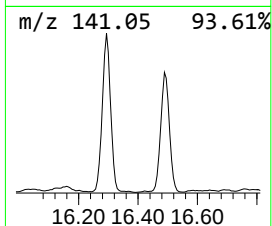
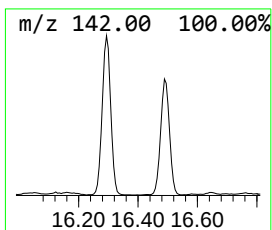
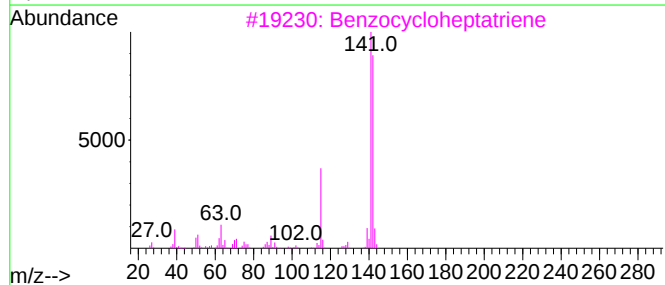
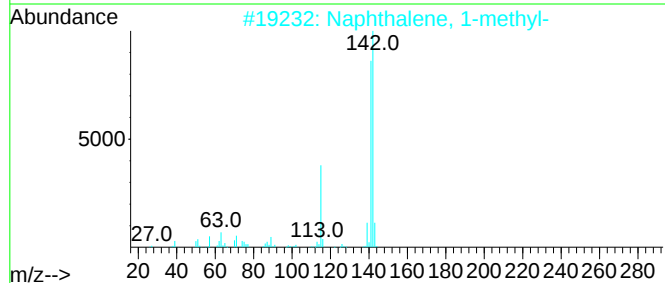
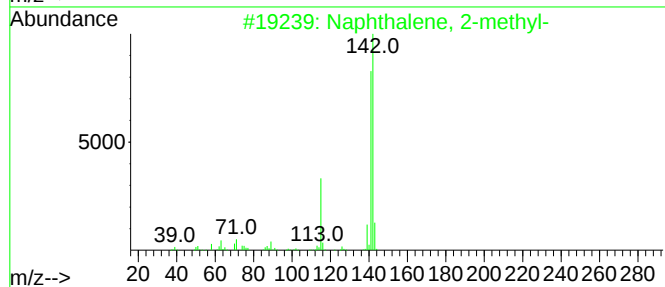
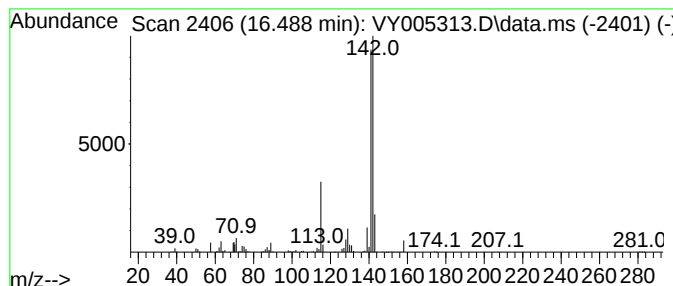
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Quant Title : SW846 8260

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

Peak Number 10 Benzocycloheptatriene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.488	45.46 ug/l	689589	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	95
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	93
3			Benzocycloheptatriene	142	C11H10	000264-09-5	90
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	87
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY070121\
Data File : VY005313.D
Acq On : 01 Jul 2021 14:52
Operator : SY/MD
Sample : M2896-05
Misc : 9.26G/5ML/MSVOA_Y/SOIL
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
FS-03

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y062221S.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
Heptane, 3-ethy...	12.239	46.5	ug/l	960235	3	11.496	1031690	50.0
unknown13.629	13.629	51.1	ug/l	774845	4	13.428	758535	50.0
Undecane	13.751	39.2	ug/l	594834	4	13.428	758535	50.0
Benzene, pentyl-	14.592	52.4	ug/l	794494	4	13.428	758535	50.0
Benzene, 2-ethe...	14.690	61.8	ug/l	936936	4	13.428	758535	50.0
1H-Benzimidazol...	15.318	103.3	ug/l	1566810	4	13.428	758535	50.0
1H-Indene, 1,3-...	15.861	67.7	ug/l	1027540	4	13.428	758535	50.0
1H-Inden-1-one,...	16.202	46.2	ug/l	701436	4	13.428	758535	50.0
Naphthalene, 1-...	16.287	62.5	ug/l	948287	4	13.428	758535	50.0
Benzocyclohepta...	16.488	45.5	ug/l	689589	4	13.428	758535	50.0