

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033122\
 Data File : VY008143.D
 Acq On : 31 Mar 2022 17:06
 Operator : SY/MD
 Sample : N2123-02
 Misc : 4.49g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 20220339-AMENDED-COMP-CS-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: VY008143.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.716	958	967	972	rBV2	63360	164361	1.19%	0.159%
2	7.789	972	979	987	rVB	196379	496561	3.59%	0.481%
3	8.002	1003	1014	1021	rBV	69710	171687	1.24%	0.166%
4	8.142	1029	1037	1049	rVB	97159	215307	1.55%	0.209%
5	8.545	1094	1103	1114	rBV	84297	171341	1.24%	0.166%
6	8.691	1118	1127	1141	rBV	290180	585673	4.23%	0.567%
7	9.185	1192	1208	1214	rBV3	172451	501831	3.62%	0.486%
8	9.465	1244	1254	1261	rVB	105853	227089	1.64%	0.220%
9	9.606	1270	1277	1288	rVB3	110772	211014	1.52%	0.204%
10	9.819	1306	1312	1316	rBV	223495	408917	2.95%	0.396%
11	9.959	1324	1335	1346	rBV4	157890	449650	3.25%	0.436%
12	10.179	1364	1371	1386	rBV	690060	1444566	10.43%	1.399%
13	10.368	1394	1402	1409	rVB5	278720	648567	4.68%	0.628%
14	10.520	1423	1427	1435	rVB	171675	308391	2.23%	0.299%
15	10.618	1435	1443	1449	rBV3	127946	311951	2.25%	0.302%
16	10.801	1468	1473	1479	rVB	308503	483107	3.49%	0.468%
17	10.904	1479	1490	1497	rBV3	130685	390618	2.82%	0.378%
18	10.971	1497	1501	1504	rVV2	100188	175117	1.26%	0.170%
19	11.038	1508	1512	1515	rVV	216179	341765	2.47%	0.331%
20	11.093	1515	1521	1526	rVV	506867	960498	6.94%	0.930%
21	11.148	1526	1530	1533	rVV2	112794	182644	1.32%	0.177%
22	11.197	1533	1538	1543	rVV3	183269	364919	2.64%	0.353%
23	11.288	1543	1553	1563	rVB5	288359	935621	6.76%	0.906%
24	11.380	1563	1568	1581	rBV	307227	647026	4.67%	0.627%
25	11.489	1581	1586	1591	rBV	392480	640217	4.62%	0.620%
26	11.593	1597	1603	1607	rBV	612286	998224	7.21%	0.967%
27	11.697	1612	1620	1624	rVV3	488736	1271683	9.18%	1.232%
28	11.739	1624	1627	1631	rVV	292485	545332	3.94%	0.528%
29	11.782	1631	1634	1639	rVB	225990	347216	2.51%	0.336%
30	11.843	1639	1644	1648	rBV3	105435	204985	1.48%	0.199%
31	12.026	1663	1674	1682	rBV3	460879	1505144	10.87%	1.458%
32	12.123	1686	1690	1694	rVB	482071	658464	4.76%	0.638%
33	12.203	1698	1703	1706	rBV3	289698	495877	3.58%	0.480%
34	12.239	1706	1709	1715	rVB2	374781	650224	4.70%	0.630%
35	12.331	1719	1724	1728	rBV	347299	554765	4.01%	0.537%
36	12.379	1728	1732	1735	rBV5	105815	183805	1.33%	0.178%

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Stop Thrs : 0

Peak Location: TOP

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37	12.483	1745	1749	1753	rBV2	334106	509654	3.68%	0.494%
38	12.648	1771	1776	1784	rVB2	405970	865180	6.25%	0.838%
39	12.733	1784	1790	1793	rBV	848354	1466525	10.59%	1.420%
40	12.770	1793	1796	1799	rVV	656021	1100854	7.95%	1.066%
41	12.812	1799	1803	1809	rVV5	684005	1402423	10.13%	1.358%
42	12.861	1809	1811	1816	rVB2	196822	251867	1.82%	0.244%
43	12.977	1822	1830	1837	rVV2	416534	1094007	7.90%	1.060%
44	13.044	1837	1841	1847	rVB	525018	767657	5.54%	0.744%
45	13.123	1848	1854	1859	rVV	1419770	2188944	15.81%	2.120%
46	13.172	1859	1862	1867	rVB2	281677	392884	2.84%	0.381%
47	13.227	1867	1871	1873	rBV2	156356	232046	1.68%	0.225%
48	13.318	1879	1886	1890	rBV2	634026	1218965	8.80%	1.181%
49	13.367	1890	1894	1899	rVV2	675838	1109868	8.02%	1.075%
50	13.428	1901	1904	1906	rVV	308192	420277	3.04%	0.407%
51	13.459	1906	1909	1913	rVB	597777	830789	6.00%	0.805%
52	13.501	1913	1916	1921	rVB4	142472	185044	1.34%	0.179%
53	13.593	1923	1931	1933	rBV2	343800	687455	4.96%	0.666%
54	13.629	1933	1937	1941	rVV2	835747	1126749	8.14%	1.091%
55	13.672	1941	1944	1952	rVB2	462894	761479	5.50%	0.738%
56	13.751	1952	1957	1962	rBV3	605921	1023641	7.39%	0.992%
57	13.806	1962	1966	1972	rBV2	902902	1522049	10.99%	1.474%
58	13.885	1976	1979	1981	rBV	212877	287300	2.07%	0.278%
59	13.916	1981	1984	1989	rVB	477983	688830	4.97%	0.667%
60	13.977	1989	1994	1998	rVB	1154934	1663923	12.02%	1.612%
61	14.019	1998	2001	2005	rVB2	271787	361840	2.61%	0.350%
62	14.080	2005	2011	2015	rBV3	764847	1408142	10.17%	1.364%
63	14.215	2027	2033	2036	rBV3	186152	422937	3.05%	0.410%
64	14.263	2037	2041	2044	rVV4	357084	761550	5.50%	0.738%
65	14.300	2044	2047	2050	rVV	378685	629228	4.54%	0.609%
66	14.336	2050	2053	2057	rVB	661618	917797	6.63%	0.889%
67	14.385	2057	2061	2064	rBV2	346384	426931	3.08%	0.414%
68	14.428	2064	2068	2070	rBV2	302307	453554	3.28%	0.439%
69	14.452	2070	2072	2076	rVB	239139	266902	1.93%	0.259%
70	14.568	2086	2091	2096	rBV	639957	1003239	7.25%	0.972%
71	14.617	2096	2099	2103	rVB	411286	525628	3.80%	0.509%
72	14.678	2103	2109	2115	rBV2	1127734	2192579	15.83%	2.124%
73	14.751	2115	2121	2128	rVB3	1833372	3435847	24.81%	3.328%
74	14.836	2129	2135	2144	rVB	2812514	4447089	32.12%	4.307%
75	14.946	2146	2153	2157	rBV4	766793	1503918	10.86%	1.457%
76	15.056	2166	2171	2177	rBV3	393247	838595	6.06%	0.812%

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77	15.239	2193	2201	2208	rVB	8399518	13847143	100.00%	13.412%
78	15.342	2213	2218	2223	rBV3	419154	844621	6.10%	0.818%
79	15.440	2231	2234	2242	rVB5	176144	439125	3.17%	0.425%
80	15.531	2243	2249	2251	rBV3	260983	535023	3.86%	0.518%
81	15.720	2274	2280	2284	rVV3	538296	1039619	7.51%	1.007%
82	15.763	2284	2287	2292	rVB	325068	526152	3.80%	0.510%
83	15.836	2292	2299	2304	rBV2	607421	1330815	9.61%	1.289%
84	16.031	2326	2331	2336	rBV5	146915	298300	2.15%	0.289%
85	16.104	2337	2343	2349	rVV2	327470	764239	5.52%	0.740%
86	16.159	2349	2352	2357	rVV2	153859	238225	1.72%	0.231%
87	16.232	2358	2364	2368	rVV2	264856	546912	3.95%	0.530%
88	16.300	2368	2375	2389	rVB	4850161	9693139	70.00%	9.389%
89	16.415	2389	2394	2399	rBV2	121667	263648	1.90%	0.255%
90	16.495	2399	2407	2418	rVB	6149596	12433484	89.79%	12.043%
91	16.598	2420	2424	2434	rVB4	78736	192028	1.39%	0.186%

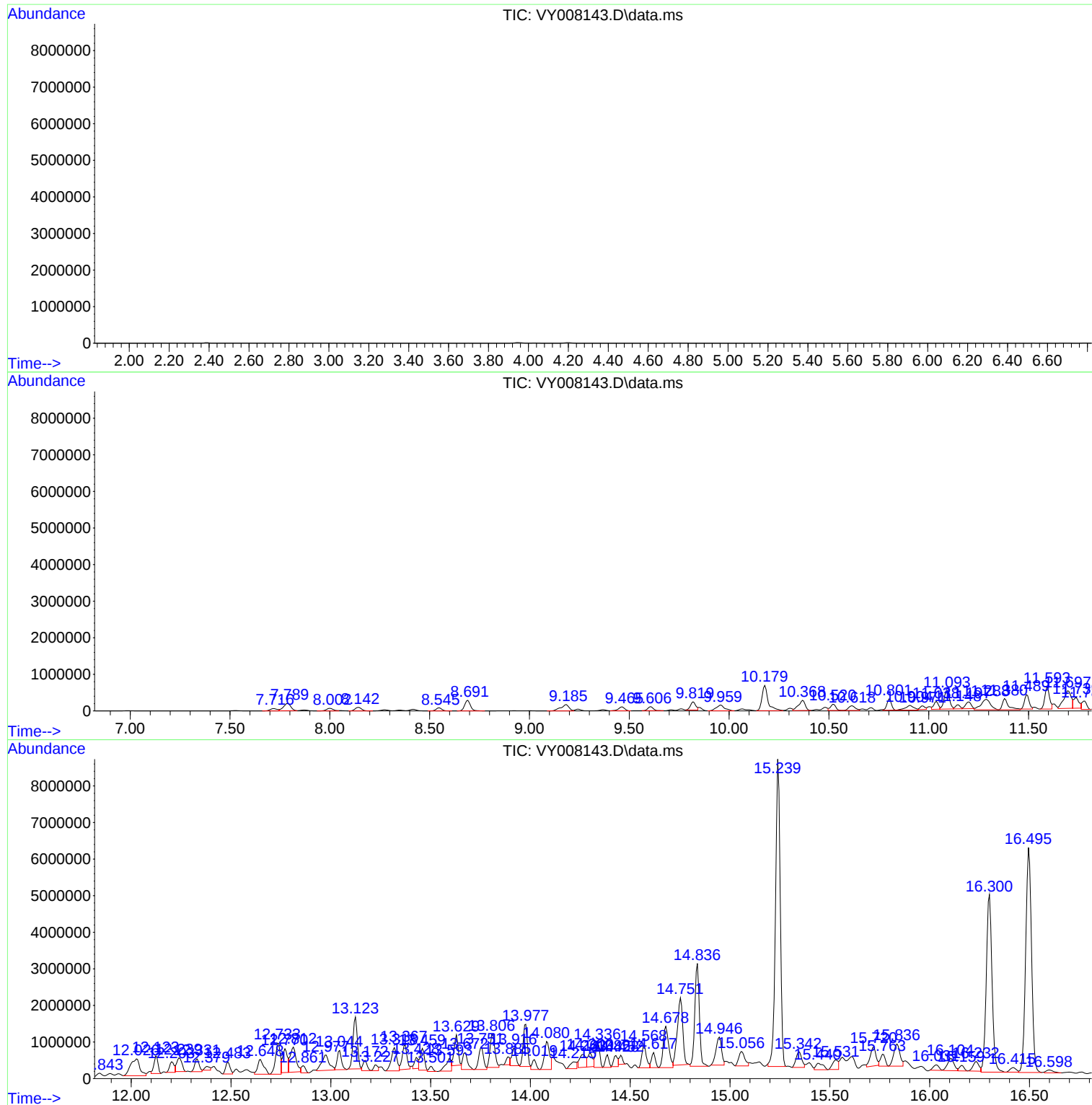
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032522S.M
Quant Title : SW846 8260

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



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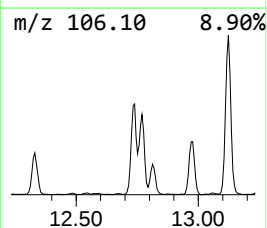
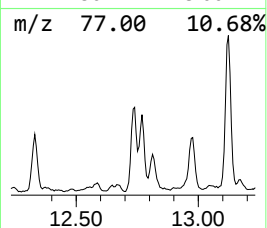
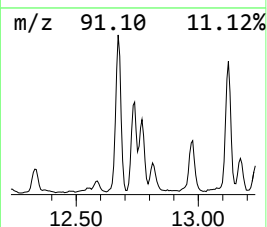
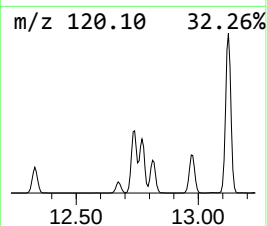
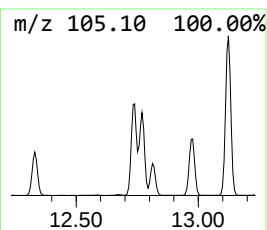
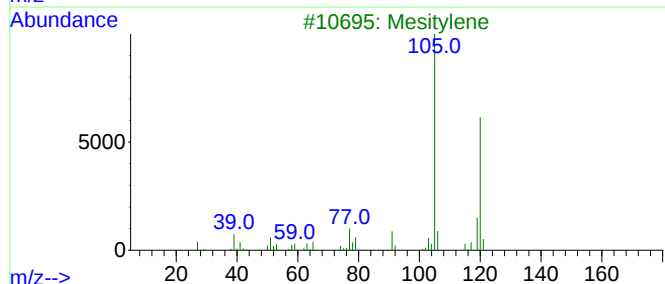
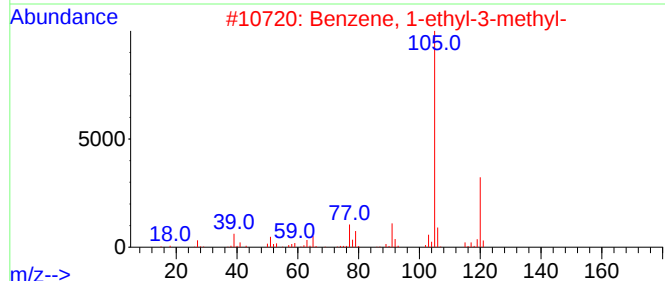
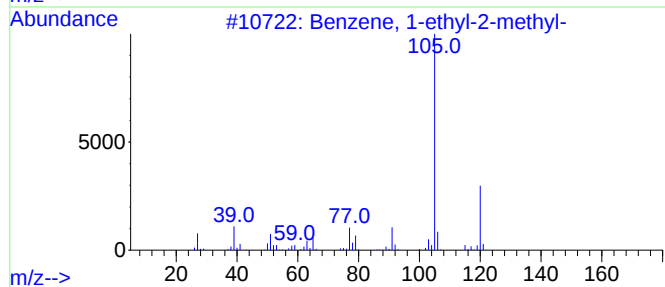
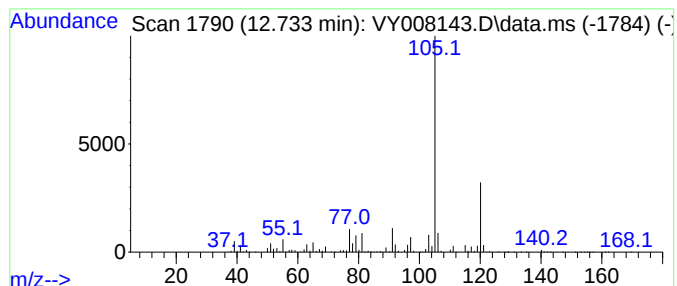
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032522S.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.733	174.47 ug/l	1466530	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	94
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3			Mesitylene	120	C9H12	000108-67-8	87
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	87
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	87



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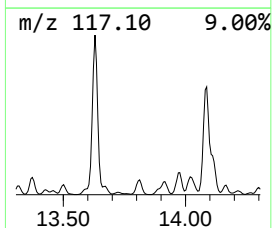
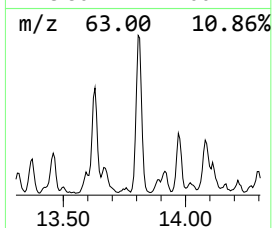
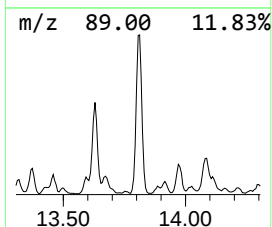
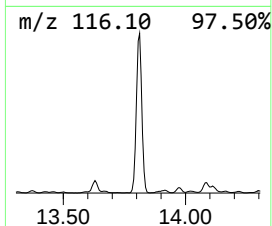
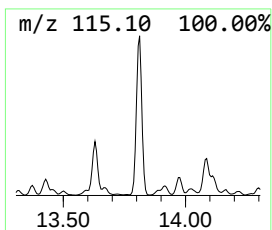
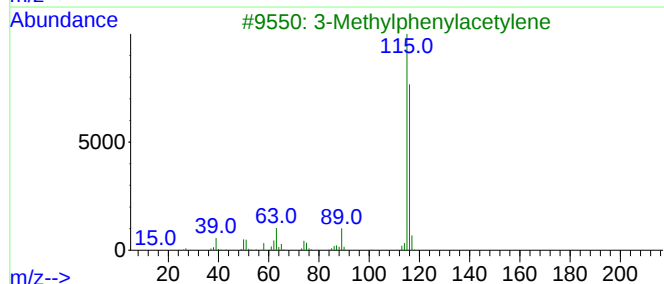
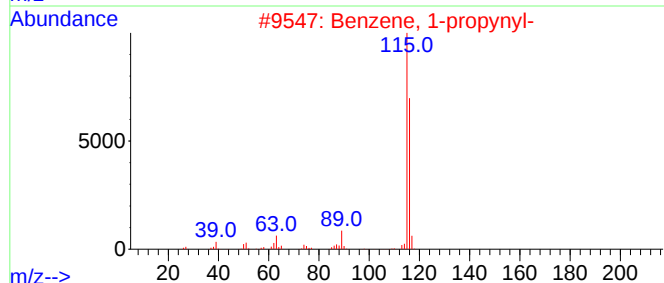
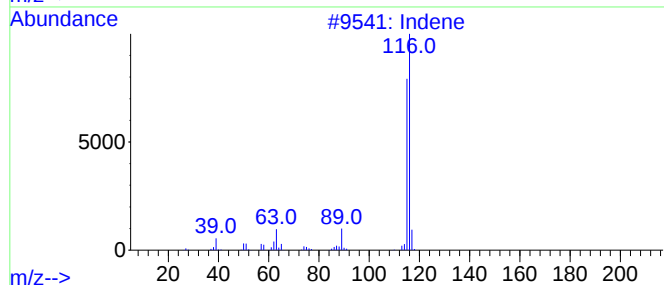
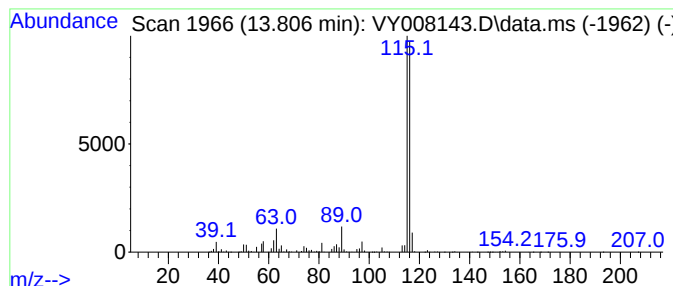
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032522S.M
Quant Title : SW846 8260

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

Peak Number 2 Indene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.806	181.08 ug/l	1522050	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indene	116	C9H8	000095-13-6	97
2			Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3			3-Methylphenylacetylene	116	C9H8	000766-82-5	91
4			Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	90
5			2-Methylphenylacetylene	116	C9H8	000766-47-2	83



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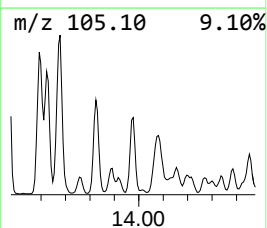
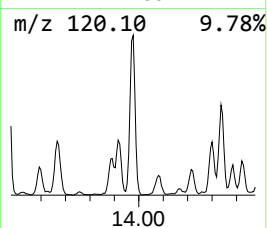
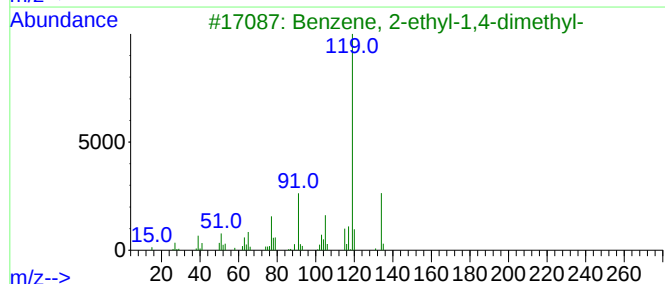
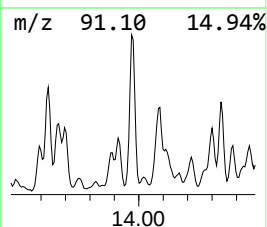
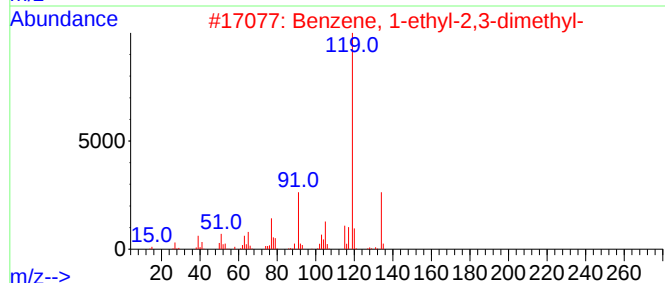
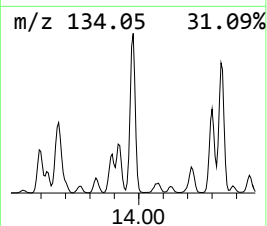
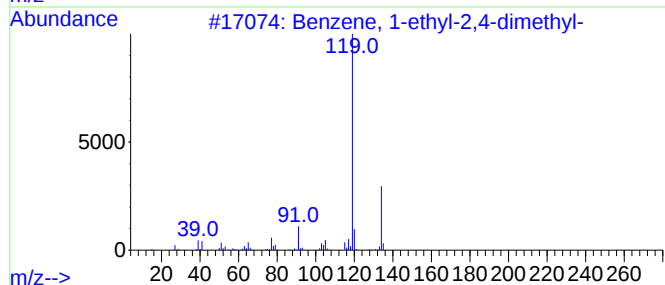
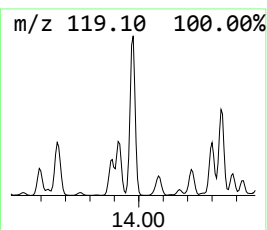
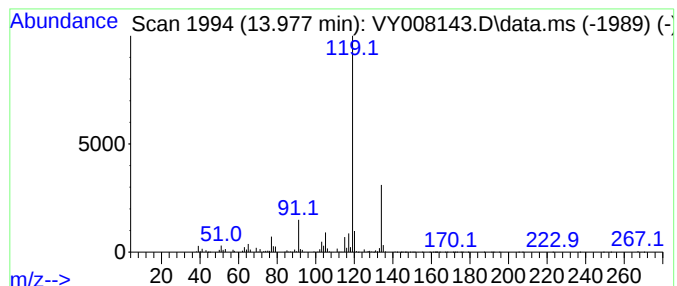
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y032522S.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.977	197.96 ug/l	1663920	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033122\
Data File : VY008143.D
Acq On : 31 Mar 2022 17:06
Operator : SY/MD
Sample : N2123-02
Misc : 4.49g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
20220339-AMENDED-COMP-CS-1

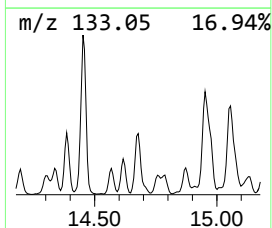
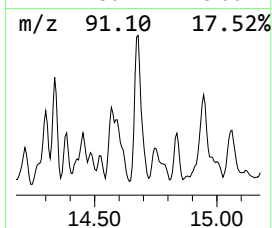
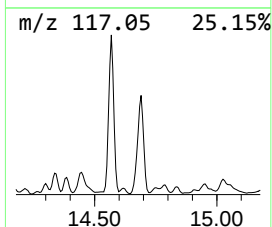
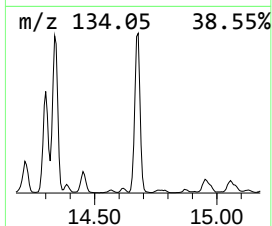
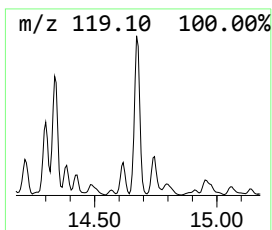
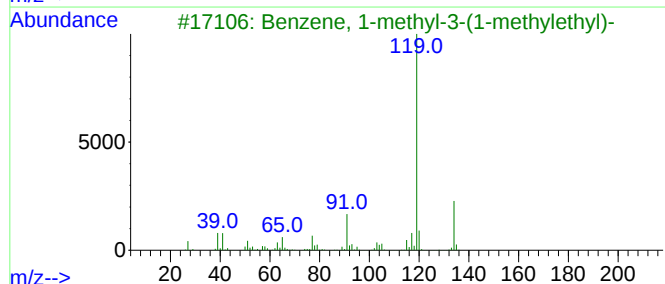
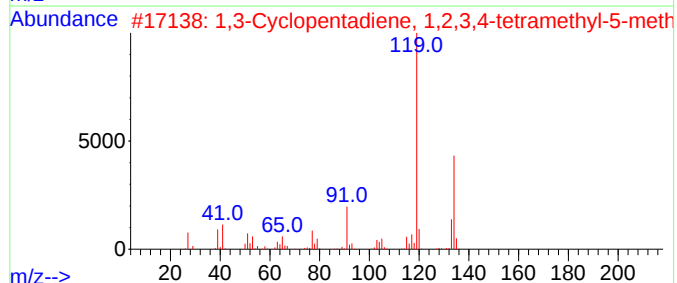
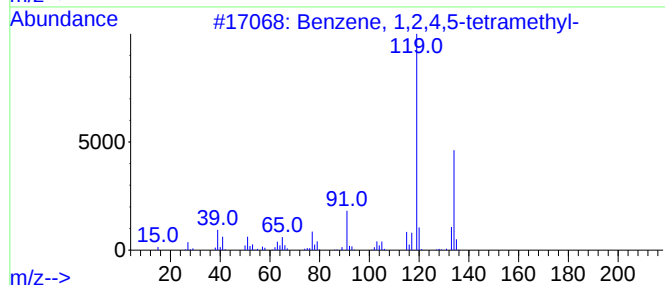
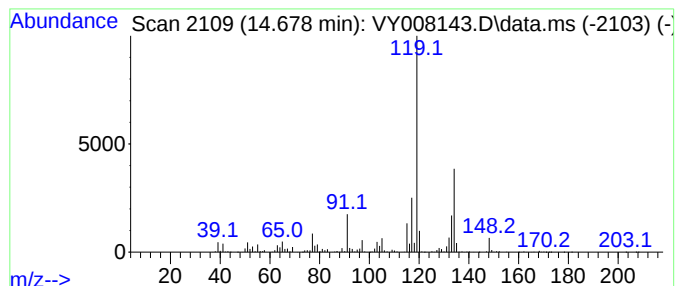
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.678	260.85 ug/l	2192580	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
2			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	93
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	81
4			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	81
5			o-Cymene	134	C10H14	000527-84-4	81



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033122\
Data File : VY008143.D
Acq On : 31 Mar 2022 17:06
Operator : SY/MD
Sample : N2123-02
Misc : 4.49g/5.0mL/MSVOA_Y/Soil
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
20220339-AMENDED-COMP-CS-1

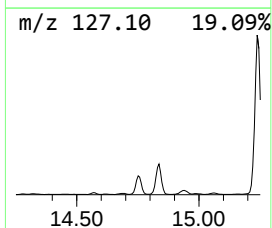
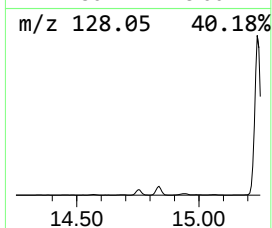
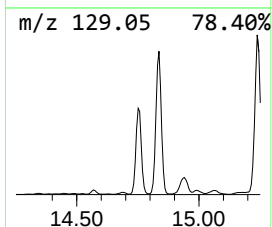
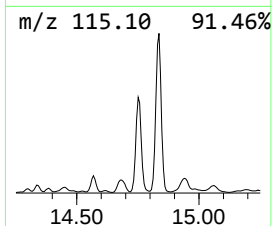
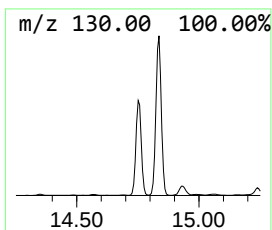
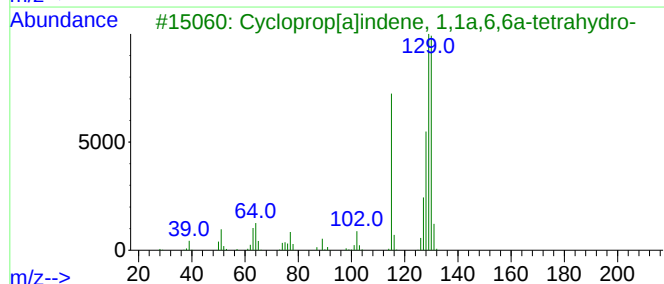
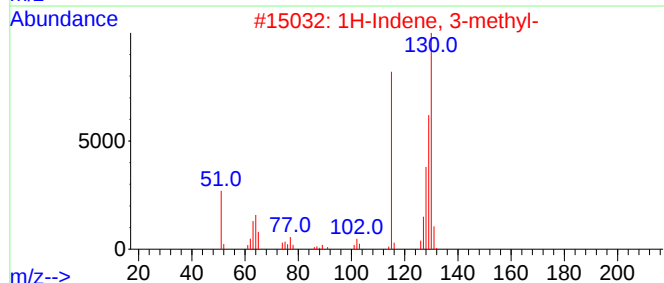
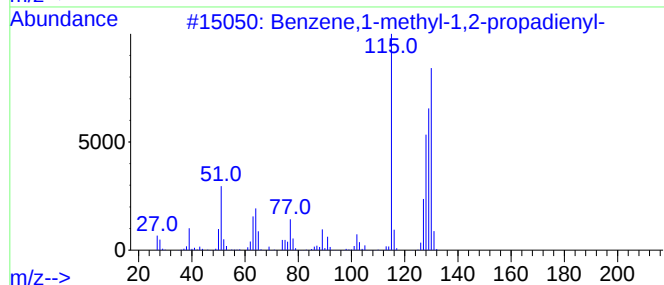
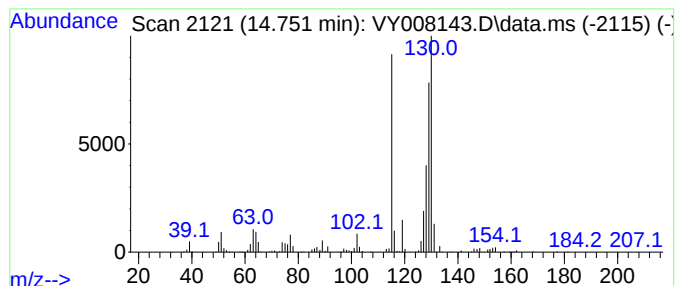
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Quant Title : SW846 8260

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

Peak Number 5 Benzene,1-methyl-1,2-propad... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.751	408.76 ug/l	3435850	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	93
2			1H-Indene, 3-methyl-	130	C10H10	000767-60-2	93
3			Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	92
4			Benzene, 1-butyryl-	130	C10H10	000622-76-4	91
5			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033122\
Data File : VY008143.D
Acq On : 31 Mar 2022 17:06
Operator : SY/MD
Sample : N2123-02
Misc : 4.49g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
20220339-AMENDED-COMP-CS-1

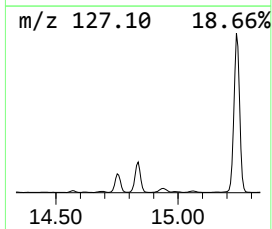
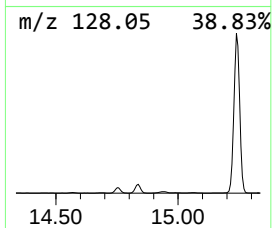
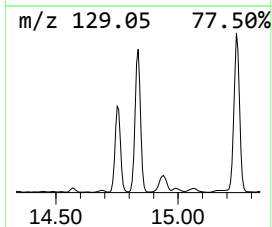
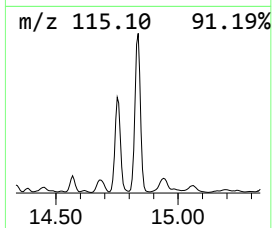
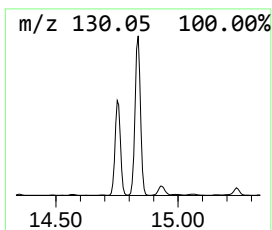
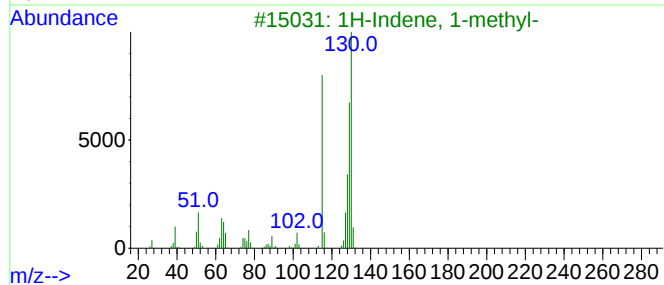
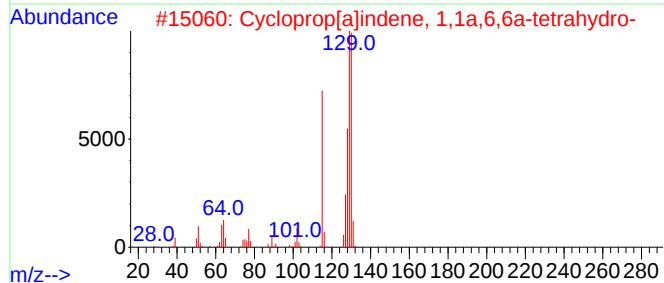
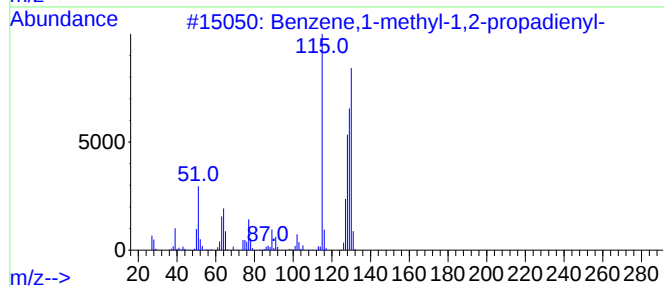
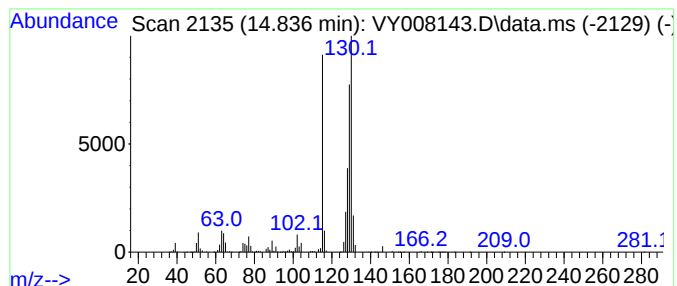
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Quant Title : SW846 8260

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

Peak Number 6 Cycloprop[a]indene, 1,1a,6,... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.836	529.07 ug/l	4447090	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	94
2			Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	93
3			1H-Indene, 1-methyl-	130	C10H10	000767-59-9	93
4			Benzene, 1-butyryl-	130	C10H10	000622-76-4	91
5			Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033122\
Data File : VY008143.D
Acq On : 31 Mar 2022 17:06
Operator : SY/MD
Sample : N2123-02
Misc : 4.49g/5.0mL/MSVOA_Y/SoIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
20220339-AMENDED-COMP-CS-1

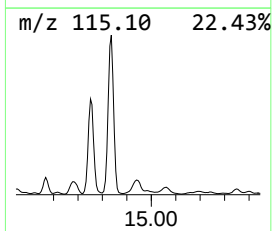
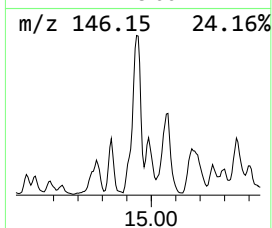
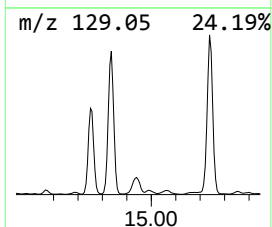
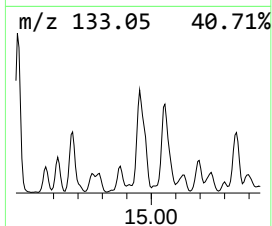
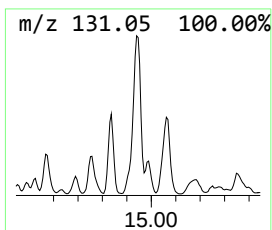
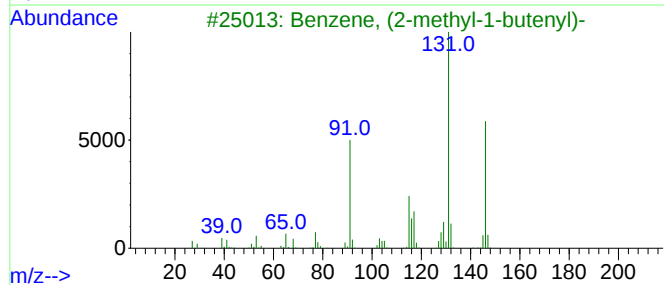
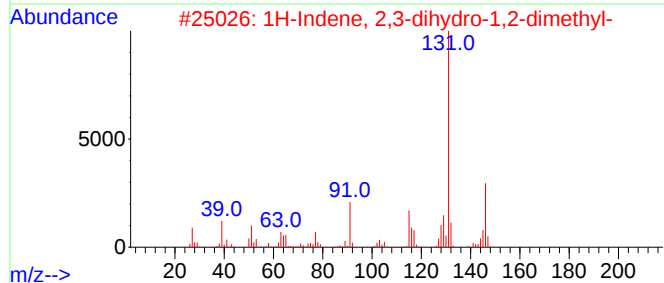
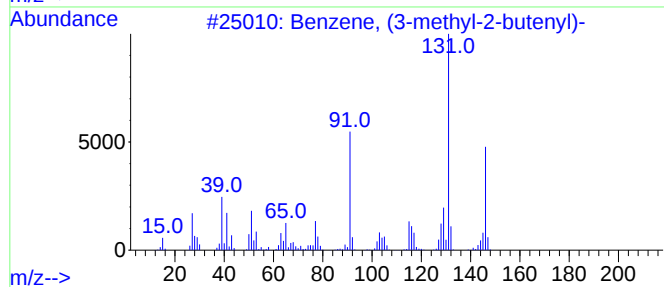
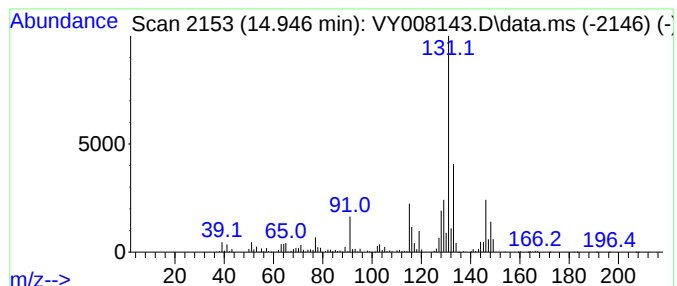
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, (3-methyl-2-butenyl)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.946	178.92 ug/l	1503920	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	64
3			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	58
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	52
5			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033122\
Data File : VY008143.D
Acq On : 31 Mar 2022 17:06
Operator : SY/MD
Sample : N2123-02
Misc : 4.49g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
20220339-AMENDED-COMP-CS-1

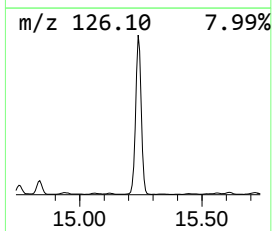
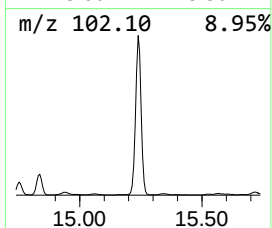
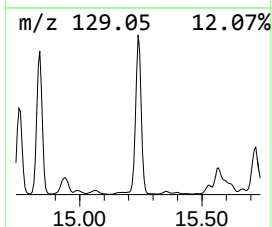
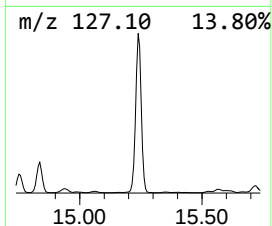
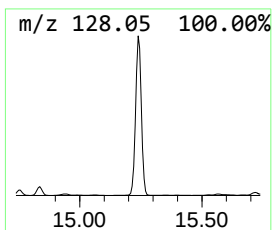
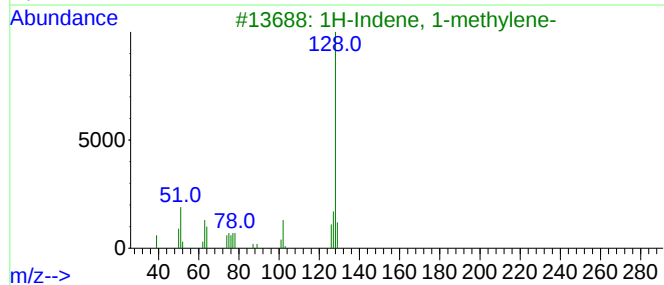
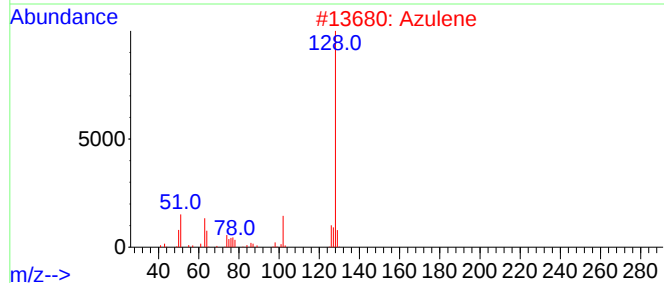
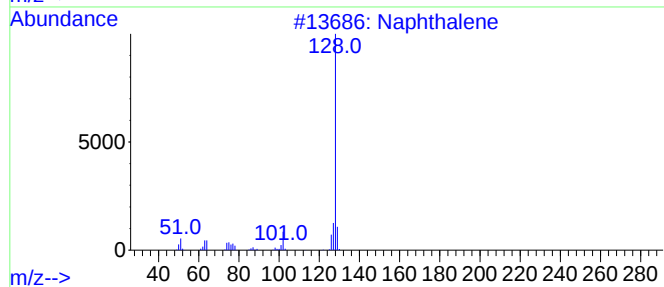
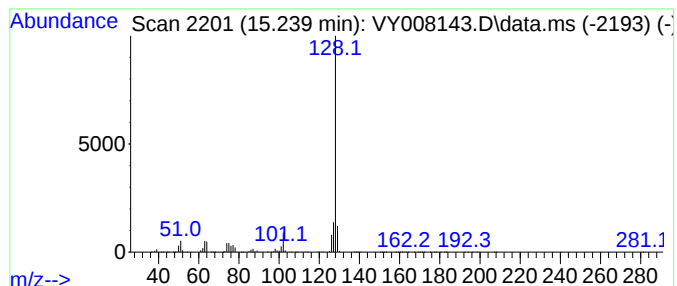
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Quant Title : SW846 8260

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

Peak Number 8 Azulene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.239	1647.38 ug/l	13847100	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Azulene	128	C10H8	000275-51-4	91
3			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	91
4			4a,8a-(Methaniminomethano)naphth...	275	C18H13NO2	069915-10-2	78
5			1,3-Dicyanobenzene	128	C8H4N2	000626-17-5	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033122\
Data File : VY008143.D
Acq On : 31 Mar 2022 17:06
Operator : SY/MD
Sample : N2123-02
Misc : 4.49g/5.0mL/MSVOA_Y/soil
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
20220339-AMENDED-COMP-CS-1

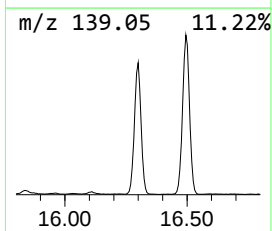
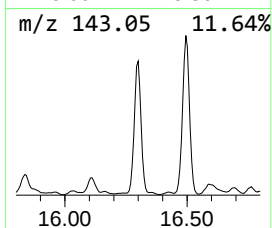
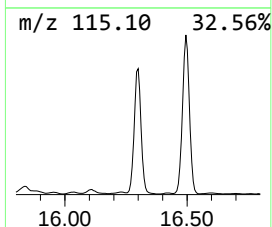
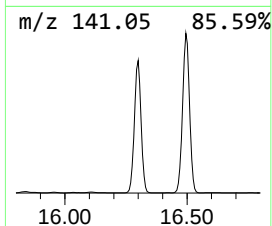
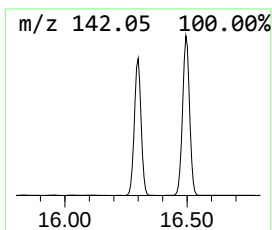
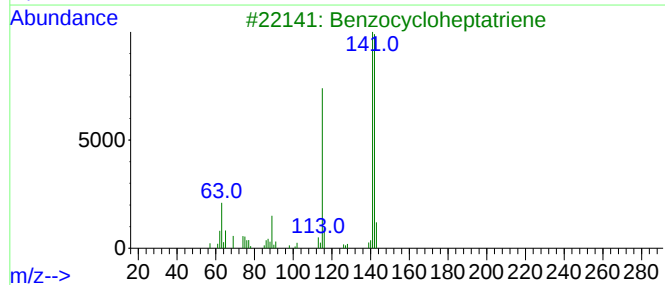
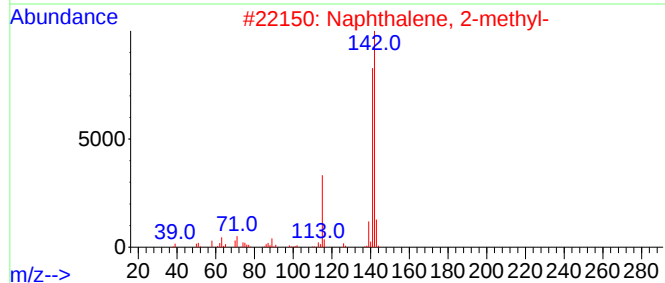
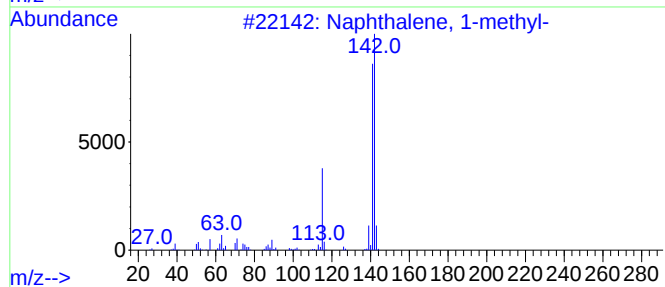
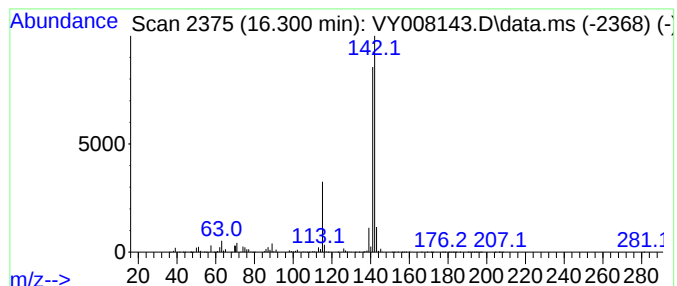
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Quant Title : SW846 8260

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

Peak Number 9 Benzocycloheptatriene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.299	1153.18 ug/l	9693140	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	91
5			1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033122\
Data File : VY008143.D
Acq On : 31 Mar 2022 17:06
Operator : SY/MD
Sample : N2123-02
Misc : 4.49g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
20220339-AMENDED-COMP-CS-1

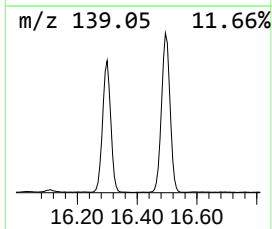
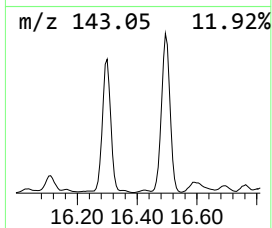
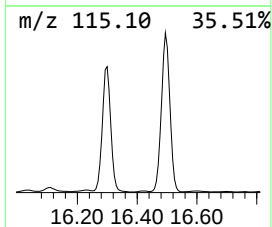
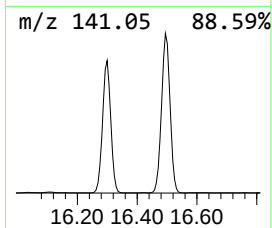
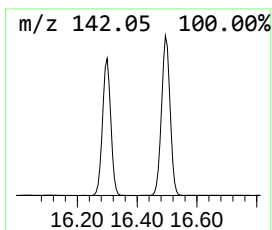
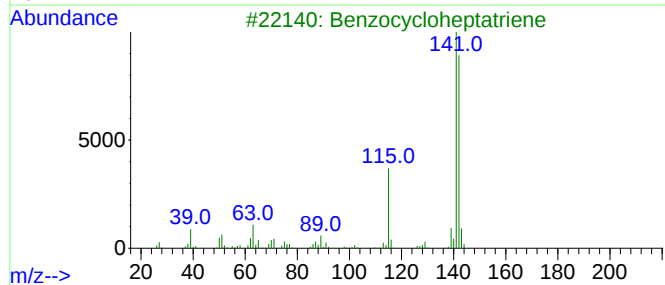
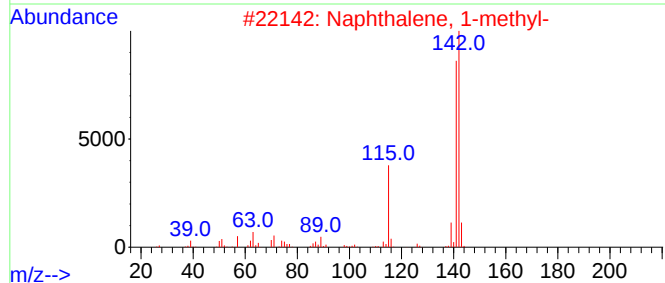
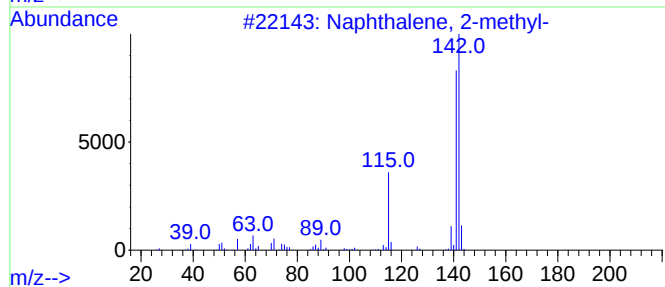
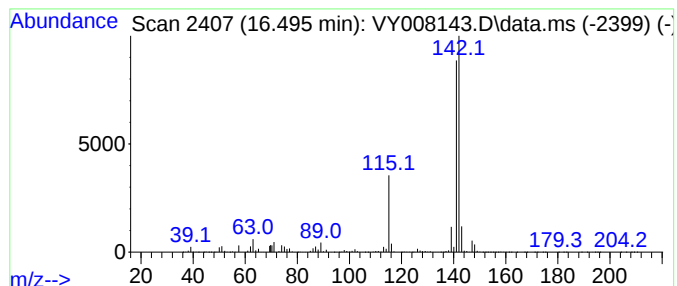
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Quant Title : SW846 8260

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

Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.495	1479.20 ug/l	12433500	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	96
3			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
5			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY033122\
Data File : VY008143.D
Acq On : 31 Mar 2022 17:06
Operator : SY/MD
Sample : N2123-02
Misc : 4.49g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
20220339-AMENDED-COMP-CS-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
Benzene, 1-ethy...	12.733	174.5	ug/l	1466530	4	13.428	420277	50.0
Indene	13.806	181.1	ug/l	1522050	4	13.428	420277	50.0
Benzene, 1-ethy...	13.977	198.0	ug/l	1663920	4	13.428	420277	50.0
Benzene, 1,2,4,...	14.678	260.9	ug/l	2192580	4	13.428	420277	50.0
Benzene,1-methy...	14.751	408.8	ug/l	3435850	4	13.428	420277	50.0
Cycloprop[a]ind...	14.836	529.1	ug/l	4447090	4	13.428	420277	50.0
Benzene, (3-met...	14.946	178.9	ug/l	1503920	4	13.428	420277	50.0
Azulene	15.239	1647.4	ug/l	13847100	4	13.428	420277	50.0
Benzocyclohepta...	16.299	1153.2	ug/l	9693140	4	13.428	420277	50.0
1,4-Methanonaph...	16.495	1479.2	ug/l	12433500	4	13.428	420277	50.0