LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011723\
Data File : VY012219.D
Acq On : 17 Jan 2023 13:03
Operator : KP/MD
Sample : 01159-04
Misc : 4.60g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
MH-6

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Title : SW846 8260

Signal : TIC: VY012219.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.722	463	476	485	rBV2	20619	63723	4.81%	0.307%
2	5.753	636	645	657	rVB3	6302	19426	1.47%	0.094%
3	6.972	833	845	860	rBV	36536	111942	8.44%	0.539%
4	7.716	957	967	973	rBV	182866	437784	33.02%	2.108%
5	7.789	973	979	997	rVB	321396	736112	55.53%	3.544%
6	8.142	1023	1037	1053	rBV2	191316	488951	36.88%	2.354%
7	8.691	1118	1127	1141	rVB	492458	949703	71.64%	4.573%
8	9.185	1195	1208	1215	rBV4	12640	32406	2.44%	0.156%
9	10.179	1364	1371	1387	rVB	772105	1325718	100.00%	6.384%
10	10.374	1394	1403	1407	rVB6	7064	17828	1.34%	0.086%
11	10.581	1431	1437	1449	rVB	40830	81719	6.16%	0.393%
12	10.947	1491	1497	1504	rBV	40365	71751	5.41%	0.345%
13	11.038	1507	1512	1516	rVV	19622	30757	2.32%	0.148%
14	11.093	1516	1521	1526	rVB	17582	31046	2.34%	0.149%
15	11.203	1534	1539	1543	rBV3	10662	18271	1.38%	0.088%
16	11.282	1543	1552	1555	rVV4	20381	48252	3.64%	0.232%
17	11.319	1555	1558	1563	rVB	24254	38245	2.88%	0.184%
18	11.380	1563	1568	1573	rBV	16095	26610	2.01%	0.128%
19	11.489	1579	1586	1596	rBV	741542	1199272	90.46%	5.775%
20	11.593	1597	1603	1606	rBV3	14559	26376	1.99%	0.127%
21	11.703	1612	1621	1625	rBV2	38406	108195	8.16%	0.521%
22	11.782	1630	1634	1639	rVB2	21790	37836	2.85%	0.182%
23	12.007	1663	1671	1678	rBV4	37542	122101	9.21%	0.588%
24	12.123	1686	1690	1694	rVB	38271	51821	3.91%	0.250%
25	12.203	1699	1703	1706	rBV4	32356	54793	4.13%	0.264%
26	12.245	1706	1710	1715	rVB3	53392	96334	7.27%	0.464%
27	12.331	1719	1724	1727	rBV2	26589	42635	3.22%	0.205%
28	12.385	1728	1733	1735	rBV5	17659	31775	2.40%	0.153%
29	12.416	1735	1738	1743	rVB2	34814	48030	3.62%	0.231%
30	12.483	1743	1749	1755	rBV	633099	993267	74.92%	4.783%
31	12.672	1772	1780	1784	rVB2	86904	197916	14.93%	0.953%
32	12.733	1784	1790	1793	rBV4	62684	110773	8.36%	0.533%
33	12.812	1793	1803	1808	rVV5	145648	383536	28.93%	1.847%
34	12.861	1808	1811	1816	rVB2	53978	79482	6.00%	0.383%
35	13.044	1837	1841	1848	rVB	123248	186031	14.03%	0.896%
36	13.123	1848	1854	1859	rBV	132541	202335	15.26%	0.974%

LSC Area Percent Report

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Operator : KP/MD
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Misc : 4.60g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
MH-6

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Title : SW846 8260

37	13.172	1859	1862	1866	rVV4	23305	30366	2.29%	0.146%
38	13.251	1868	1875	1879	rVV4	62595	118791	8.96%	0.572%
39	13.318	1879	1886	1890	rVV4	116364	241170	18.19%	1.161%
40	13.361	1890	1893	1897	rVV3	72114	109997	8.30%	0.530%
41	13.428	1897	1904	1913	rVV	776167	1281965	96.70%	6.173%
42	13.501	1913	1916	1919	rVB3	18873	22390	1.69%	0.108%
43	13.568	1924	1927	1928	rBV3	16888	20019	1.51%	0.096%
44	13.593	1928	1931	1933	rBV	42207	51949	3.92%	0.250%
45	13.629	1933	1937	1941	rVB	178723	261497	19.72%	1.259%
46	13.678	1941	1945	1946	rBV2	67570	93706	7.07%	0.451%
47	13.751	1952	1957	1961	rBV3	195872	323934	24.43%	1.560%
48	13.794	1961	1964	1968	rVB2	33833	47713	3.60%	0.230%
49	13.916	1981	1984	1988	rVB2	48637	58154	4.39%	0.280%
50	13.971	1989	1993	1998	rVB	240154	335461	25.30%	1.615%
51	14.025	1998	2002	2006	rVB2	72486	126082	9.51%	0.607%
52	14.080	2006	2011	2016	rVB2	201523	331745	25.02%	1.597%
53	14.153	2016	2023	2027	rBV7	61000	138631	10.46%	0.668%
54	14.196	2027	2030	2032	rBV2	32295	42661	3.22%	0.205%
55	14.263	2036	2041	2044	rBV3	136356	234795	17.71%	1.131%
56	14.379	2056	2060	2063	rBV	81238	112563	8.49%	0.542%
57	14.422	2063	2067	2071	rVV3	112992	154273	11.64%	0.743%
58	14.483	2074	2077	2080	rVV3	35323	46498	3.51%	0.224%
59	14.568	2086	2091	2096	rBV2	188928	351855	26.54%	1.694%
60	14.617	2096	2099	2103	rVV3	98983	130615	9.85%	0.629%
61	14.684	2103	2110	2114	rVV4	418070	860668	64.92%	4.144%
62	14.733	2114	2118	2125	rVB3	269295	492335	37.14%	2.371%
63	14.836	2131	2135	2140	rVB	165948	240177	18.12%	1.156%
64	14.909	2143	2147	2148	rBV2	61176	71709	5.41%	0.345%
65	14.946	2149	2153	2156	rVV2	191030	280032	21.12%	1.348%
66	14.983	2156	2159	2165	rVB2	73743	96408	7.27%	0.464%
67	15.056	2165	2171	2177	rVB2	166629	330789	24.95%	1.593%
68	15.129	2179	2183	2184	rBV2	30772	47933	3.62%	0.231%
69	15.220	2192	2198	2199	rBV3	179182	294208	22.19%	1.417%
70	15.293	2206	2210	2214	rBV	127606	182588	13.77%	0.879%
71	15.348	2214	2219	2225	rBV4	106344	261394	19.72%	1.259%
72	15.434	2231	2233	2242	rVB8	79225	167480	12.63%	0.806%
73	15.531	2242	2249	2255	rBV	141583	324160	24.45%	1.561%
74	15.611	2257	2262	2267	rVB5	82316	148326	11.19%	0.714%
75	15.690	2267	2275	2278	rBV3	107884	267415	20.17%	1.288%
76	15.732	2278	2282	2291	rVB	245145	439452	33.15%	2.116%

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Misc : 4.60g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
MH-6

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Title : SW846 8260

77	15.818	2293	2296	2301	rVB3	47878	78172	5.90%	0.376%
78	15.891	2302	2308	2313	rBV3	80084	167673	12.65%	0.807%
79	16.037	2325	2332	2338	rBV2	143593	309344	23.33%	1.490%
80	16.159	2348	2352	2356	rBV2	55042	88488	6.67%	0.426%
81	16.232	2358	2364	2368	rVV2	145244	313902	23.68%	1.511%
82	16.293	2368	2374	2387	rVB	630183	1291106	97.39%	6.217%
83	16.488	2399	2406	2414	rBV	321603	730981	55.14%	3.520%
84	16.629	2421	2429	2438	rVB7	36484	113438	8.56%	0.546%

Sum of corrected areas: 20767760

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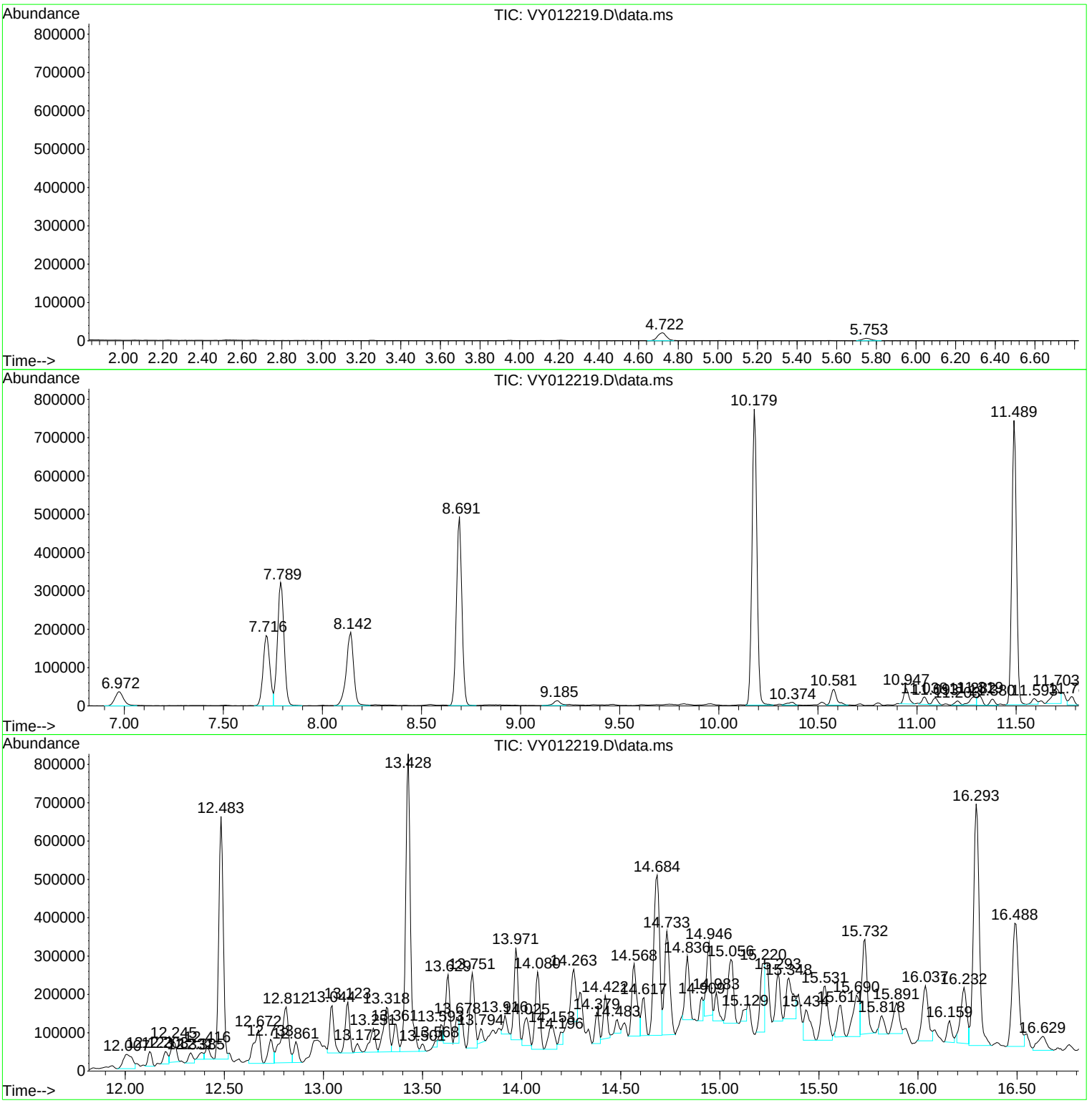
Instrument :
MSVOA_Y
ClientSampleId :
MH-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011723\
Data File : VY012219.D
Acq On : 17 Jan 2023 13:03
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Instrument :
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ClientSampleId :
MH-6

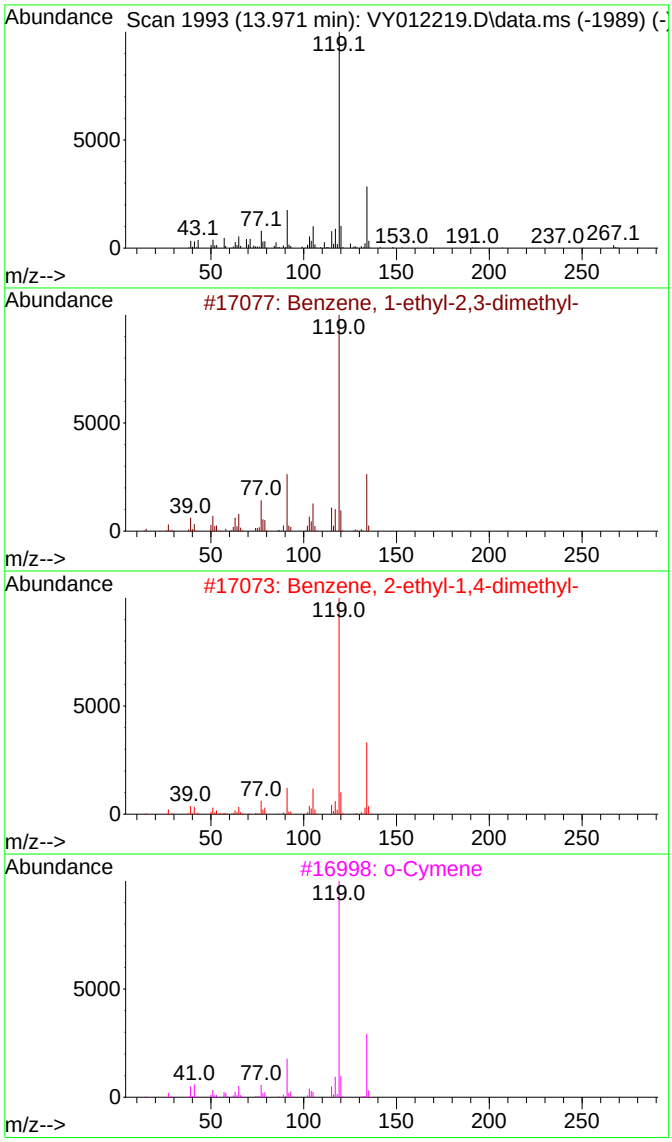
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Quant Title : SW846 8260

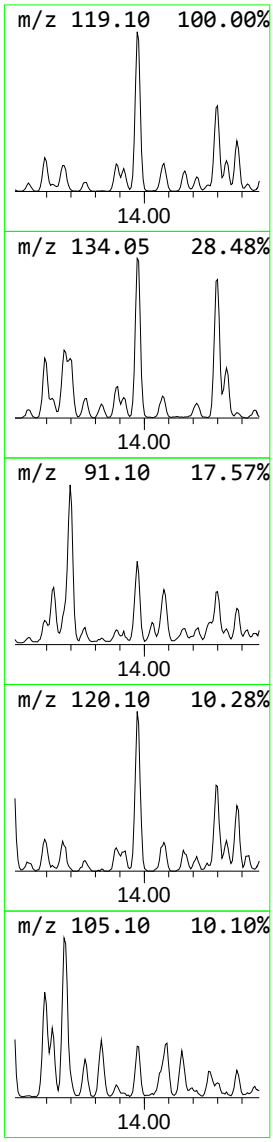
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.971	13.08 ug/l	335461	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011723\
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Acq On : 17 Jan 2023 13:03
Operator : KP/MD
Sample : 01159-04
Misc : 4.60g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
MH-6

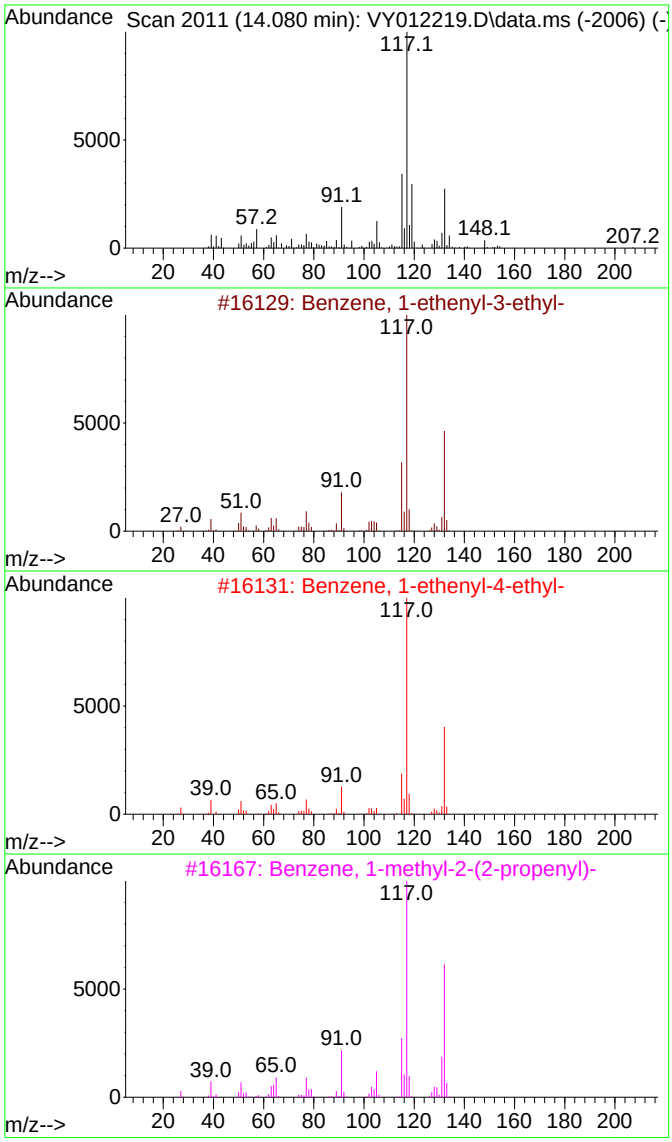
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Quant Title : SW846 8260

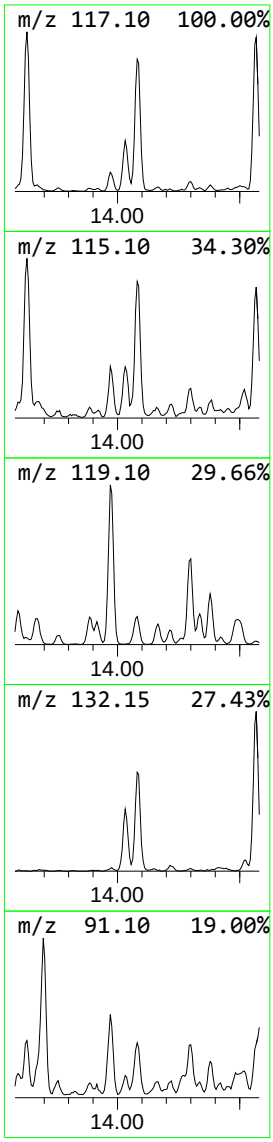
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.080	12.94 ug/l	331745	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	74
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	74
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	74
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	72





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Operator : KP/MD
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Misc : 4.60g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
MH-6

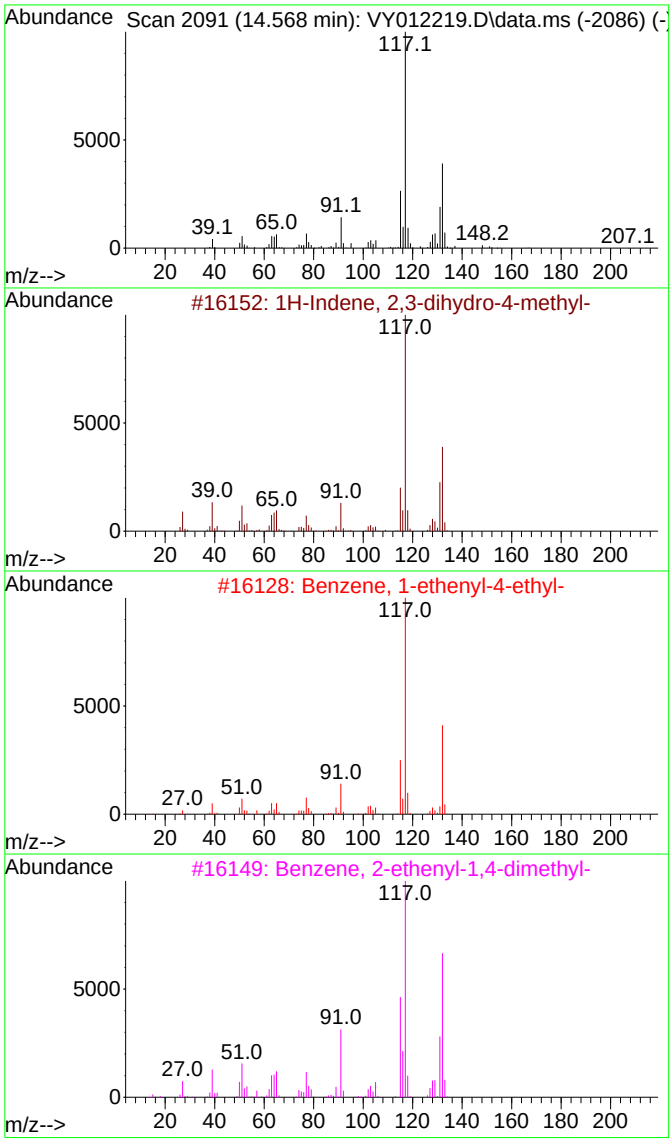
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Quant Title : SW846 8260

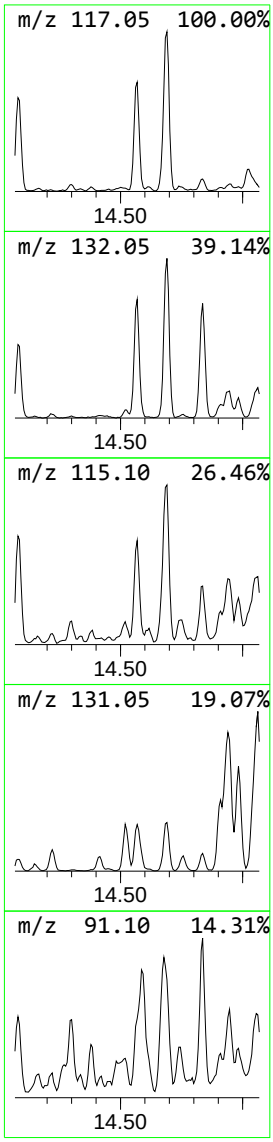
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.568	13.72 ug/l	351855	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91





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Operator : KP/MD
Sample : 01159-04
Misc : 4.60g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
MH-6

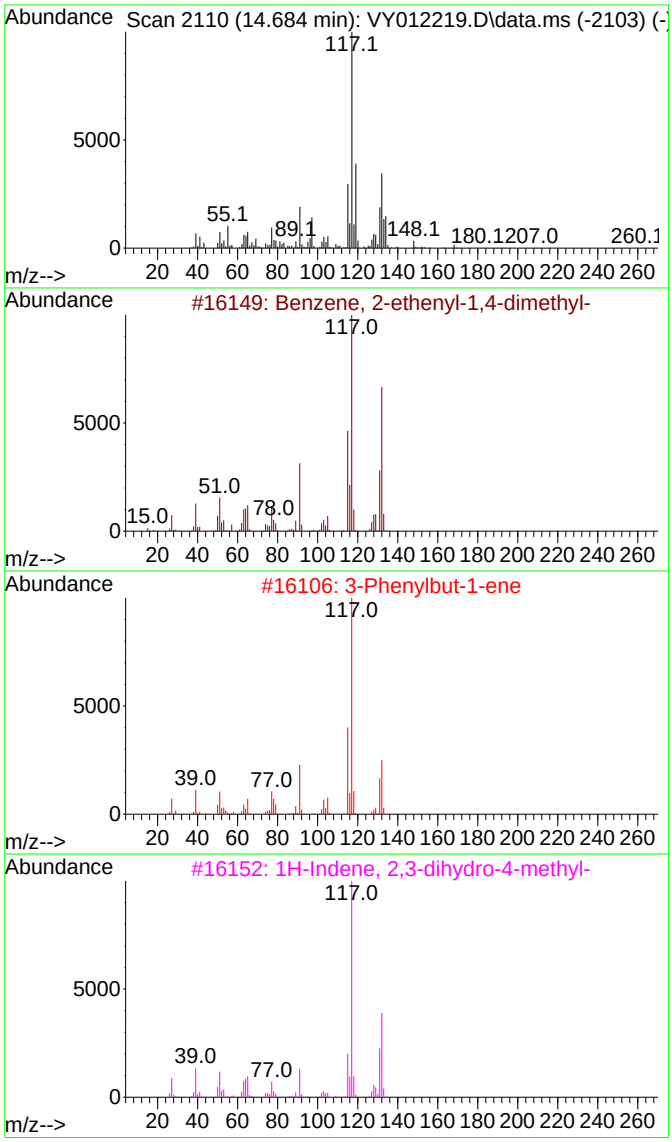
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Quant Title : SW846 8260

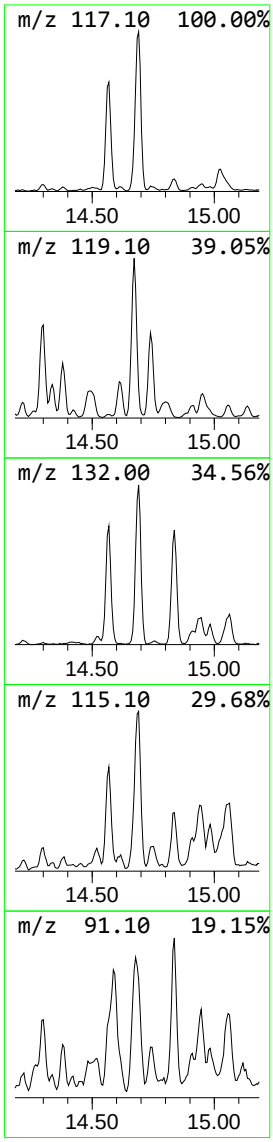
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	33.57 ug/l	860668	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	90
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64





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Misc : 4.60g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
MH-6

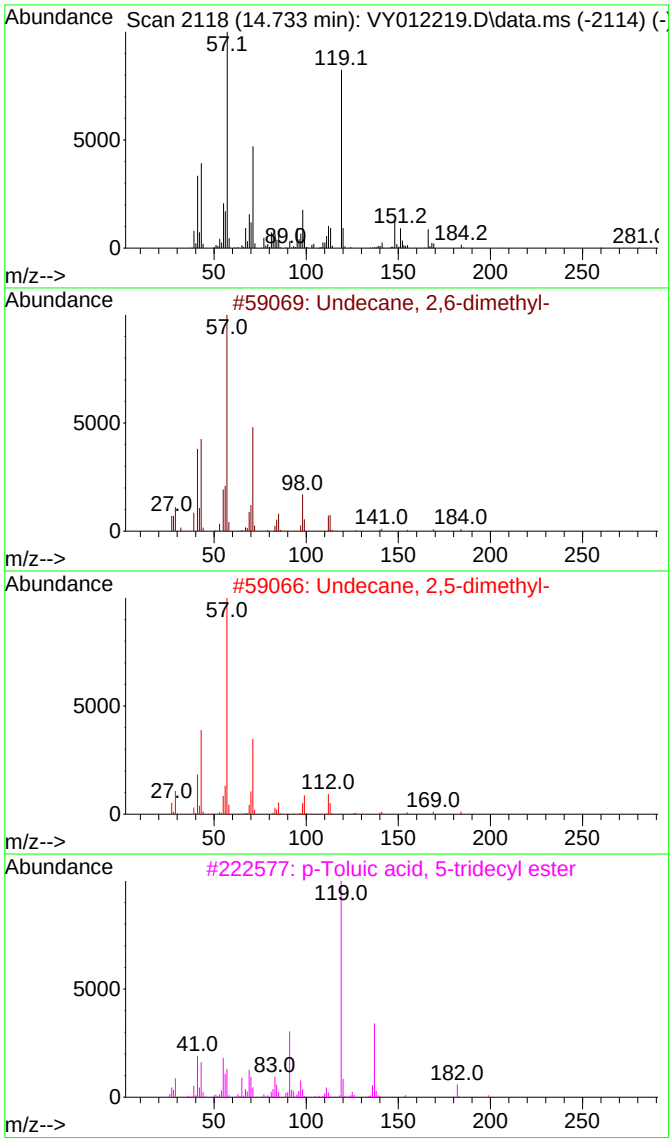
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Quant Title : SW846 8260

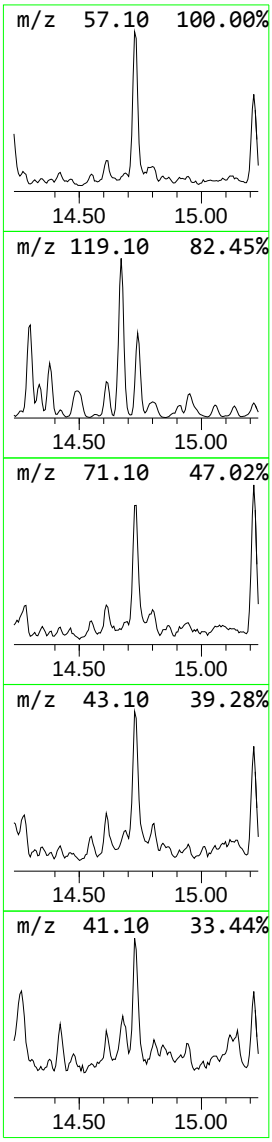
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Undecane, 2,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.733	19.20 ug/l	492335	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	81
2	Undecane, 2,5-dimethyl-	184	C13H28	017301-22-3	42
3	p-Toluic acid, 5-tridecyl ester	318	C21H34O2	1000299-80-5	35
4	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	35
5	Sulfurous acid, butyl octyl ester	250	C12H26O3S	1000309-17-5	25





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011723\
Data File : VY012219.D
Acq On : 17 Jan 2023 13:03
Operator : KP/MD
Sample : 01159-04
Misc : 4.60g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
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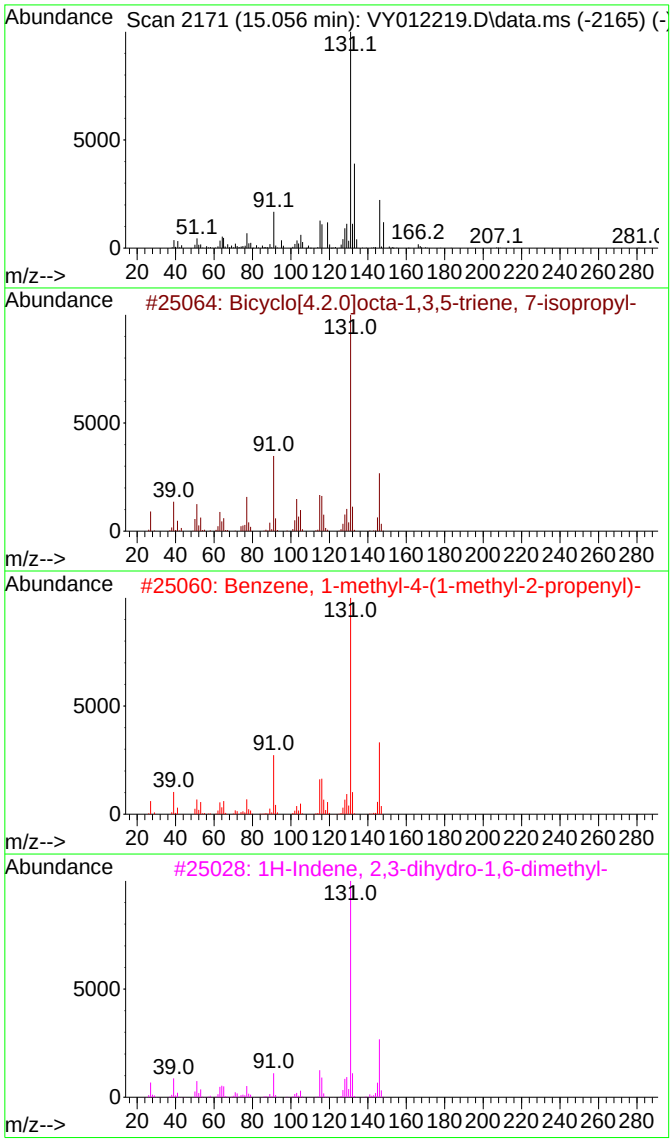
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Quant Title : SW846 8260

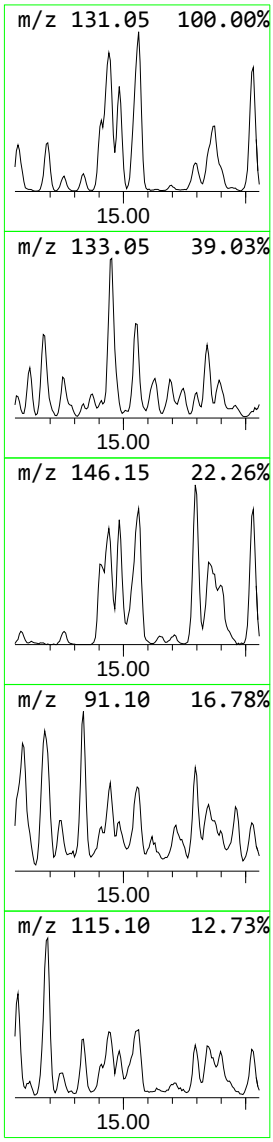
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.056	12.90 ug/l	330789	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	91
2	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	70
3	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	64
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	64
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011723\
Data File : VY012219.D
Acq On : 17 Jan 2023 13:03
Operator : KP/MD
Sample : 01159-04
Misc : 4.60g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
MH-6

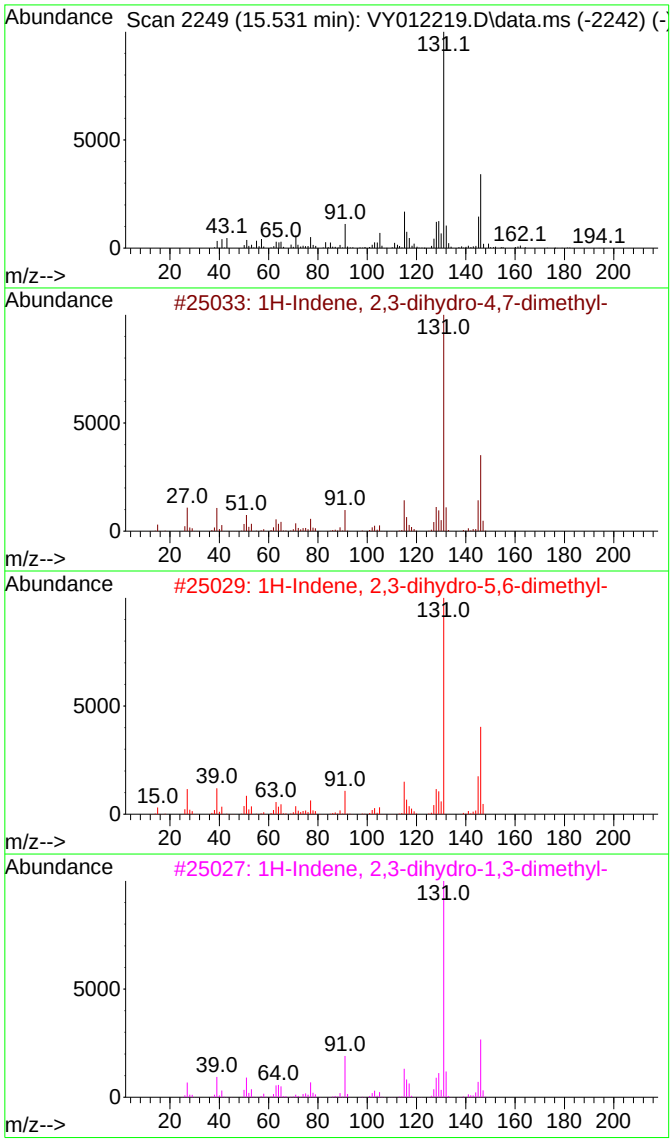
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Quant Title : SW846 8260

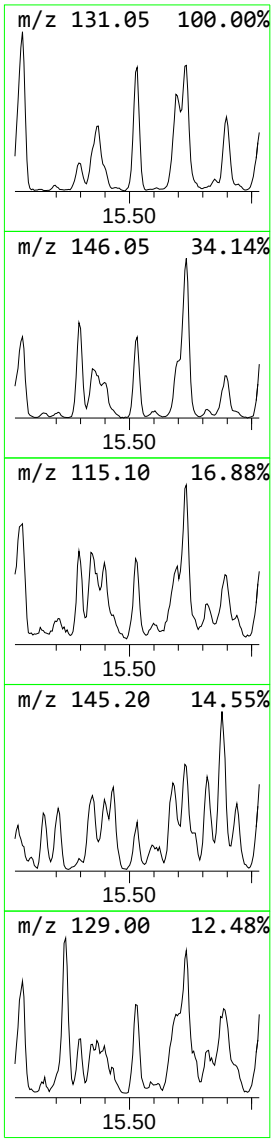
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.531	12.64 ug/l	324160	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
2	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	93
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011723\
Data File : VY012219.D
Acq On : 17 Jan 2023 13:03
Operator : KP/MD
Sample : 01159-04
Misc : 4.60g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
MH-6

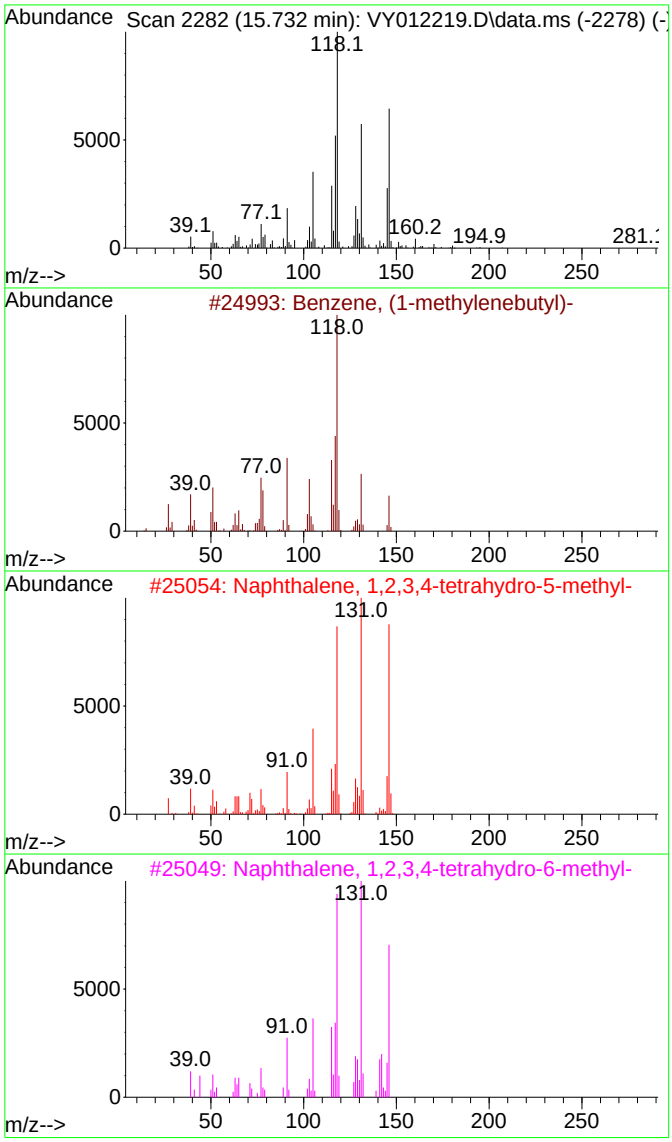
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Quant Title : SW846 8260

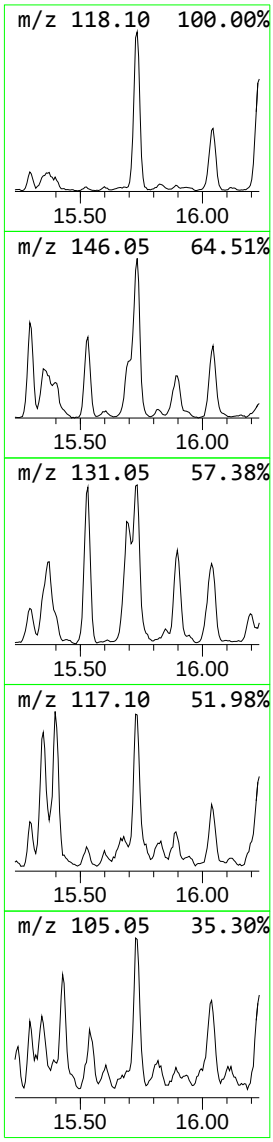
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, (1-methylenebutyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.732	17.14 ug/l	439452	1,4-Dichlorobenzene-d4	13.428

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	74
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	59
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	58
4	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	58
5	3-Phenylcyclopentanecarboxylic acid	190	C12H14O2	091495-75-9	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011723\
Data File : VY012219.D
Acq On : 17 Jan 2023 13:03
Operator : KP/MD
Sample : 01159-04
Misc : 4.60g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
MH-6

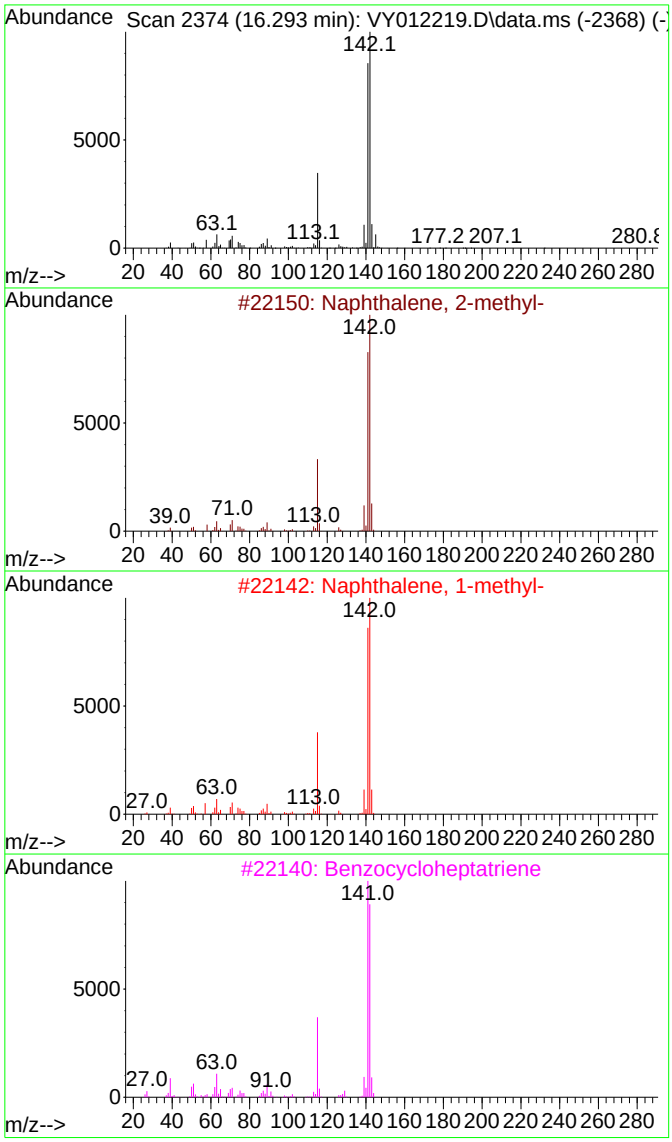
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Quant Title : SW846 8260

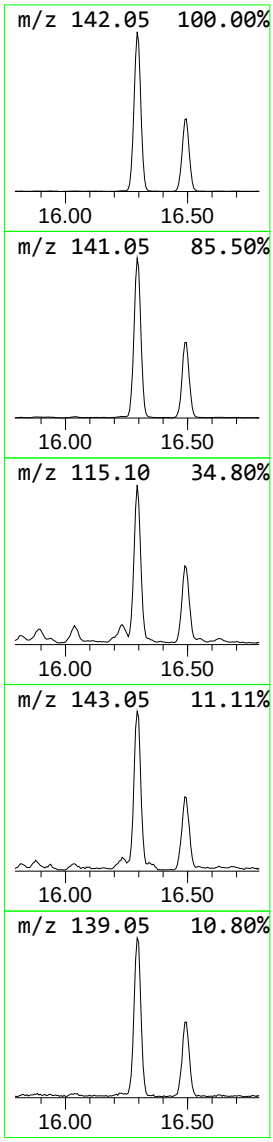
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzocycloheptatriene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.293	50.36 ug/l	1291110	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	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	59





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Data File : VY012219.D
Acq On : 17 Jan 2023 13:03
Operator : KP/MD
Sample : 01159-04
Misc : 4.60g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
MH-6

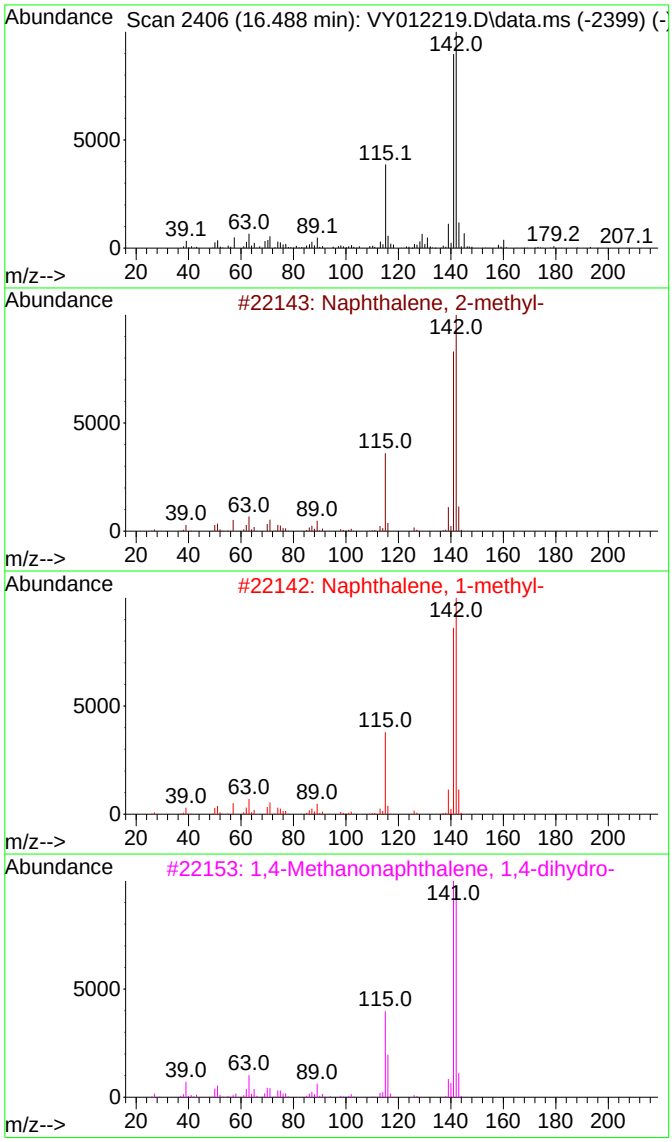
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Quant Title : SW846 8260

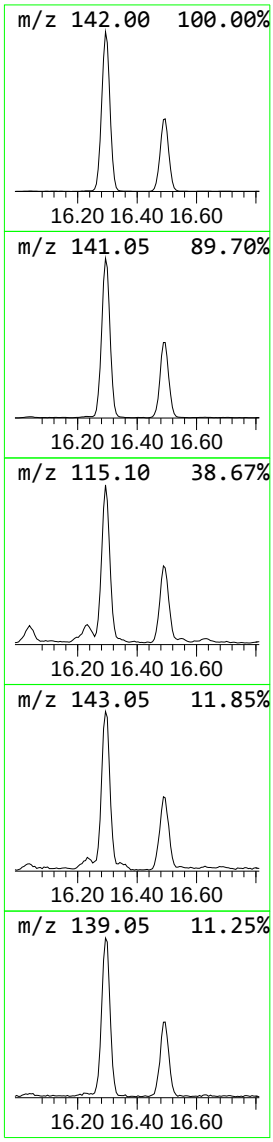
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1,4-Methanonaphthalene, 1,4... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.488	28.51 ug/l	730981	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	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	90
4	Benzocycloheptatriene	142	C11H10	000264-09-5	90
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	58





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY011723\
Data File : VY012219.D
Acq On : 17 Jan 2023 13:03
Operator : KP/MD
Sample : 01159-04
Misc : 4.60g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
MH-6

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y011623S.M
Quant Title : SW846 8260

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1-ethy...	13.971	13.1	ug/l	335461	4	13.428	1281970	50.0
Benzene, 1-ethe...	14.080	12.9	ug/l	331745	4	13.428	1281970	50.0
1H-Indene, 2,3-...	14.568	13.7	ug/l	351855	4	13.428	1281970	50.0
Benzene, 2-ethe...	14.684	33.6	ug/l	860668	4	13.428	1281970	50.0
Undecane, 2,6-d...	14.733	19.2	ug/l	492335	4	13.428	1281970	50.0
Bicyclo[4.2.0]o...	15.056	12.9	ug/l	330789	4	13.428	1281970	50.0
1H-Indene, 2,3-...	15.531	12.6	ug/l	324160	4	13.428	1281970	50.0
Benzene, (1-met...	15.732	17.1	ug/l	439452	4	13.428	1281970	50.0
Benzocyclohepta...	16.293	50.4	ug/l	1291110	4	13.428	1281970	50.0
1,4-Methanonaph...	16.488	28.5	ug/l	730981	4	13.428	1281970	50.0