LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102022\
Data File : VY011034.D
Acq On : 20 Oct 2022 18:37
Operator : KP/MD
Sample : N5193-09
Misc : 5.53g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
LSP-SS-04-15-3

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\82Y101722S.M
Title : SW846 8260

Signal : TIC: VY011034.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.863	3	7	14	rVB3	10174	16044	1.36%	0.120%
2	3.948	339	349	362	rBV	10578	28890	2.45%	0.216%
3	4.716	464	475	496	rVB5	9907	37765	3.21%	0.282%
4	5.747	634	644	654	rBV4	4880	16038	1.36%	0.120%
5	6.192	708	717	729	rBV2	5311	17032	1.45%	0.127%
6	7.716	955	967	972	rBV	124472	303454	25.77%	2.268%
7	7.789	972	979	990	rVB	386107	884724	75.13%	6.613%
8	8.136	1027	1036	1046	rBV	87929	198476	16.85%	1.484%
9	8.685	1118	1126	1140	rBV	454816	894165	75.93%	6.683%
10	8.886	1151	1159	1169	rBV	10334	24417	2.07%	0.183%
11	9.179	1192	1207	1213	rBV2	19713	50686	4.30%	0.379%
12	9.606	1271	1277	1285	rBV6	6291	12423	1.05%	0.093%
13	9.819	1306	1312	1315	rBV	19237	34726	2.95%	0.260%
14	9.959	1324	1335	1346	rBV2	23918	56701	4.81%	0.424%
15	10.179	1364	1371	1384	rBV	676313	1177637	100.00%	8.802%
16	10.520	1422	1427	1432	rBV2	12789	19710	1.67%	0.147%
17	10.606	1433	1441	1449	rBV5	20714	50267	4.27%	0.376%
18	10.709	1453	1458	1464	rVB2	16884	28947	2.46%	0.216%
19	10.801	1464	1473	1478	rBV2	17577	33932	2.88%	0.254%
20	10.904	1478	1490	1496	rBV2	23716	69069	5.87%	0.516%
21	10.965	1496	1500	1503	rBV3	8139	12238	1.04%	0.091%
22	11.032	1508	1511	1516	rVB	22666	31897	2.71%	0.238%
23	11.087	1516	1520	1525	rVB2	20343	32789	2.78%	0.245%
24	11.142	1525	1529	1534	rVB6	8890	12800	1.09%	0.096%
25	11.203	1534	1539	1543	rBV3	20583	33282	2.83%	0.249%
26	11.276	1543	1551	1562	rVB5	74794	194004	16.47%	1.450%
27	11.380	1562	1568	1577	rBV	91142	153749	13.06%	1.149%
28	11.489	1580	1586	1595	rBV	404332	652661	55.42%	4.878%
29	11.587	1598	1602	1605	rBV6	7960	14019	1.19%	0.105%
30	11.672	1611	1616	1621	rBV8	19732	39882	3.39%	0.298%
31	11.733	1621	1626	1630	rBV	33069	55869	4.74%	0.418%
32	11.776	1630	1633	1638	rVB	27102	39938	3.39%	0.299%
33	11.886	1646	1651	1654	rBV4	10057	20001	1.70%	0.149%
34	11.928	1654	1658	1663	rVB4	21838	37676	3.20%	0.282%
35	12.002	1664	1670	1676	rBV3	56318	118991	10.10%	0.889%
36	12.056	1676	1679	1682	rVB2	13997	16107	1.37%	0.120%

LSC Area Percent Report

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ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
LSP-SS-04-15-3

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\82Y101722S.M
Title : SW846 8260

37	12.124	1682	1690	1694	rVB	55380	84515	7.18%	0.632%
38	12.197	1697	1702	1705	rVV4	38517	67440	5.73%	0.504%
39	12.245	1705	1710	1714	rVB4	45137	78716	6.68%	0.588%
40	12.331	1719	1724	1726	rBV2	32755	56846	4.83%	0.425%
41	12.380	1727	1732	1733	rVV3	41341	71494	6.07%	0.534%
42	12.410	1734	1737	1743	rVB	95152	142989	12.14%	1.069%
43	12.483	1743	1749	1759	rVB3	292735	547108	46.46%	4.089%
44	12.569	1759	1763	1768	rVB4	17355	25856	2.20%	0.193%
45	12.666	1772	1779	1785	rVB	164714	273618	23.23%	2.045%
46	12.727	1785	1789	1793	rVB5	8372	13278	1.13%	0.099%
47	12.806	1795	1802	1806	rBV4	65983	143502	12.19%	1.073%
48	12.861	1807	1811	1815	rVB2	41363	61211	5.20%	0.458%
49	12.947	1816	1825	1830	rBV4	51627	117275	9.96%	0.877%
50	13.001	1830	1834	1837	rBV3	36110	58240	4.95%	0.435%
51	13.038	1837	1840	1847	rVB	86922	118874	10.09%	0.889%
52	13.099	1847	1850	1857	rVB5	32666	62292	5.29%	0.466%
53	13.166	1857	1861	1866	rBV	45126	62815	5.33%	0.470%
54	13.227	1866	1871	1872	rBV2	57141	76274	6.48%	0.570%
55	13.318	1878	1886	1890	rBV4	83042	151477	12.86%	1.132%
56	13.361	1890	1893	1897	rVV2	59222	86394	7.34%	0.646%
57	13.422	1899	1903	1909	rVB	169298	255340	21.68%	1.909%
58	13.593	1925	1931	1933	rBV	145168	226315	19.22%	1.692%
59	13.623	1933	1936	1941	rVV	169513	264226	22.44%	1.975%
60	13.696	1941	1948	1952	rVV2	145072	298736	25.37%	2.233%
61	13.751	1952	1957	1961	rVV2	99712	180957	15.37%	1.353%
62	13.788	1961	1963	1967	rVV3	46242	64891	5.51%	0.485%
63	13.849	1968	1973	1978	rVV4	34022	78013	6.62%	0.583%
64	13.910	1979	1983	1987	rVV3	31773	71608	6.08%	0.535%
65	13.971	1987	1993	1998	rVV	483230	729515	61.95%	5.453%
66	14.026	1998	2002	2006	rVV2	77838	134264	11.40%	1.004%
67	14.080	2006	2011	2016	rVB2	238349	373438	31.71%	2.791%
68	14.148	2016	2022	2027	rVB5	44812	81220	6.90%	0.607%
69	14.202	2027	2031	2032	rBV4	16100	23360	1.98%	0.175%
70	14.263	2033	2041	2043	rVV6	123746	265994	22.59%	1.988%
71	14.294	2043	2046	2054	rVB	224257	337148	28.63%	2.520%
72	14.379	2055	2060	2063	rBV	84611	133503	11.34%	0.998%
73	14.416	2063	2066	2070	rVV4	55930	88227	7.49%	0.659%
74	14.477	2073	2076	2080	rVV	42964	63899	5.43%	0.478%
75	14.513	2080	2082	2085	rVV	20937	22947	1.95%	0.172%
76	14.562	2085	2090	2096	rVV2	176689	301407	25.59%	2.253%

LSC Area Percent Report

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Instrument :
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ClientSampleId :
LSP-SS-04-15-3

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
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Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101722S.M
Title : SW846 8260

77	14.611	2096	2098	2102	rVB3	39554	47224	4.01%	0.353%
78	14.678	2102	2109	2114	rBV3	261130	537306	45.63%	4.016%
79	14.733	2114	2118	2124	rVB2	124894	215199	18.27%	1.609%
80	14.904	2143	2146	2148	rBV3	25291	36560	3.10%	0.273%
81	14.940	2148	2152	2156	rVV3	101847	162046	13.76%	1.211%
82	14.977	2156	2158	2164	rVV	28802	38966	3.31%	0.291%
83	15.056	2164	2171	2176	rVB2	71450	146412	12.43%	1.094%
84	15.141	2182	2185	2190	rVB2	31443	50774	4.31%	0.380%
85	15.208	2190	2196	2202	rBV4	36819	73331	6.23%	0.548%
86	15.342	2213	2218	2222	rBV3	30801	54058	4.59%	0.404%
87	15.525	2242	2248	2255	rBV3	27090	59770	5.08%	0.447%
88	15.599	2257	2260	2265	rVB6	14164	24300	2.06%	0.182%
89	15.684	2268	2274	2278	rBV5	14143	30288	2.57%	0.226%
90	15.812	2293	2295	2301	rVB4	9230	12841	1.09%	0.096%
91	16.025	2326	2330	2334	rVB7	8829	14063	1.19%	0.105%
92	16.153	2347	2351	2357	rBV7	6604	12957	1.10%	0.097%
93	16.287	2367	2373	2388	rVB	38320	86149	7.32%	0.644%
94	16.483	2399	2405	2412	rBV	29038	60309	5.12%	0.451%

Sum of corrected areas: 13378781

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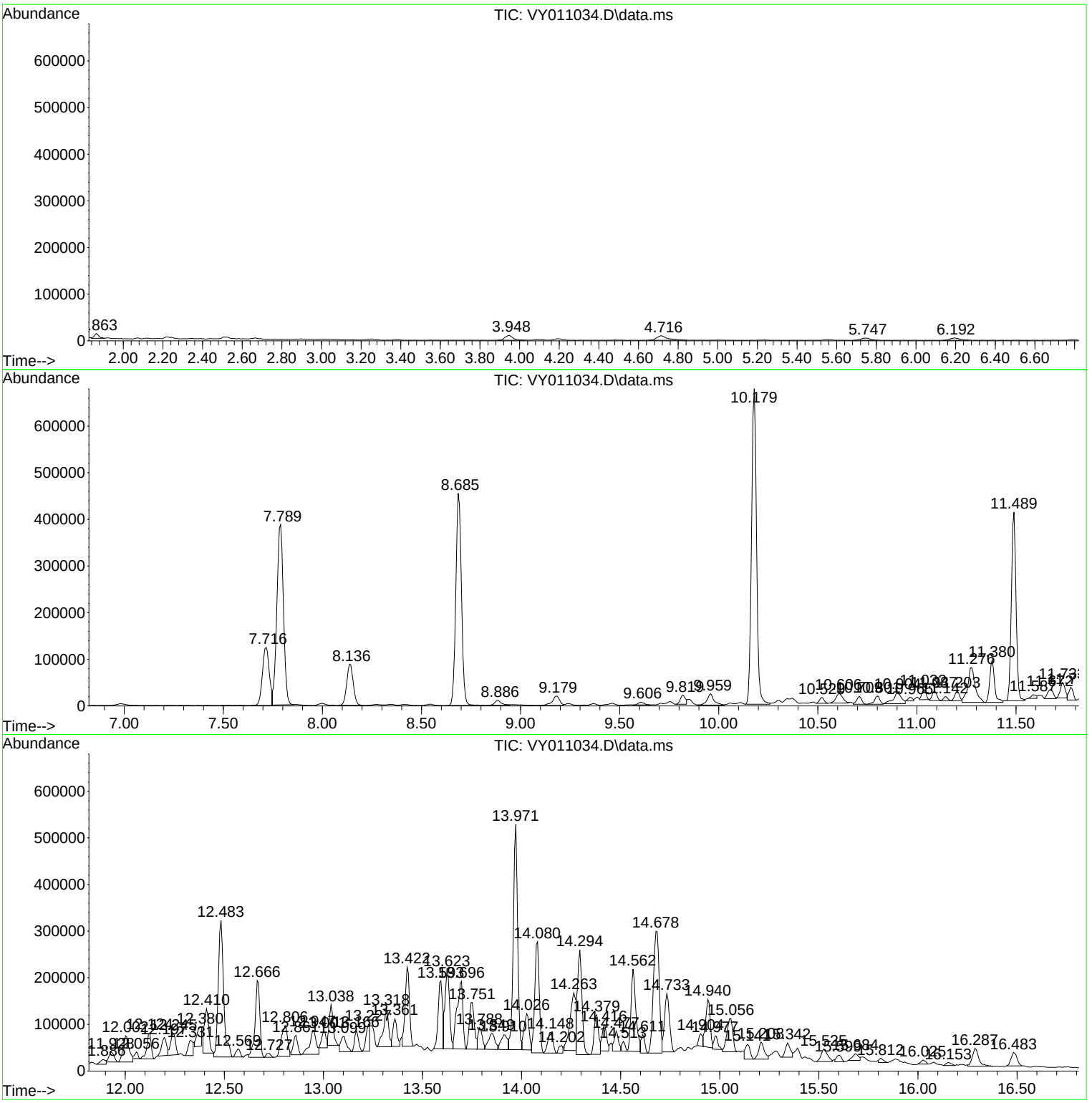
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
LSP-SS-04-15-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101722S.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102022\
Data File : VY011034.D
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MSVOA_Y
ClientSampleId :
LSP-SS-04-15-3

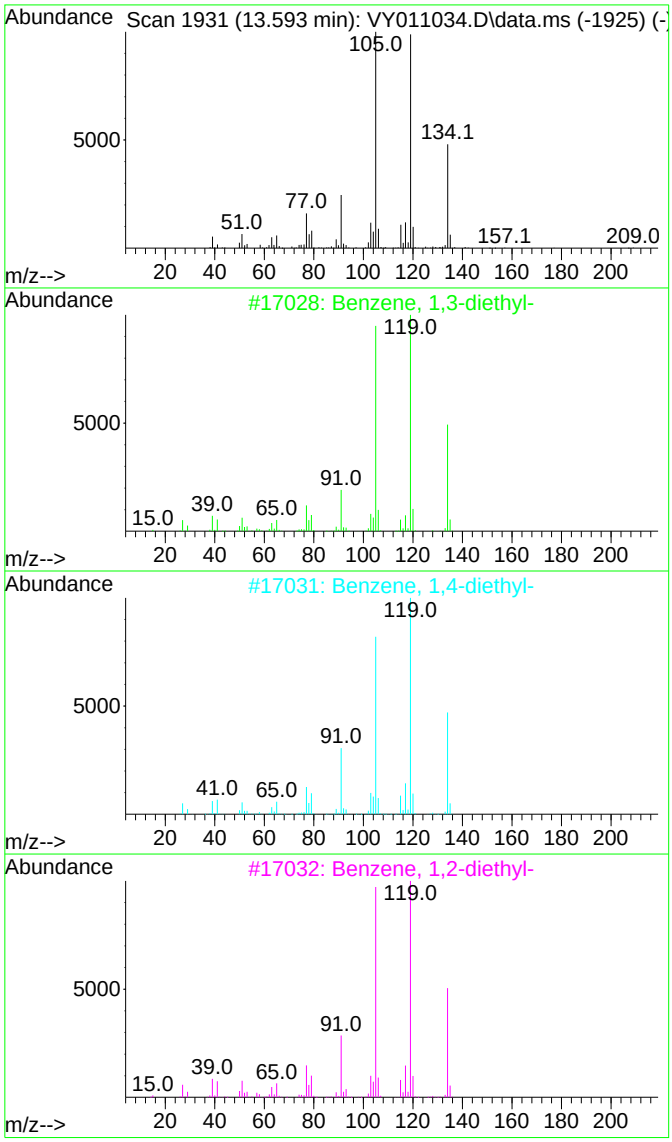
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101722S.M
Quant Title : SW846 8260

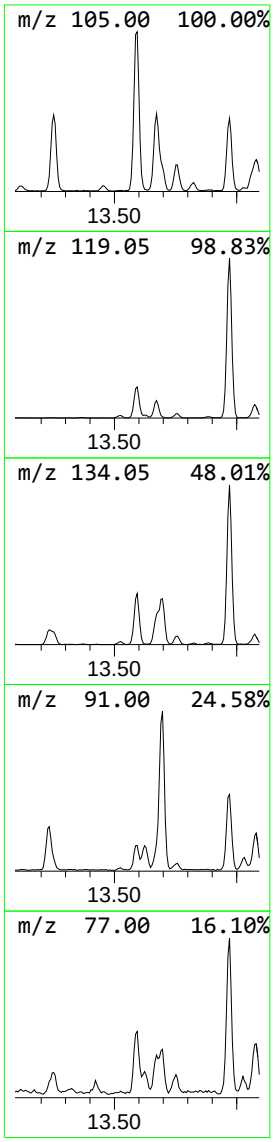
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Benzene, 1,3-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.593	44.32 ug/l	226315	1,4-Dichlorobenzene-d4	13.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
3	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102022\
Data File : VY011034.D
Acq On : 20 Oct 2022 18:37
Operator : KP/MD
Sample : N5193-09
Misc : 5.53g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
LSP-SS-04-15-3

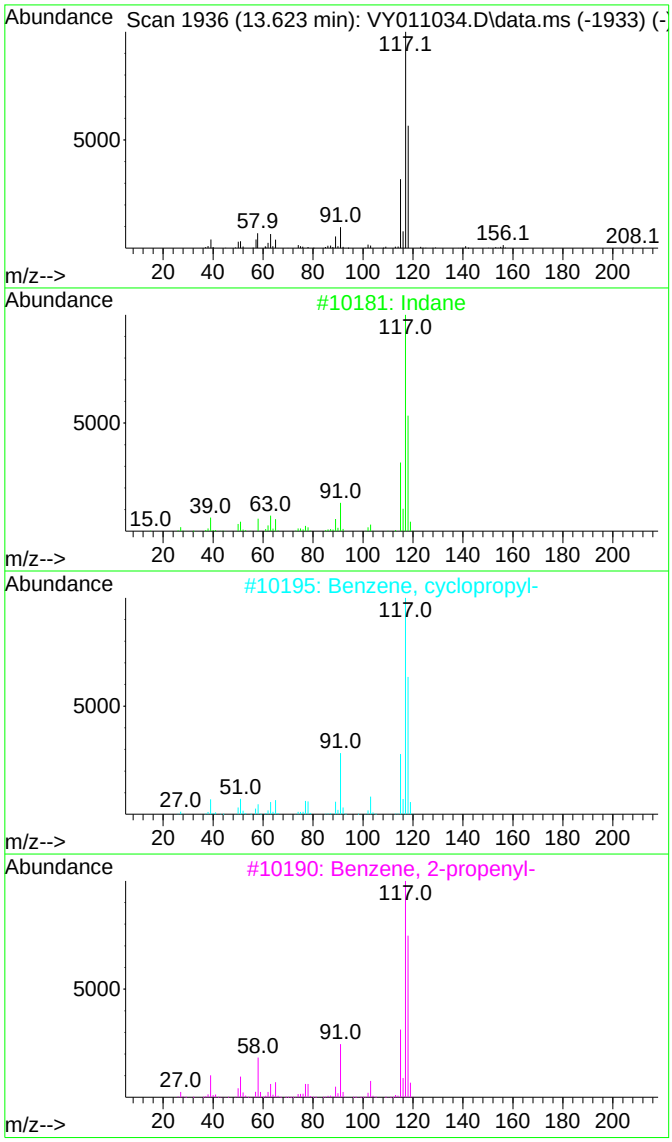
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101722S.M
Quant Title : SW846 8260

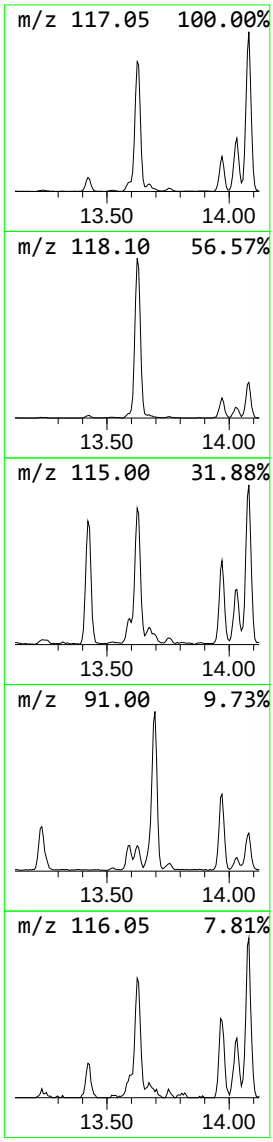
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Indane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.623	51.74 ug/l	264226	1,4-Dichlorobenzene-d4	13.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	81
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
4	Indane-5-carboxylic acid	162	C10H10O2	065898-38-6	59
5	Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	50





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Data File : VY011034.D
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Instrument :
MSVOA_Y
ClientSampleId :
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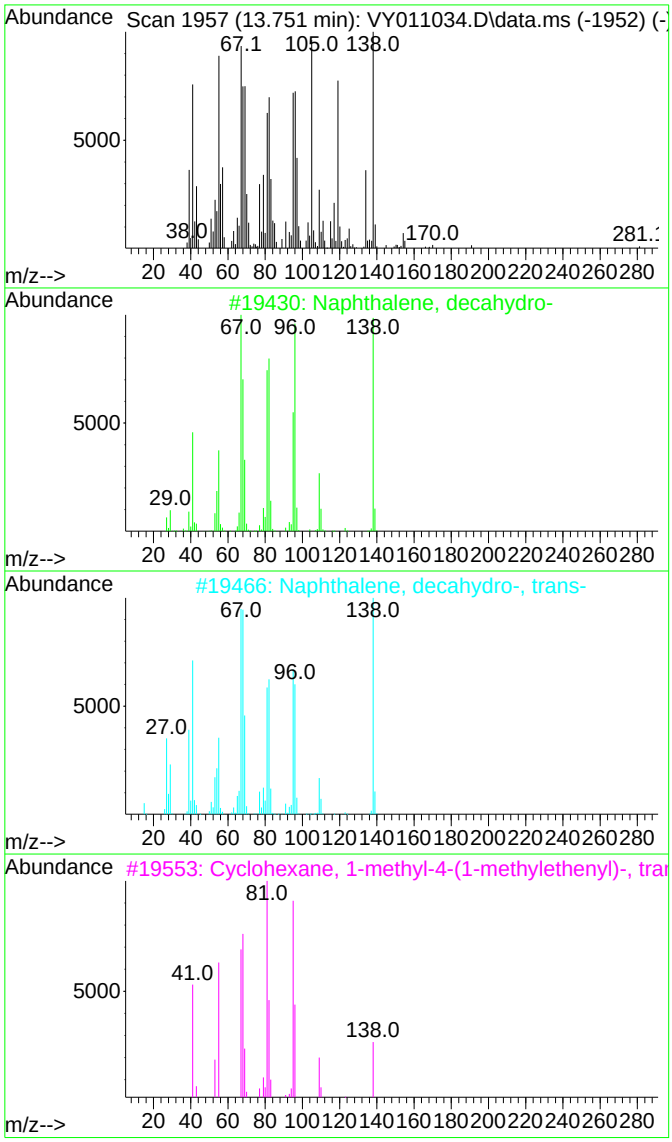
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101722S.M
Quant Title : SW846 8260

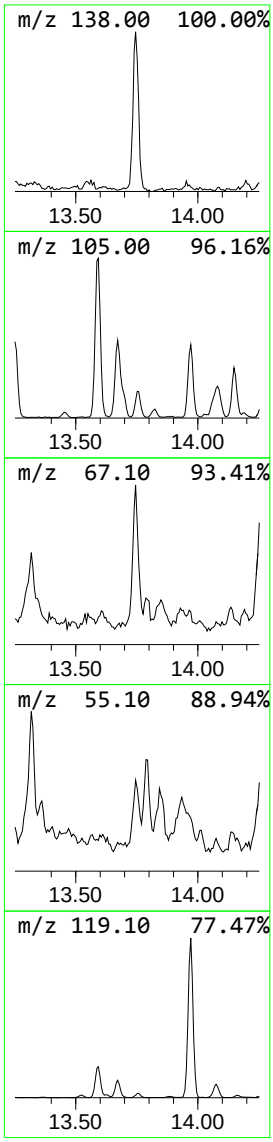
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Naphthalene, decahydro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.751	35.43 ug/l	180957	1,4-Dichlorobenzene-d4	13.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, decahydro-	138	C10H18	000091-17-8	95
2	Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	91
3	Cyclohexane, 1-methyl-4-(1-methy...	138	C10H18	001124-25-0	86
4	9-Borabicyclo[3.3.1]nonane, 9-hy...	138	C8H15BO	063366-65-4	58
5	1,1'-Bicyclopentyl	138	C10H18	001636-39-1	50





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Instrument :
MSVOA_Y
ClientSampleId :
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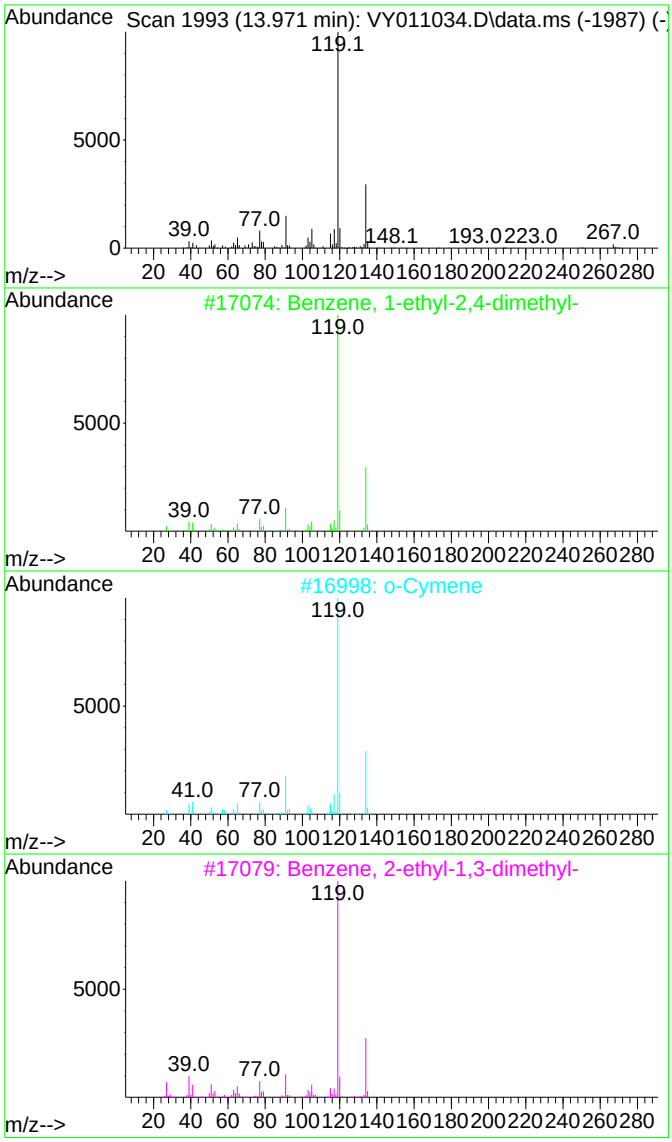
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101722S.M
Quant Title : SW846 8260

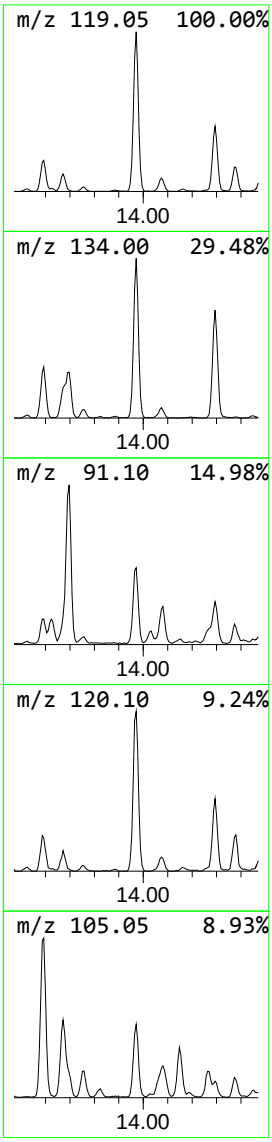
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.971	142.85 ug/l	729515	1,4-Dichlorobenzene-d4	13.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95





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Data File : VY011034.D
Acq On : 20 Oct 2022 18:37
Operator : KP/MD
Sample : N5193-09
Misc : 5.53g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
LSP-SS-04-15-3

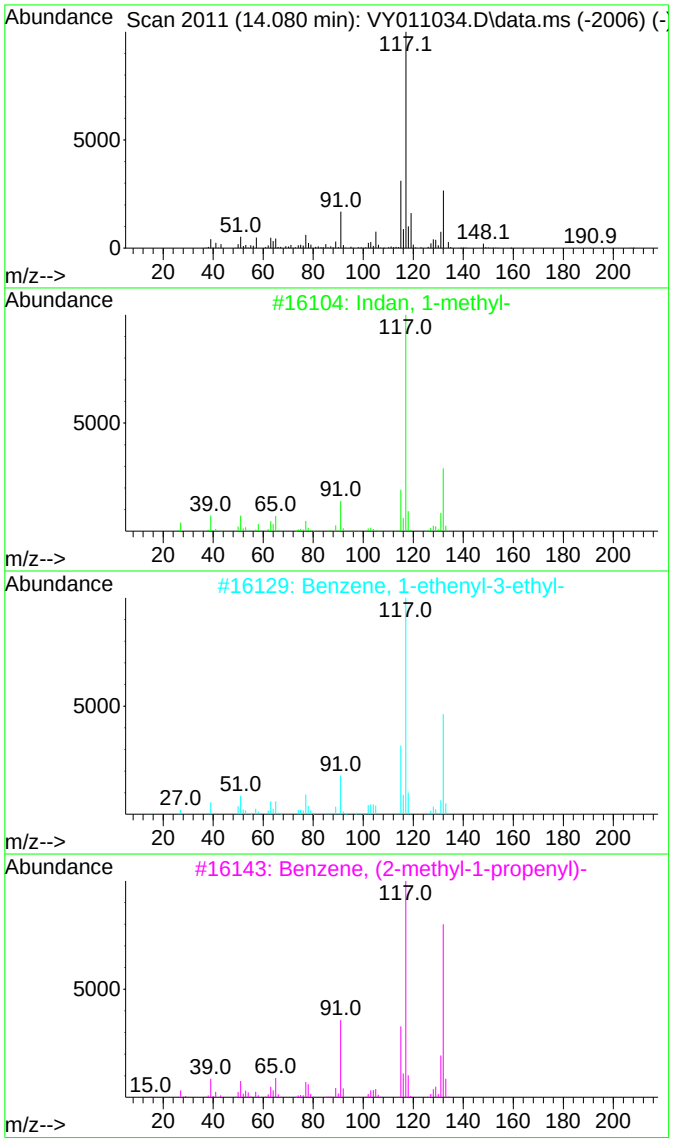
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101722S.M
Quant Title : SW846 8260

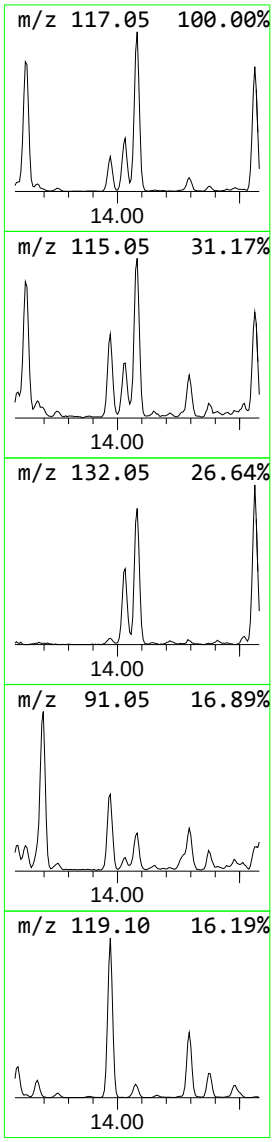
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.081	73.13 ug/l	373438	1,4-Dichlorobenzene-d4	13.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
4	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	80
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102022\
Data File : VY011034.D
Acq On : 20 Oct 2022 18:37
Operator : KP/MD
Sample : N5193-09
Misc : 5.53g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
LSP-SS-04-15-3

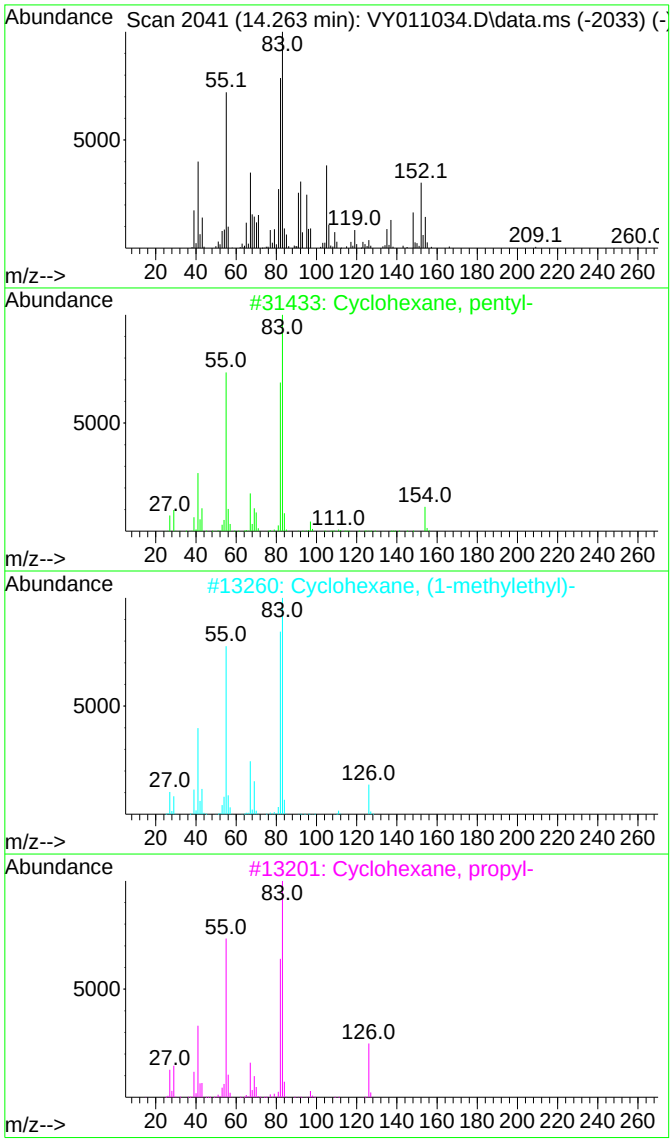
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101722S.M
Quant Title : SW846 8260

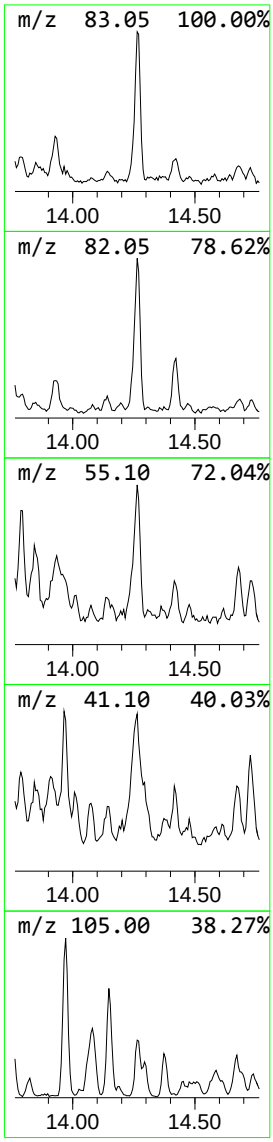
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Cyclohexane, pentyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.263	52.09 ug/l	265994	1,4-Dichlorobenzene-d4	13.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, pentyl-	154	C11H22	004292-92-6	91
2	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	60
3	Cyclohexane, propyl-	126	C9H18	001678-92-8	49
4	Cyclohexane, octyl-	196	C14H28	001795-15-9	47
5	n-Heptadecylcyclohexane	322	C23H46	019781-73-8	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102022\
Data File : VY011034.D
Acq On : 20 Oct 2022 18:37
Operator : KP/MD
Sample : N5193-09
Misc : 5.53g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
LSP-SS-04-15-3

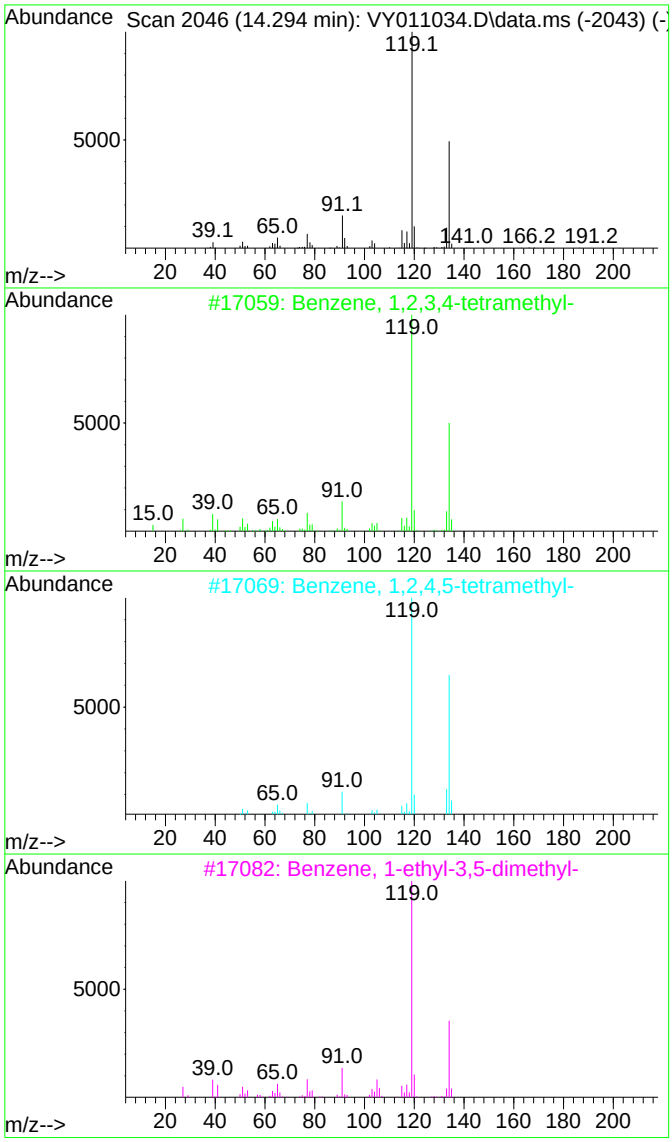
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101722S.M
Quant Title : SW846 8260

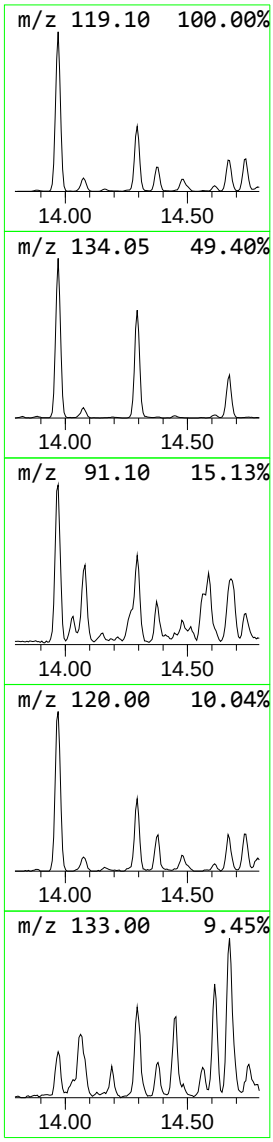
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.294	66.02 ug/l	337148	1,4-Dichlorobenzene-d4	13.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91
3	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102022\
Data File : VY011034.D
Acq On : 20 Oct 2022 18:37
Operator : KP/MD
Sample : N5193-09
Misc : 5.53g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
LSP-SS-04-15-3

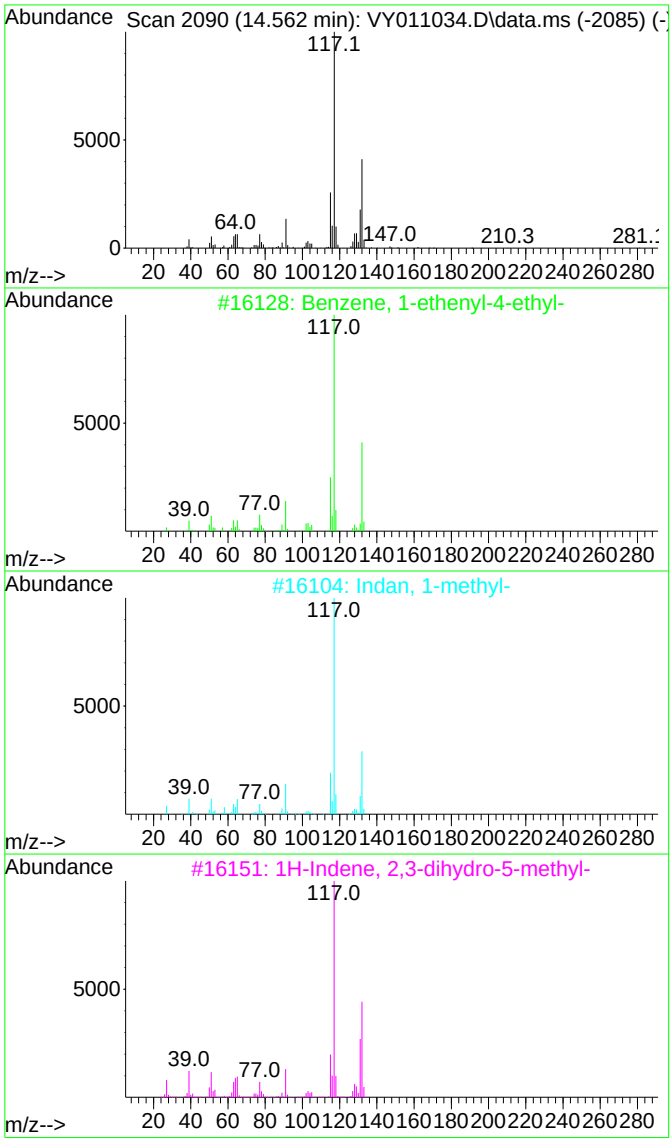
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101722S.M
Quant Title : SW846 8260

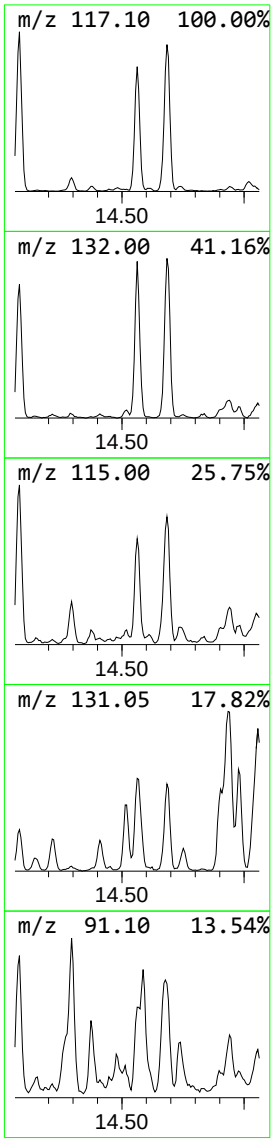
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.562	59.02 ug/l	301407	1,4-Dichlorobenzene-d4	13.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	Indan, 1-methyl-	132	C10H12	000767-58-8	93
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
4	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102022\
Data File : VY011034.D
Acq On : 20 Oct 2022 18:37
Operator : KP/MD
Sample : N5193-09
Misc : 5.53g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
LSP-SS-04-15-3

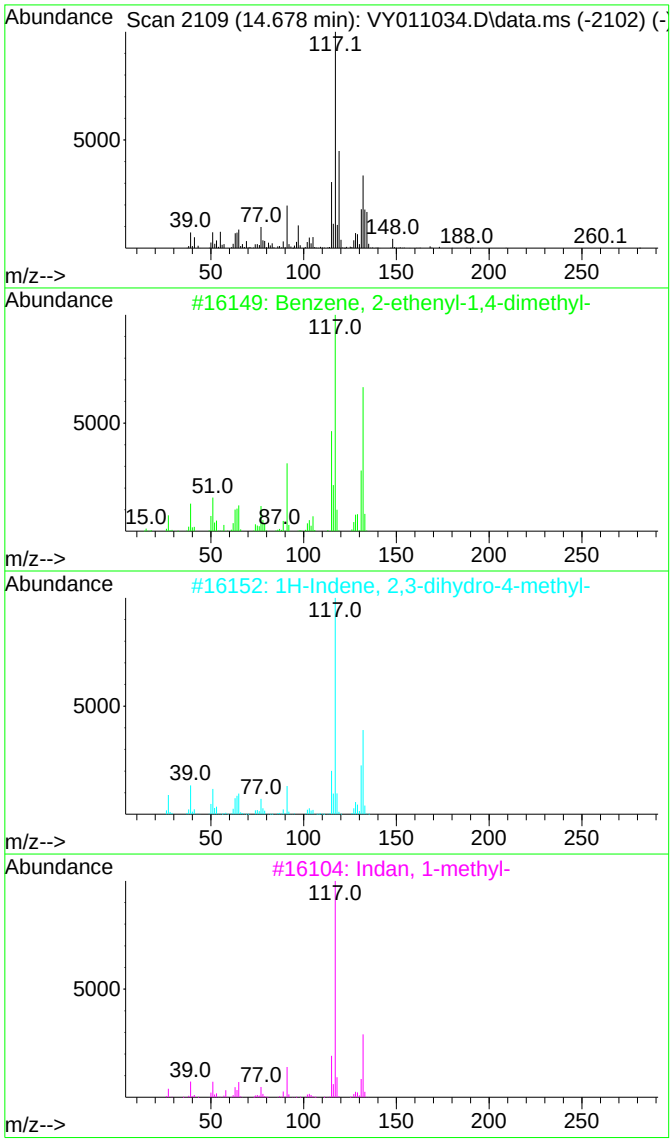
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101722S.M
Quant Title : SW846 8260

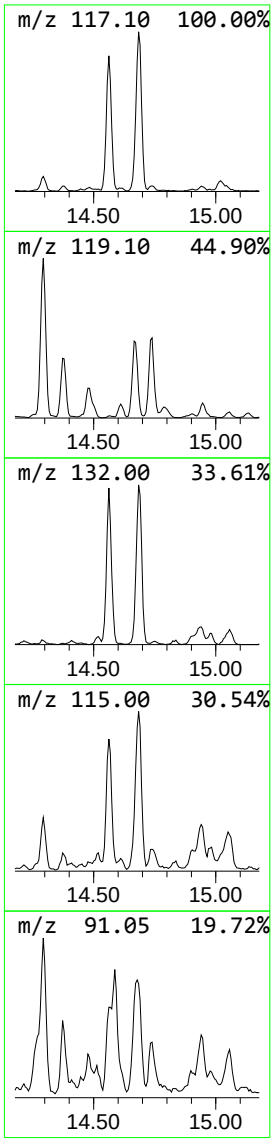
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.678	105.21 ug/l	537306	1,4-Dichlorobenzene-d4	13.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70
3	Indan, 1-methyl-	132	C10H12	000767-58-8	70
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102022\
Data File : VY011034.D
Acq On : 20 Oct 2022 18:37
Operator : KP/MD
Sample : N5193-09
Misc : 5.53g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
LSP-SS-04-15-3

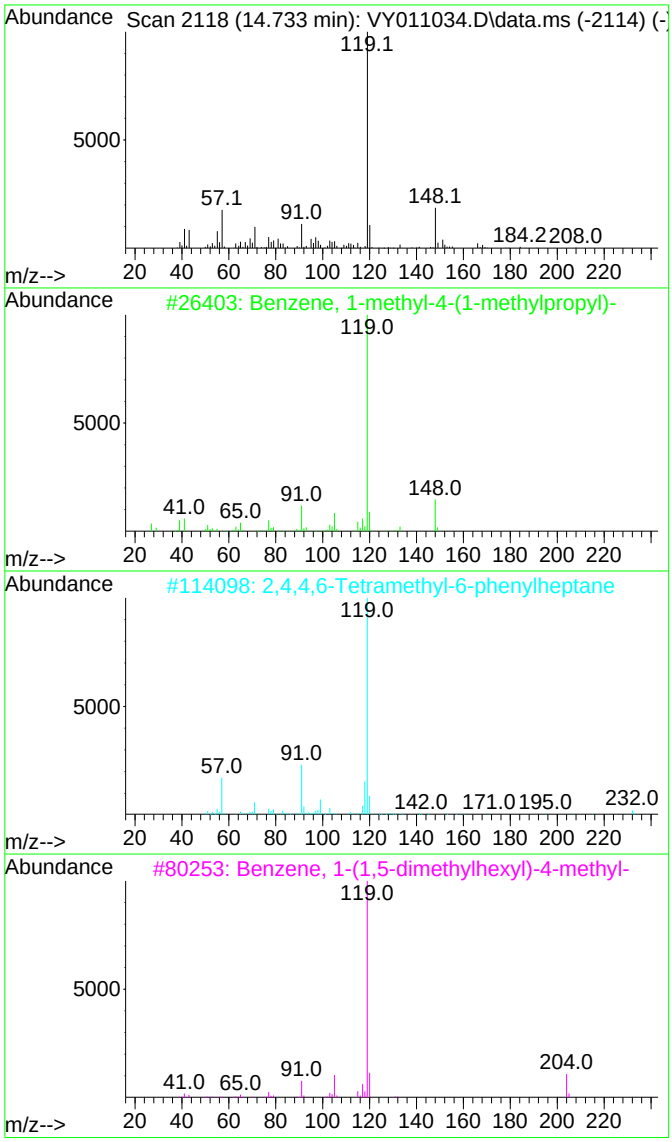
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101722S.M
Quant Title : SW846 8260

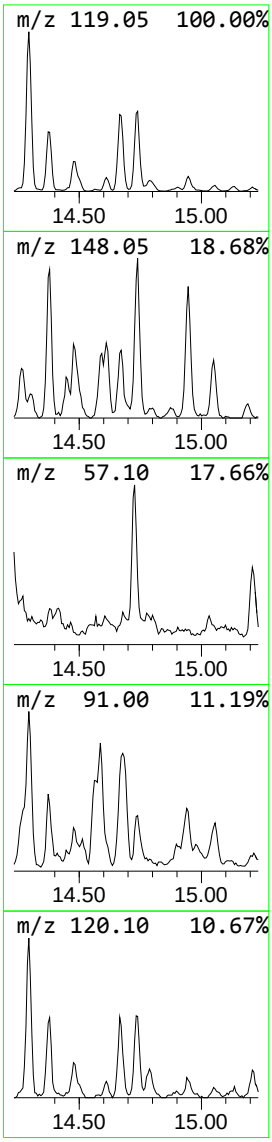
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-methyl-4-(1-meth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.733	42.14 ug/l	215199	1,4-Dichlorobenzene-d4	13.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	72
2	2,4,4,6-Tetramethyl-6-phenylheptane	232	C17H28	1010293-28-6	59
3	Benzene, 1-(1,5-dimethylhexyl)-4...	204	C15H24	001461-02-5	53
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	53
5	m-Toluic acid, undec-2-enyl ester	288	C19H28O2	1000292-61-3	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY102022\
Data File : VY011034.D
Acq On : 20 Oct 2022 18:37
Operator : KP/MD
Sample : N5193-09
Misc : 5.53g/5.0mL/MSVOA_Y/SOIL
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
LSP-SS-04-15-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y101722S.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, 1,3-di...	13.593	44.3	ug/l	226315	4	13.422	255340	50.0
Indane	13.623	51.7	ug/l	264226	4	13.422	255340	50.0
Naphthalene, de...	13.751	35.4	ug/l	180957	4	13.422	255340	50.0
Benzene, 1-ethy...	13.971	142.8	ug/l	729515	4	13.422	255340	50.0
Indan, 1-methyl-	14.081	73.1	ug/l	373438	4	13.422	255340	50.0
Cyclohexane, pe...	14.263	52.1	ug/l	265994	4	13.422	255340	50.0
Benzene, 1,2,3,...	14.294	66.0	ug/l	337148	4	13.422	255340	50.0
Benzene, 1-ethe...	14.562	59.0	ug/l	301407	4	13.422	255340	50.0
Benzene, 2-ethe...	14.678	105.2	ug/l	537306	4	13.422	255340	50.0
Benzene, 1-meth...	14.733	42.1	ug/l	215199	4	13.422	255340	50.0