LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX110918\
Data File : VX005889.D
Acq On : 10 Nov 2018 04:24
Operator : JC/MD
Sample : J5889-11
Misc : 12.72g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BOTTOM-B

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_X\METHOD\82X110718W.M
Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.659	731	739	761	rVB2	124518	399537	1.45%	0.178%
2	6.860	926	936	959	rBV	173216	461163	1.67%	0.206%
3	7.451	1022	1033	1044	rBV2	238497	554063	2.01%	0.247%
4	8.665	1224	1232	1235	rBV	288641	524120	1.90%	0.234%
5	8.713	1235	1240	1254	rVB2	377224	762659	2.76%	0.341%
6	9.037	1288	1293	1304	rVB	249181	415808	1.51%	0.186%
7	9.622	1384	1389	1393	rVB	353313	495825	1.80%	0.221%
8	9.677	1393	1398	1402	rBV	321949	477620	1.73%	0.213%
9	9.890	1426	1433	1440	rBV2	221616	489365	1.77%	0.219%
10	10.116	1465	1470	1475	rBV	345850	485193	1.76%	0.217%
11	10.256	1486	1493	1498	rBV	624202	998915	3.62%	0.446%
12	10.366	1507	1511	1516	rVV2	636015	1148477	4.16%	0.513%
13	10.414	1516	1519	1530	rVB	358117	520169	1.89%	0.232%
14	10.640	1549	1556	1564	rBV3	293713	630580	2.29%	0.282%
15	10.835	1579	1588	1592	rBV2	485902	817642	2.96%	0.365%
16	10.908	1592	1600	1604	rVV2	556083	914352	3.31%	0.408%
17	11.018	1612	1618	1622	rBV	314749	444706	1.61%	0.199%
18	11.140	1632	1638	1644	rBV2	461354	786257	2.85%	0.351%
19	11.274	1657	1660	1669	rVB3	477652	720732	2.61%	0.322%
20	11.359	1670	1674	1678	rBV	651929	867737	3.15%	0.388%
21	11.439	1682	1687	1693	rVV	1623729	3300691	11.97%	1.474%
22	11.512	1693	1699	1707	rVB	1825670	3005111	10.89%	1.342%
23	11.670	1720	1725	1732	rVB	1401730	1945494	7.05%	0.869%
24	11.768	1738	1741	1744	rBV	334376	380376	1.38%	0.170%
25	11.810	1744	1748	1753	rVV	6515286	7808346	28.31%	3.487%
26	11.951	1767	1771	1775	rVB	505629	724871	2.63%	0.324%
27	11.999	1775	1779	1781	rBV	456817	550860	2.00%	0.246%
28	12.030	1781	1784	1787	rVB	621111	720668	2.61%	0.322%
29	12.073	1787	1791	1797	rVB3	575928	1090107	3.95%	0.487%
30	12.146	1797	1803	1809	rBV	2462263	3156271	11.44%	1.410%
31	12.225	1812	1816	1824	rVB4	254139	562878	2.04%	0.251%
32	12.304	1824	1829	1831	rBV	1628271	2374290	8.61%	1.060%
33	12.323	1831	1832	1836	rVB	1464077	1385890	5.02%	0.619%
34	12.371	1836	1840	1847	rVB3	2832764	4534217	16.44%	2.025%

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ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BOTTOM-B

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_X\METHOD\82X110718W.M
Title : SW846 8260

35	12.463	1851	1855	1860	rBV3	213386	409096	1.48%	0.183%
36	12.524	1860	1865	1870	rVB	931467	1336515	4.85%	0.597%
37	12.591	1870	1876	1877	rBV	1244933	1456803	5.28%	0.651%
38	12.615	1877	1880	1885	rVB	1398296	1650068	5.98%	0.737%
39	12.670	1885	1889	1892	rBV	1896093	2087090	7.57%	0.932%
40	12.700	1892	1894	1898	rVB	498838	588610	2.13%	0.263%
41	12.761	1898	1904	1911	rBV4	1399790	3190510	11.57%	1.425%
42	12.853	1915	1919	1921	rVV2	324197	374960	1.36%	0.167%
43	12.877	1921	1923	1925	rVV2	338478	421672	1.53%	0.188%
44	12.902	1925	1927	1931	rVB2	772323	915047	3.32%	0.409%
45	12.944	1931	1934	1936	rBV2	353391	476805	1.73%	0.213%
46	12.981	1936	1940	1943	rVB	1412379	1589879	5.76%	0.710%
47	13.024	1943	1947	1953	rVB	2138699	3087613	11.19%	1.379%
48	13.085	1953	1957	1959	rBV2	565229	822120	2.98%	0.367%
49	13.194	1972	1975	1977	rVV	483541	633300	2.30%	0.283%
50	13.225	1977	1980	1984	rVV3	1278990	1769118	6.41%	0.790%
51	13.267	1984	1987	1991	rVB3	409229	475747	1.72%	0.212%
52	13.310	1991	1994	1997	rBV2	575604	742767	2.69%	0.332%
53	13.353	1997	2001	2006	rVV2	4230437	6140275	22.26%	2.742%
54	13.426	2010	2013	2017	rVB	449136	520047	1.89%	0.232%
55	13.481	2018	2022	2031	rVB2	2213299	3384570	12.27%	1.512%
56	13.566	2031	2036	2039	rBV2	764603	1096391	3.97%	0.490%
57	13.603	2039	2042	2045	rVV	930090	1159497	4.20%	0.518%
58	13.645	2045	2049	2053	rVV2	1566982	2636793	9.56%	1.178%
59	13.700	2053	2058	2059	rVV4	615017	1040981	3.77%	0.465%
60	13.725	2059	2062	2067	rVB2	1230536	1841496	6.68%	0.822%
61	13.841	2072	2081	2087	rVB	5291251	7334676	26.59%	3.276%
62	13.914	2088	2093	2098	rVB	1615847	2187196	7.93%	0.977%
63	13.987	2098	2105	2109	rBV3	1797256	2840280	10.30%	1.268%
64	14.030	2109	2112	2115	rBV2	460032	516504	1.87%	0.231%
65	14.066	2115	2118	2124	rVB3	484927	813512	2.95%	0.363%
66	14.133	2124	2129	2133	rBV2	1164078	1596529	5.79%	0.713%
67	14.225	2141	2144	2147	rVB	270838	294214	1.07%	0.131%
68	14.292	2147	2155	2161	rBV3	2824211	5477730	19.86%	2.446%
69	14.383	2166	2170	2174	rVV	1043102	1341858	4.86%	0.599%
70	14.432	2174	2178	2182	rVB	1521933	1828212	6.63%	0.816%
71	14.481	2182	2186	2192	rVB	614645	927812	3.36%	0.414%

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Instrument : MSVOA_X
ClientSampleId : BOTTOM-B

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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Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Title : SW846 8260

72	14.542	2192	2196	2202	rBV2	2036725	3072390	11.14%	1.372%
73	14.706	2215	2223	2227	rBV	18381752	27584295	100.00%	12.319%
74	14.737	2227	2228	2232	rVB2	1506648	1223801	4.44%	0.547%
75	14.792	2232	2237	2241	rBV2	708225	1074401	3.89%	0.480%
76	14.846	2241	2246	2250	rVV	10985570	13989941	50.72%	6.248%
77	14.883	2250	2252	2254	rVV2	771119	873739	3.17%	0.390%
78	14.907	2254	2256	2261	rVV2	856927	1006551	3.65%	0.450%
79	14.968	2261	2266	2272	rVV4	484187	1091217	3.96%	0.487%
80	15.042	2273	2278	2283	rVB5	345798	805078	2.92%	0.360%
81	15.115	2283	2290	2294	rBV2	388441	806111	2.92%	0.360%
82	15.255	2307	2313	2315	rBV	1504794	2398646	8.70%	1.071%
83	15.322	2320	2324	2332	rVB6	368943	872235	3.16%	0.390%
84	15.450	2340	2345	2348	rBV	3690166	5228159	18.95%	2.335%
85	15.480	2348	2350	2355	rVB	1050072	1363963	4.94%	0.609%
86	15.560	2355	2363	2368	rBV	9737250	14935989	54.15%	6.670%
87	15.608	2368	2371	2376	rVV2	273621	651011	2.36%	0.291%
88	15.700	2380	2386	2390	rVV	8884290	15690476	56.88%	7.007%
89	15.737	2390	2392	2399	rVB	5839901	7398785	26.82%	3.304%
90	15.919	2413	2422	2433	rVB	2230812	5378244	19.50%	2.402%
91	16.072	2442	2447	2458	rVB	1185057	2068203	7.50%	0.924%
92	16.175	2458	2464	2469	rBV	390288	626573	2.27%	0.280%
93	16.279	2472	2481	2488	rVB2	475845	1144335	4.15%	0.511%
94	16.352	2488	2493	2496	rBV	224383	395476	1.43%	0.177%
95	16.444	2502	2508	2516	rVB	860904	1603925	5.81%	0.716%
96	16.517	2516	2520	2524	rVB	247784	383766	1.39%	0.171%
97	16.566	2524	2528	2534	rVB	211509	368401	1.34%	0.165%
98	16.657	2534	2543	2545	rBV4	253276	582710	2.11%	0.260%
99	16.694	2545	2549	2554	rVB	543408	955885	3.47%	0.427%
100	16.755	2554	2559	2564	rBV	499616	882626	3.20%	0.394%

Sum of corrected areas: 223918215

LSC Report - Integrated Chromatogram

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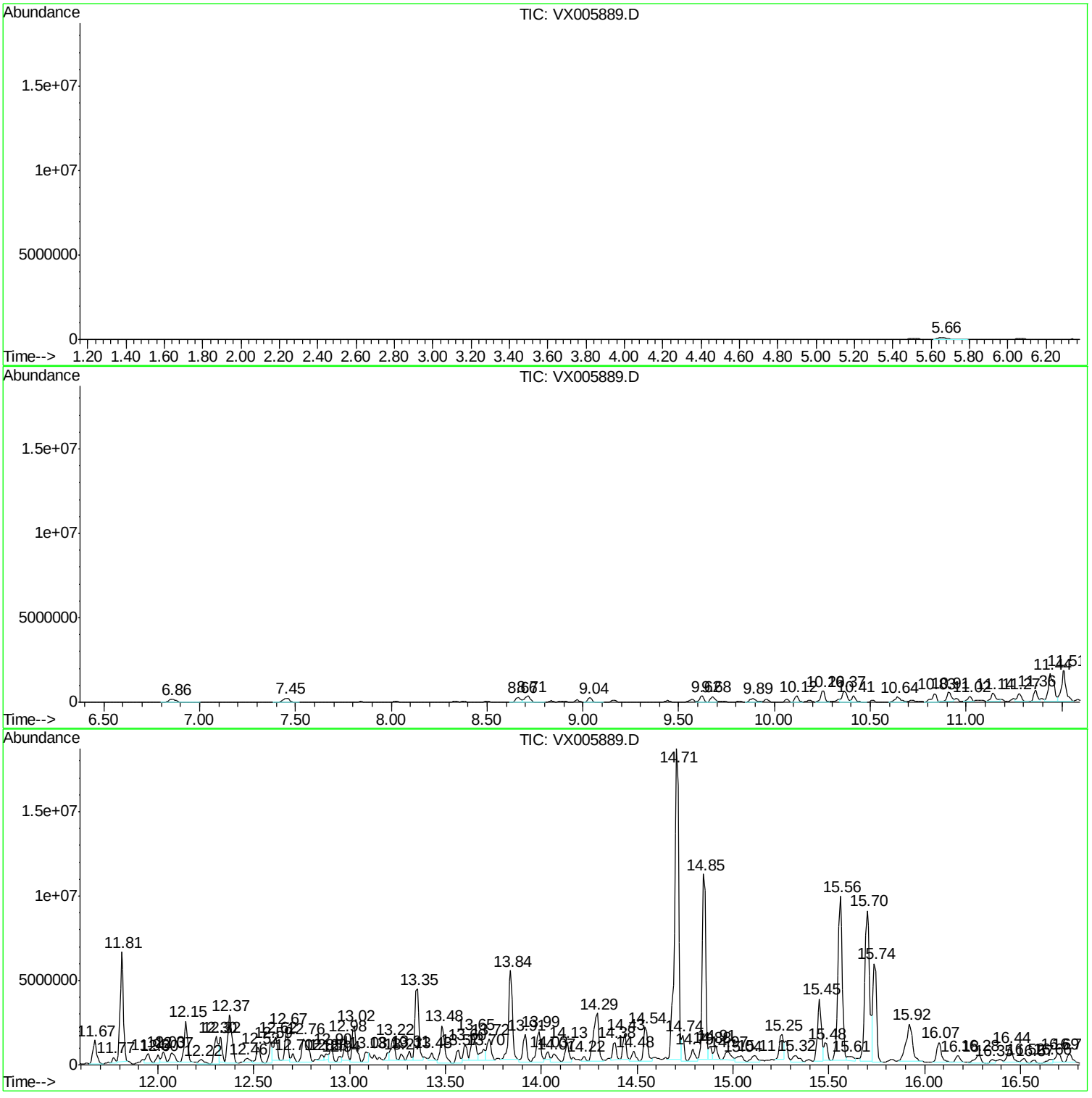
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
BOTTOM-B

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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



Library Search Compound Report

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ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BOTTOM-B

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

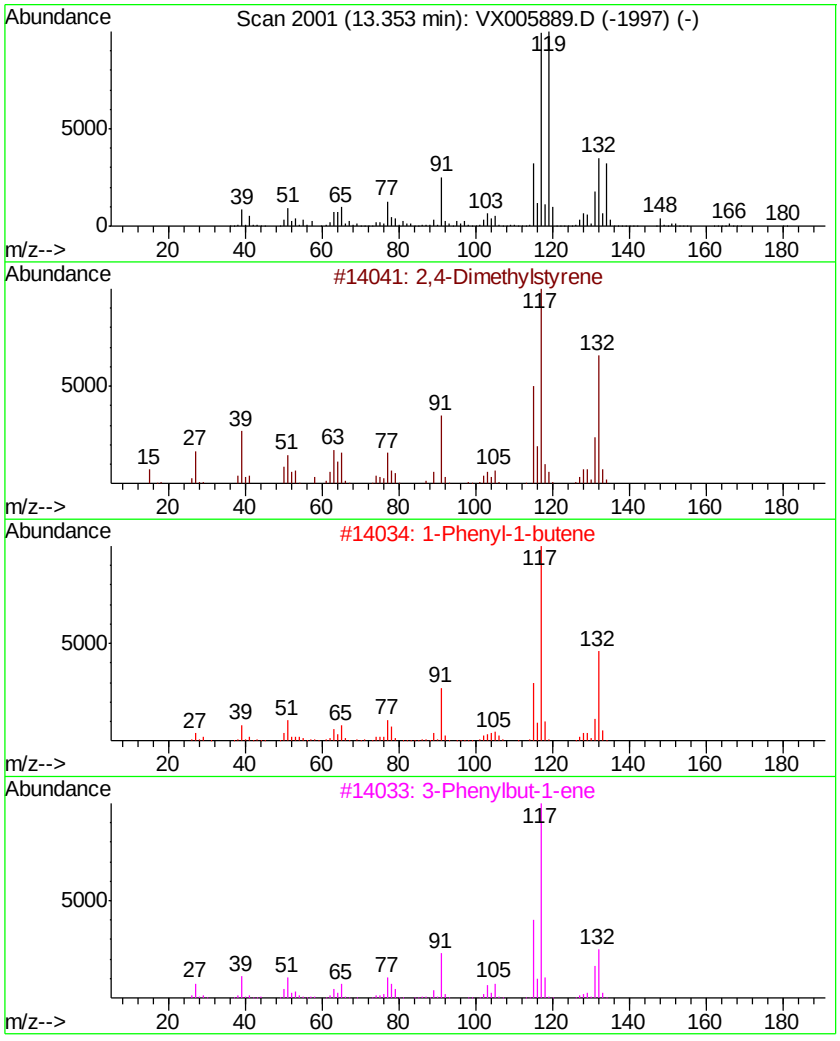
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

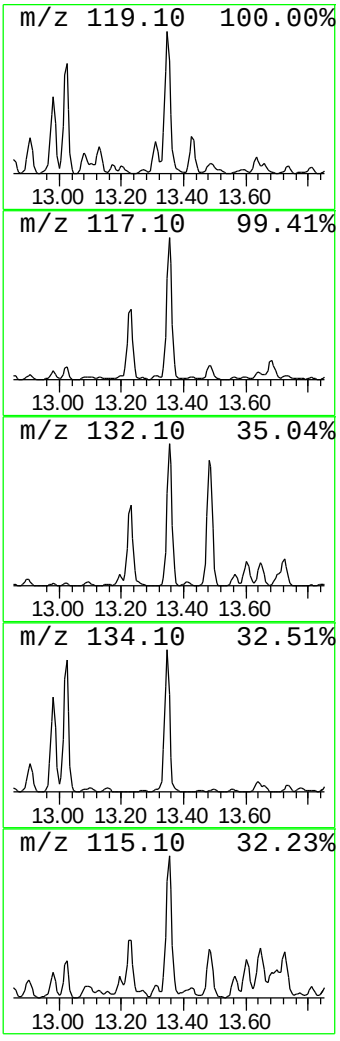
****

Peak Number 1 2,4-Dimethylstyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	281.64 ug/l	6140280	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dimethylstyrene	132	C10H12	002234-20-0	93
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	91
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	89
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	86





Library Search Compound Report

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ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BOTTOM-B

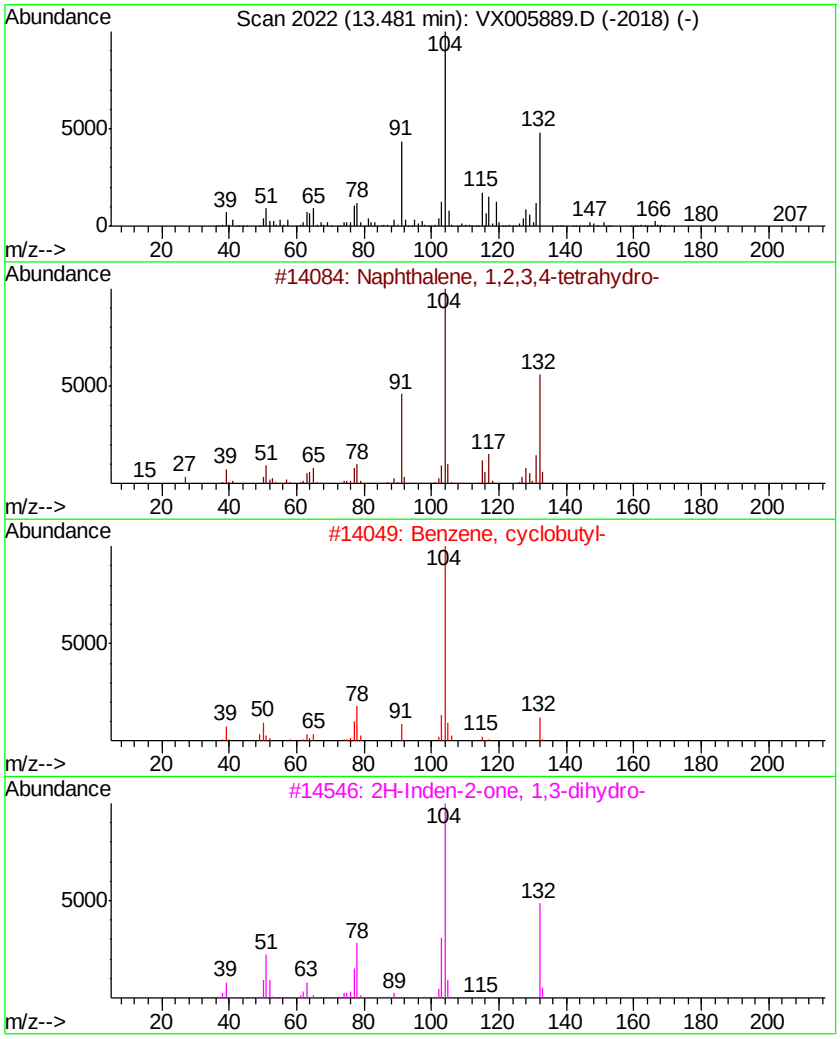
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

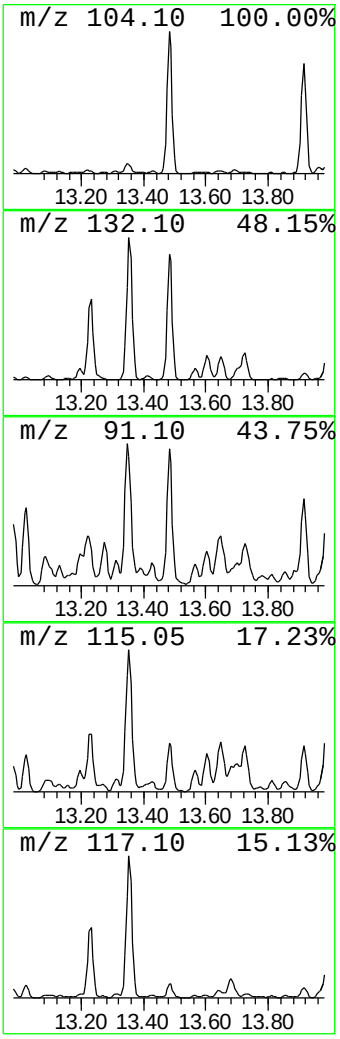
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.48	155.24 ug/l	3384570	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2	Benzene, cyclobutyl-	132	C10H12	004392-30-7	53
3	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	50
4	Naphthalene, 2-ethyl-1,2,3,4-tet...	160	C12H16	032367-54-7	47
5	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	46





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Instrument : MSVOA_X
ClientSampleId : BOTTOM-B

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

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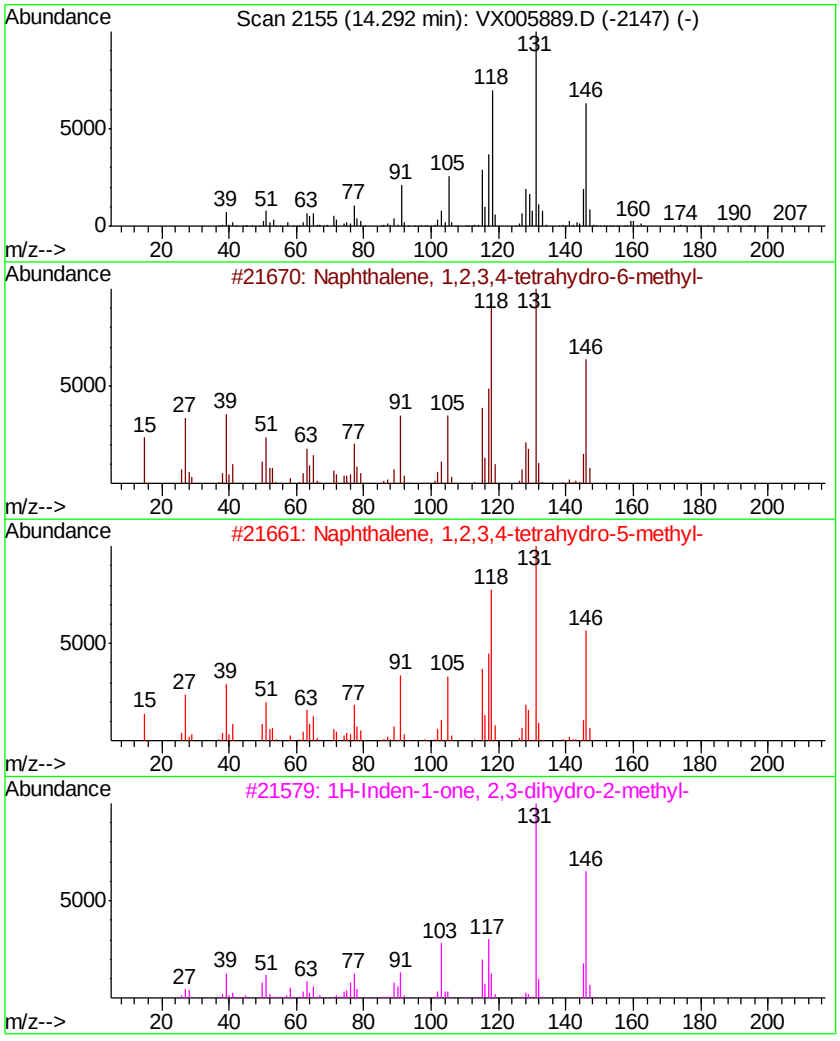
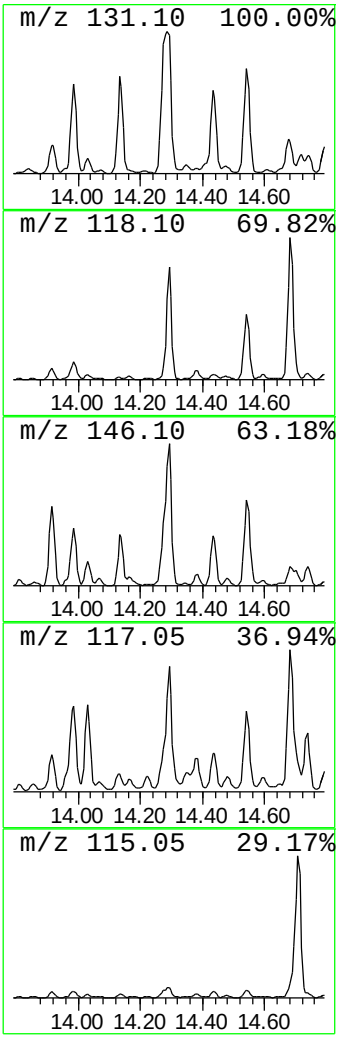
Peak Number 3 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.29	251.25 ug/l	5477730	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	96
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
3	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	76
4	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	70
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70





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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

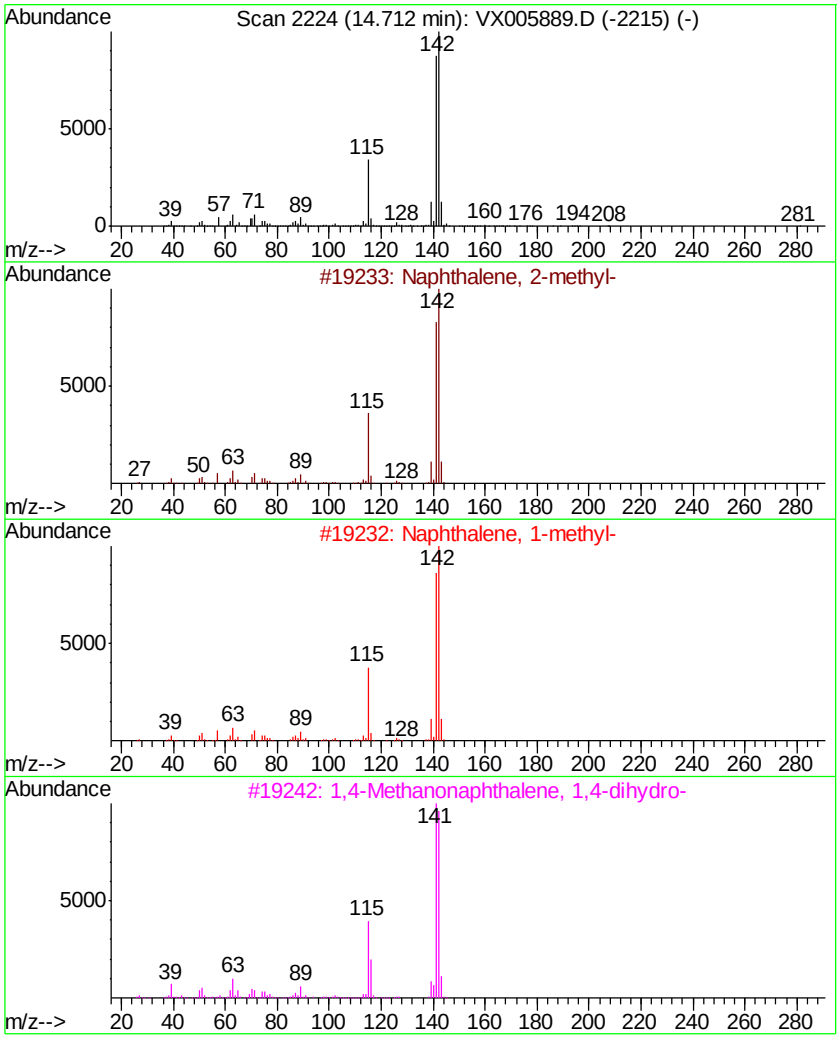
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

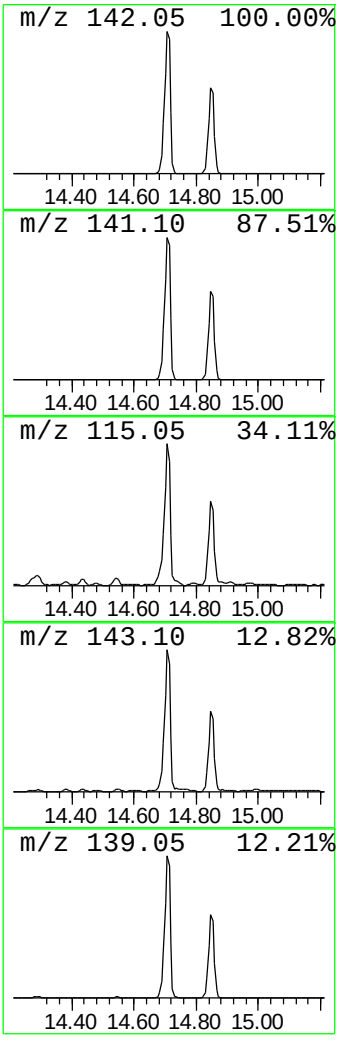
****

Peak Number 4 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.71	1265.21 ug/l	27584300	1,4-Dichlorobenzene-d4	12.08

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	93
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91





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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

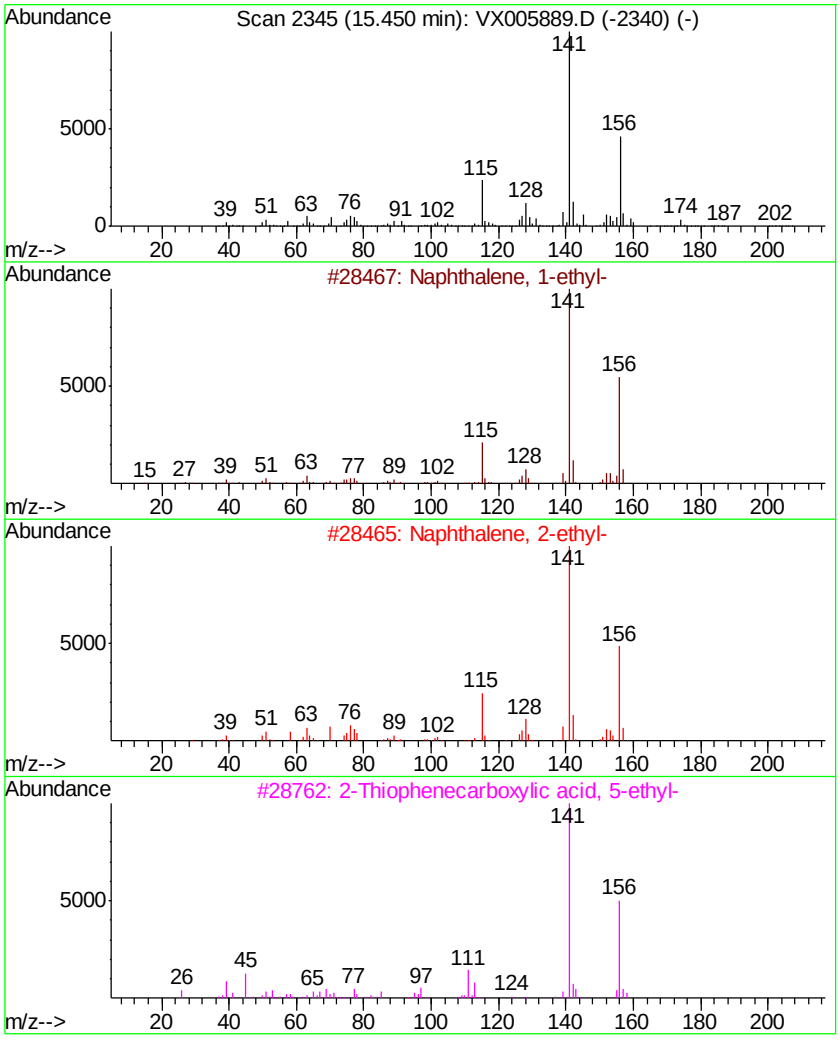
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

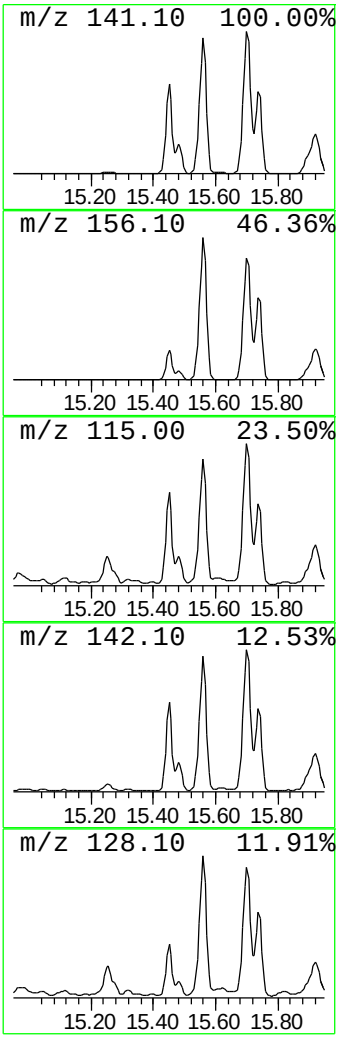
****

Peak Number 6 Naphthalene, 1-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.45	239.80 ug/l	5228160	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	96
2	Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	96
3	2-Thiophenecarboxylic acid, 5-et...	156	C7H8O2S	023229-72-3	64
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	53
5	1-Ethanone, 1-[4-(1-naphthalenyl...	276	C19H16O2	1000338-30-5	47





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX110918\
Data File : VX005889.D
Acq On : 10 Nov 2018 04:24
Operator : JC/MD
Sample : J5889-11
Misc : 12.72g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BOTTOM-B

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

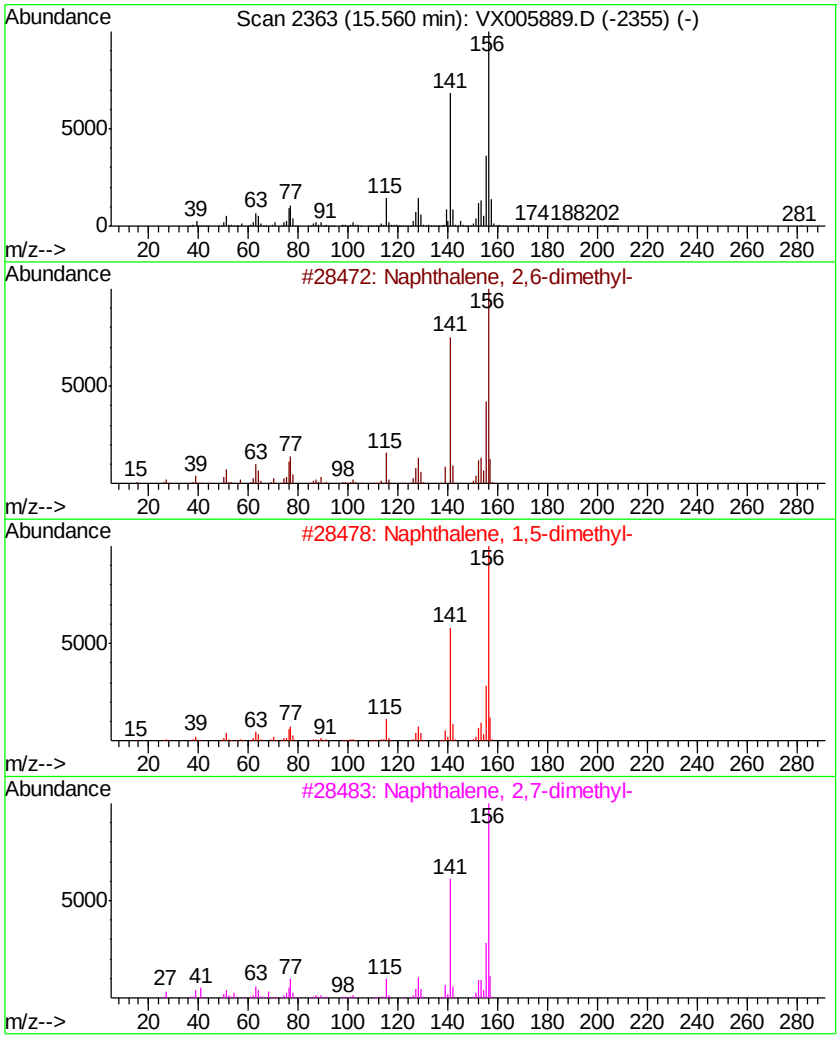
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

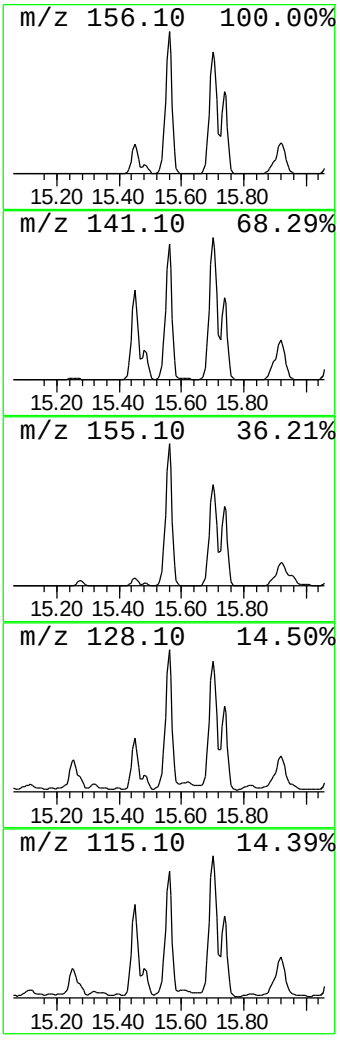
****

Peak Number 7 Naphthalene, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.56	685.07 ug/l	14936000	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX110918\
Data File : VX005889.D
Acq On : 10 Nov 2018 04:24
Operator : JC/MD
Sample : J5889-11
Misc : 12.72g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BOTTOM-B

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

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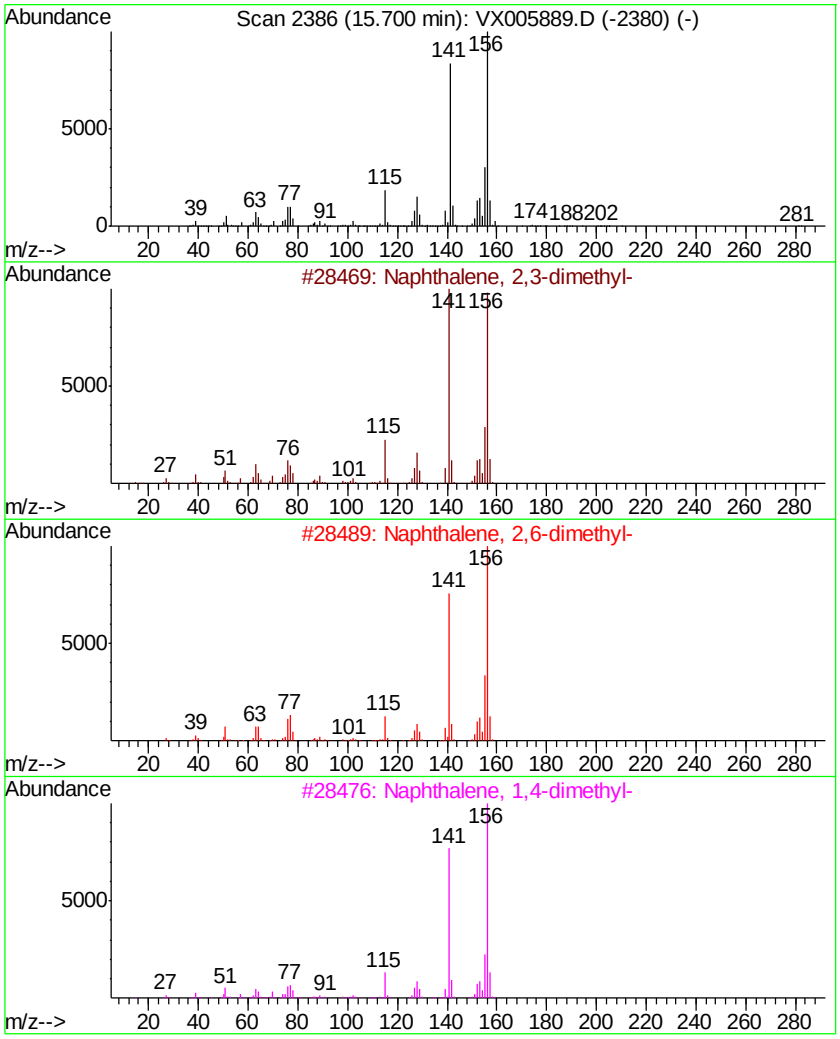
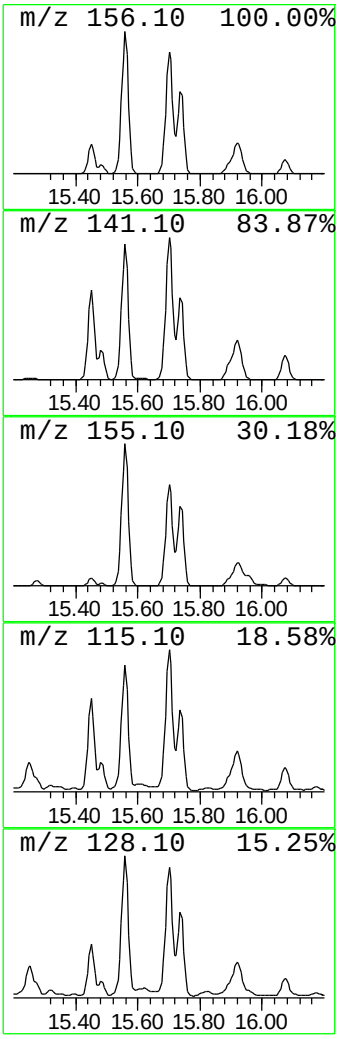
Peak Number 8 Naphthalene, 2,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.70	719.68 ug/l	15690500	1,4-Dichlorobenzene-d4	12.08

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
5	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX110918\
Data File : VX005889.D
Acq On : 10 Nov 2018 04:24
Operator : JC/MD
Sample : J5889-11
Misc : 12.72g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : BOTTOM-B

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X110718W.M
Quant Title : SW846 8260

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

					Internal Standard		
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
2,4-Dimethylstyrene	13.35	281.6	ug/l	6140280	4	12.08	1090110	50.0
Naphthalene, 1,2,...	13.48	155.2	ug/l	3384570	4	12.08	1090110	50.0
Naphthalene, 1,2,...	14.29	251.3	ug/l	5477730	4	12.08	1090110	50.0
Naphthalene, 1-me...	14.71	1265.2	ug/l	27584300	4	12.08	1090110	50.0
Naphthalene, 1-et...	15.45	239.8	ug/l	5228160	4	12.08	1090110	50.0
Naphthalene, 2,6-...	15.56	685.1	ug/l	14936000	4	12.08	1090110	50.0
Naphthalene, 2,3-...	15.70	719.7	ug/l	15690500	4	12.08	1090110	50.0