LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091319\
Data File : BM022680.D
Acq On : 13 Sep 2019 17:11
Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.034	4	8	21	rVB	533286	737638	1.71%	0.402%
2	4.040	174	179	186	rVB	70673	104027	0.24%	0.057%
3	5.358	398	403	412	rVB	214656	319242	0.74%	0.174%
4	5.493	419	426	439	rVB	514176	1048371	2.43%	0.571%
5	5.858	479	488	495	rVB	406074	695889	1.61%	0.379%
6	6.340	564	570	577	rVB	70816	102571	0.24%	0.056%
7	6.934	665	671	677	rBV	682809	1020718	2.36%	0.556%
8	6.993	677	681	688	rVV	435982	665529	1.54%	0.363%
9	7.069	688	694	703	rVB	901484	1336928	3.10%	0.729%
10	7.169	705	711	717	rBV	156894	253090	0.59%	0.138%
11	7.234	717	722	728	rBV	185995	274133	0.63%	0.149%
12	7.375	740	746	752	rVB	158847	249928	0.58%	0.136%
13	7.493	759	766	772	rBV	1956132	2990416	6.92%	1.630%
14	7.563	772	778	788	rVB	589019	927063	2.15%	0.505%
15	7.840	818	825	830	rBV	973139	1498341	3.47%	0.817%
16	7.887	830	833	839	rVV	72883	100669	0.23%	0.055%
17	7.957	839	845	853	rVV	741167	1229046	2.85%	0.670%
18	8.022	853	856	862	rVB	62979	95685	0.22%	0.052%
19	8.222	881	890	897	rBV	9290930	15002319	34.74%	8.176%
20	8.381	904	917	927	rBV	5590012	9048897	20.95%	4.932%
21	8.487	928	935	939	rVV	349265	542207	1.26%	0.296%
22	8.540	939	944	950	rVV2	140141	264119	0.61%	0.144%
23	8.675	959	967	974	rVB4	45155	108676	0.25%	0.059%
24	8.799	981	988	992	rBV	96833	152622	0.35%	0.083%
25	8.852	992	997	1003	rVV	106219	162493	0.38%	0.089%
26	8.952	1007	1014	1018	rVV	525441	842892	1.95%	0.459%
27	8.993	1018	1021	1023	rVV	200756	299217	0.69%	0.163%
28	9.028	1023	1027	1035	rVB	753856	1169039	2.71%	0.637%
29	9.199	1048	1056	1065	rVB	1050349	1722914	3.99%	0.939%
30	9.316	1065	1076	1082	rVV	2010989	4349630	10.07%	2.371%
31	9.381	1082	1087	1093	rVV	603755	964541	2.23%	0.526%
32	9.469	1096	1102	1107	rVV	280122	441117	1.02%	0.240%
33	9.528	1107	1112	1119	rVB	327209	515623	1.19%	0.281%
34	9.716	1136	1144	1150	rVB	111741	204189	0.47%	0.111%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091319\
Data File : BM022680.D
Acq On : 13 Sep 2019 17:11
Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Title : SVOA CALIBRATION

35	9.887	1168	1173	1187	rVB	854743	1391357	3.22%	0.758%
36	10.040	1187	1199	1209	rBV3	1393371	2801941	6.49%	1.527%
37	10.140	1209	1216	1228	rVV	244205	627685	1.45%	0.342%
38	10.269	1228	1238	1245	rVB2	617218	1328171	3.08%	0.724%
39	10.640	1291	1301	1304	rVV	1235920	2470544	5.72%	1.346%
40	10.704	1304	1312	1318	rVV	19455748	43189319	100.00%	23.538%
41	10.804	1322	1329	1338	rVV	4355150	7277065	16.85%	3.966%
42	10.987	1355	1360	1366	rVV2	87337	160931	0.37%	0.088%
43	11.051	1366	1371	1377	rVV2	75236	139273	0.32%	0.076%
44	11.740	1481	1488	1499	rVV2	175048	337170	0.78%	0.184%
45	12.010	1528	1534	1545	rVB	683624	1121333	2.60%	0.611%
46	12.187	1559	1564	1574	rVV	423695	682650	1.58%	0.372%
47	12.298	1574	1583	1590	rVV	9350269	15262432	35.34%	8.318%
48	12.410	1596	1602	1608	rVV	265308	476322	1.10%	0.260%
49	12.516	1608	1620	1629	rVV	5736345	9183658	21.26%	5.005%
50	12.869	1675	1680	1685	rBV	65418	102966	0.24%	0.056%
51	12.928	1685	1690	1695	rVV	135772	211391	0.49%	0.115%
52	13.187	1729	1734	1743	rVV2	154203	291649	0.68%	0.159%
53	13.304	1744	1754	1760	rVV	1235625	1842075	4.27%	1.004%
54	13.434	1768	1776	1782	rBV2	94106	161183	0.37%	0.088%
55	13.504	1782	1788	1798	rVB	170863	320200	0.74%	0.175%
56	13.598	1798	1804	1807	rBV	184297	271869	0.63%	0.148%
57	13.639	1807	1811	1820	rVB	241437	545701	1.26%	0.297%
58	13.792	1829	1837	1842	rBV	303477	527567	1.22%	0.288%
59	13.851	1842	1847	1848	rVV	170833	239294	0.55%	0.130%
60	13.881	1848	1852	1857	rVV	492454	697624	1.62%	0.380%
61	14.028	1872	1877	1881	rBV	92468	144454	0.33%	0.079%
62	14.163	1893	1900	1910	rBV2	572538	1170701	2.71%	0.638%
63	14.475	1943	1953	1958	rBV	2166009	3359104	7.78%	1.831%
64	14.534	1958	1963	1969	rVV	2708863	3998546	9.26%	2.179%
65	14.592	1969	1973	1978	rVV	243186	354246	0.82%	0.193%
66	14.651	1978	1983	1992	rVV5	107205	251288	0.58%	0.137%
67	14.869	2013	2020	2026	rVB	2068008	2858490	6.62%	1.558%
68	14.945	2026	2033	2038	rBV	112437	179716	0.42%	0.098%
69	15.181	2062	2073	2078	rBV3	75697	159378	0.37%	0.087%
70	15.386	2102	2108	2114	rVV	307084	488366	1.13%	0.266%
71	15.463	2114	2121	2125	rVV	740523	1120842	2.60%	0.611%

LSC Area Percent Report

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Data File : BM022680.D
Acq On : 13 Sep 2019 17:11
Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Title : SVOA CALIBRATION

72	15.516	2125	2130	2135	rVV	1653869	2278582	5.28%	1.242%
73	15.569	2135	2139	2143	rVV	154569	220844	0.51%	0.120%
74	15.675	2154	2157	2163	rVV3	63672	112426	0.26%	0.061%
75	15.816	2175	2181	2187	rVB3	126872	212309	0.49%	0.116%
76	15.963	2200	2206	2212	rVB4	148450	302937	0.70%	0.165%
77	16.186	2234	2244	2255	rBV2	714396	1240263	2.87%	0.676%
78	16.298	2257	2263	2266	rVV2	204245	367585	0.85%	0.200%
79	16.333	2266	2269	2274	rVV	174846	285146	0.66%	0.155%
80	16.380	2274	2277	2280	rVV2	63763	109552	0.25%	0.060%
81	16.416	2280	2283	2286	rVV2	75066	112665	0.26%	0.061%
82	16.486	2289	2295	2298	rVV4	57133	138250	0.32%	0.075%
83	16.586	2305	2312	2316	rBV2	85515	139998	0.32%	0.076%
84	16.833	2348	2354	2357	rBV2	217636	380071	0.88%	0.207%
85	16.857	2357	2358	2368	rVB	135940	153750	0.36%	0.084%
86	17.028	2383	2387	2392	rVV	138088	201293	0.47%	0.110%
87	17.080	2392	2396	2406	rVB	80734	134533	0.31%	0.073%
88	17.210	2412	2418	2421	rBV2	2763897	3838558	8.89%	2.092%
89	17.251	2421	2425	2430	rVB	1883441	2478147	5.74%	1.351%
90	17.304	2430	2434	2439	rBV2	998747	1358203	3.14%	0.740%
91	17.604	2476	2485	2491	rVV	564232	852325	1.97%	0.465%
92	18.122	2567	2573	2581	rVB4	110882	276773	0.64%	0.151%
93	18.280	2593	2600	2604	rBV	145909	229408	0.53%	0.125%
94	19.263	2762	2767	2773	rVB	548923	733242	1.70%	0.400%
95	19.592	2818	2823	2827	rBV	1056766	1585348	3.67%	0.864%
96	19.621	2827	2828	2833	rVB	332153	292421	0.68%	0.159%
97	21.110	3077	3081	3086	rBV5	65775	100259	0.23%	0.055%
98	21.386	3122	3128	3133	rBV	3442120	4644330	10.75%	2.531%
99	23.515	3485	3490	3495	rBV	633230	1096190	2.54%	0.597%
100	23.662	3508	3515	3528	rVB	2351590	4423084	10.24%	2.411%

Sum of corrected areas: 183484772

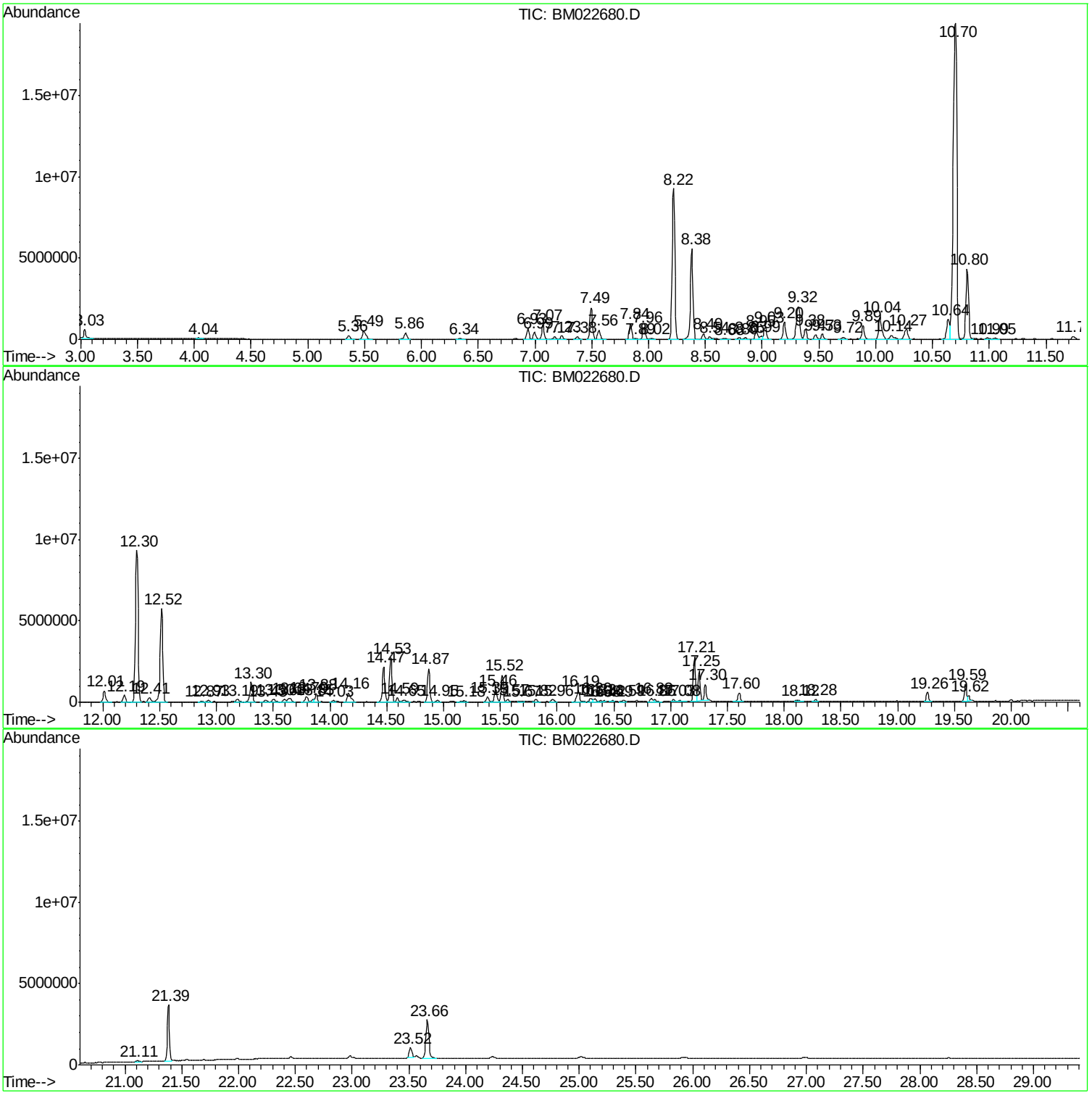
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091319\
Data File : BM022680.D
Acq On : 13 Sep 2019 17:11
Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091319\
Data File : BM022680.D
Acq On : 13 Sep 2019 17:11
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Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

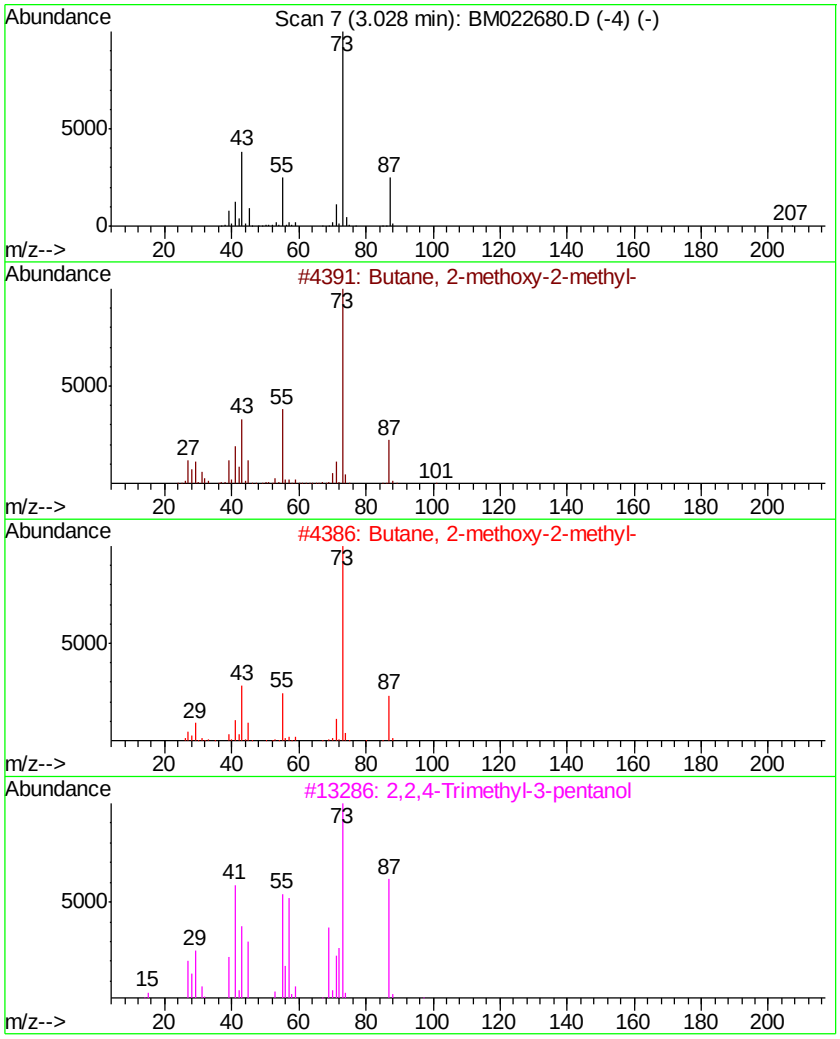
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

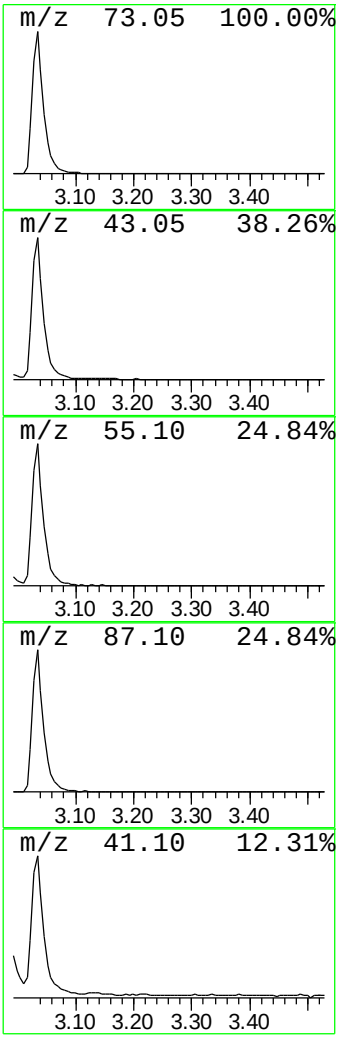
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.03	9.85 ng/ul	737638	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	56
4	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091319\
Data File : BM022680.D
Acq On : 13 Sep 2019 17:11
Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

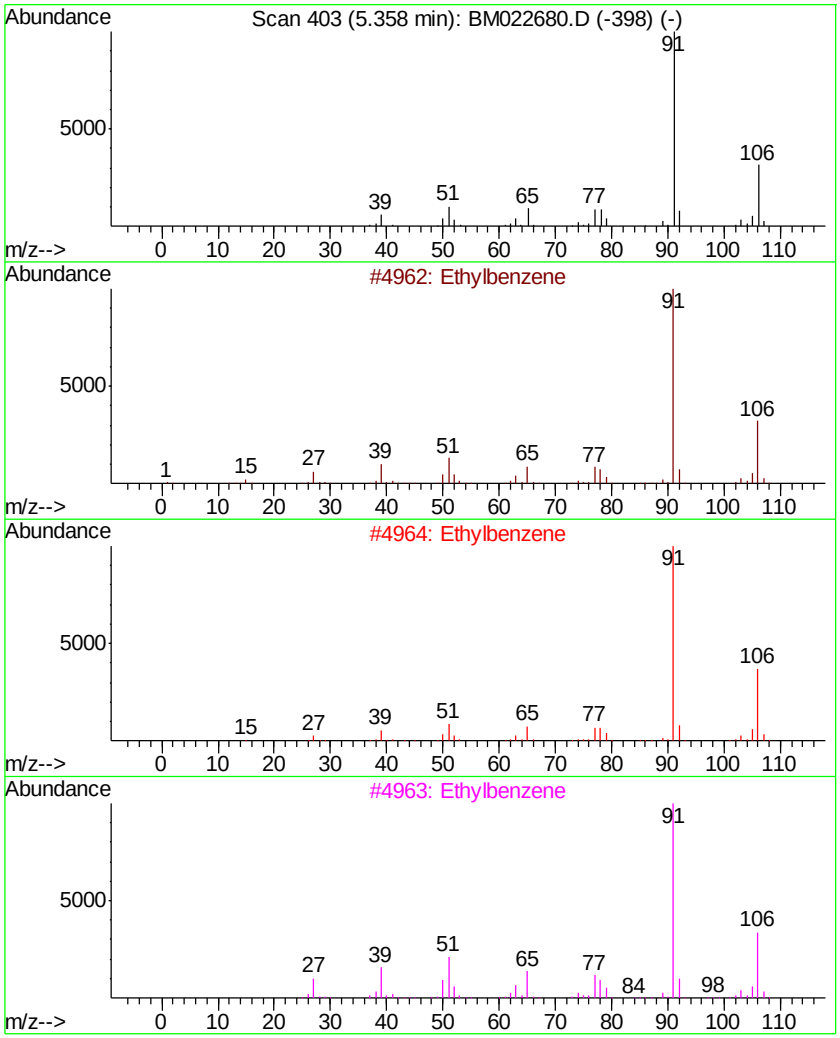
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

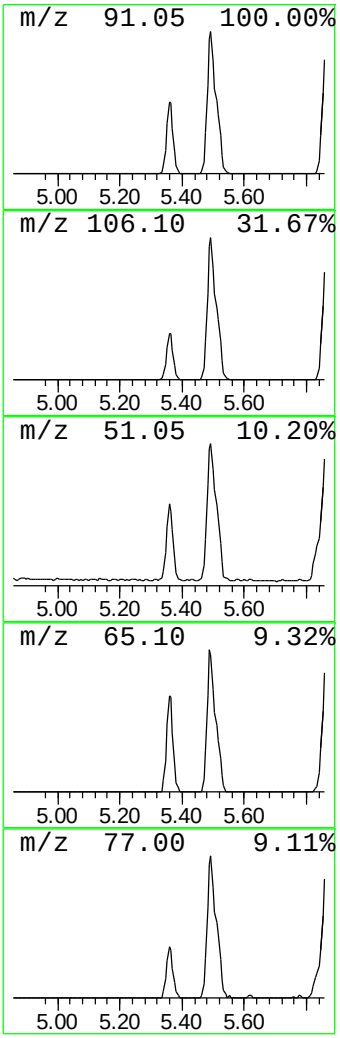
****

Peak Number 2 Ethylbenzene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.36	4.26 ng/ul	319242	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethylbenzene	106	C8H10	000100-41-4	94
2	Ethylbenzene	106	C8H10	000100-41-4	94
3	Ethylbenzene	106	C8H10	000100-41-4	91
4	Ethylbenzene	106	C8H10	000100-41-4	91
5	p-Xylene	106	C8H10	000106-42-3	72





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091319\
Data File : BM022680.D
Acq On : 13 Sep 2019 17:11
Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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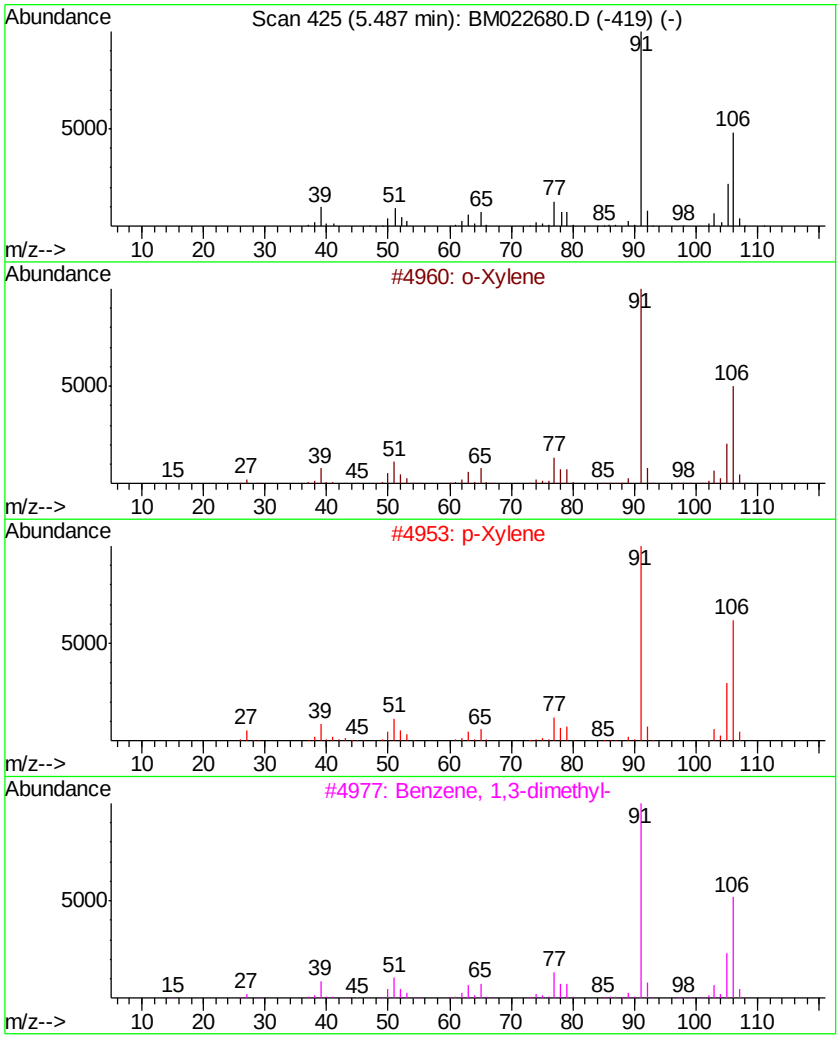
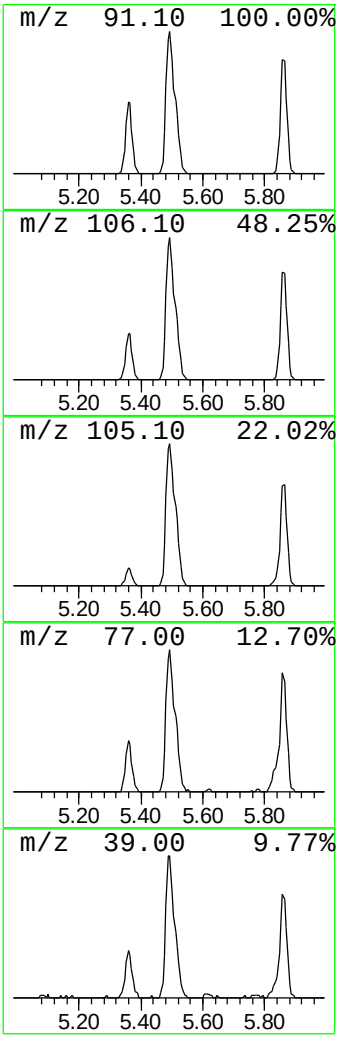
Peak Number 3 o-Xylene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.49	13.99 ng/ul	1048370	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Xylene	106	C8H10	000095-47-6	97
2	p-Xylene	106	C8H10	000106-42-3	97
3	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
4	p-Xylene	106	C8H10	000106-42-3	97
5	p-Xylene	106	C8H10	000106-42-3	95





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091319\
Data File : BM022680.D
Acq On : 13 Sep 2019 17:11
Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

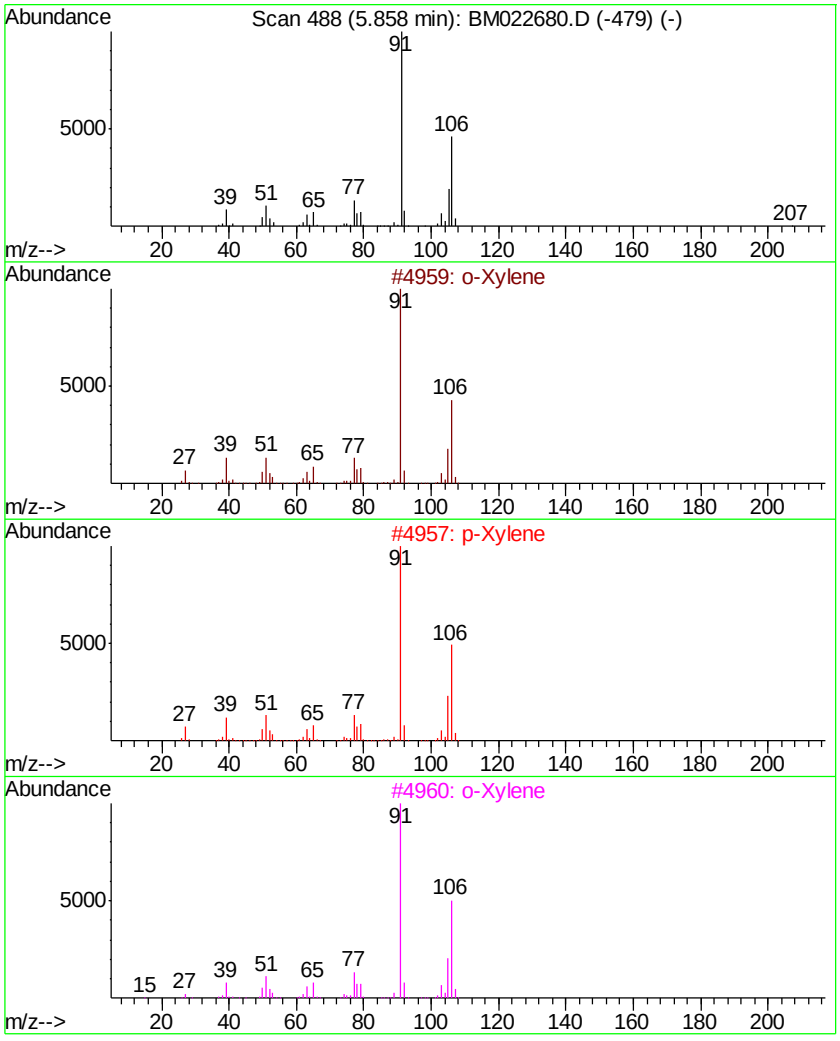
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

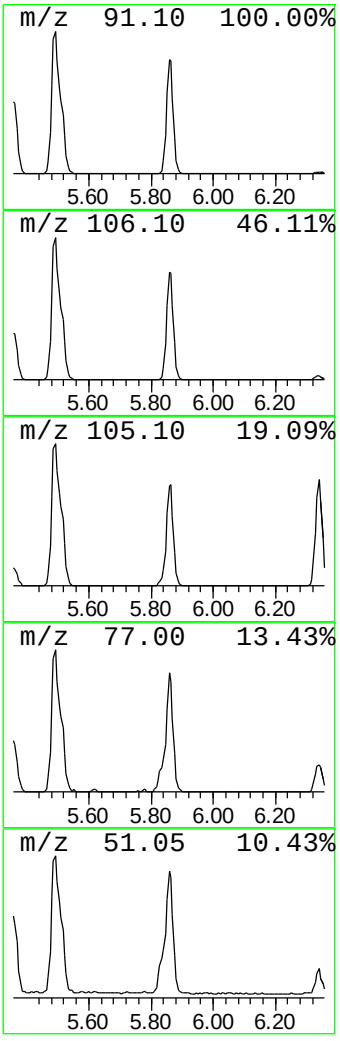
****

Peak Number 4 p-Xylene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.86	9.29 ng/ul	695889	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Xylene	106	C8H10	000095-47-6	97
2	p-Xylene	106	C8H10	000106-42-3	97
3	o-Xylene	106	C8H10	000095-47-6	97
4	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
5	o-Xylene	106	C8H10	000095-47-6	95





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091319\
Data File : BM022680.D
Acq On : 13 Sep 2019 17:11
Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

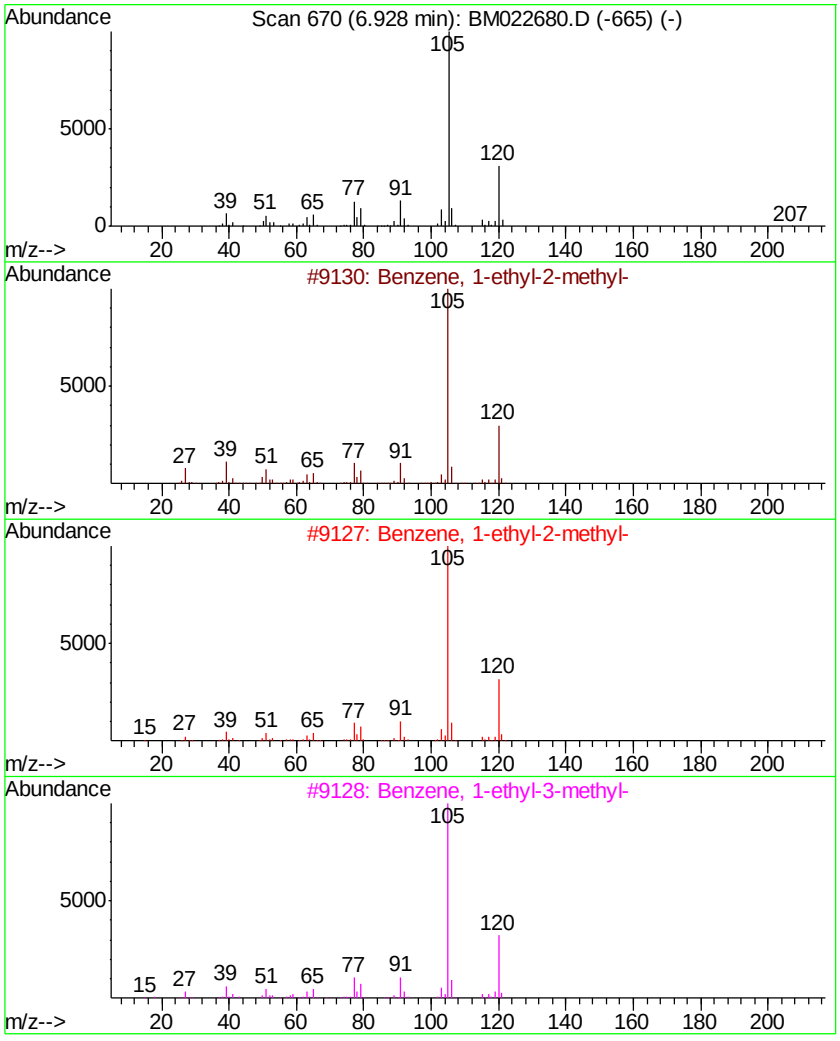
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

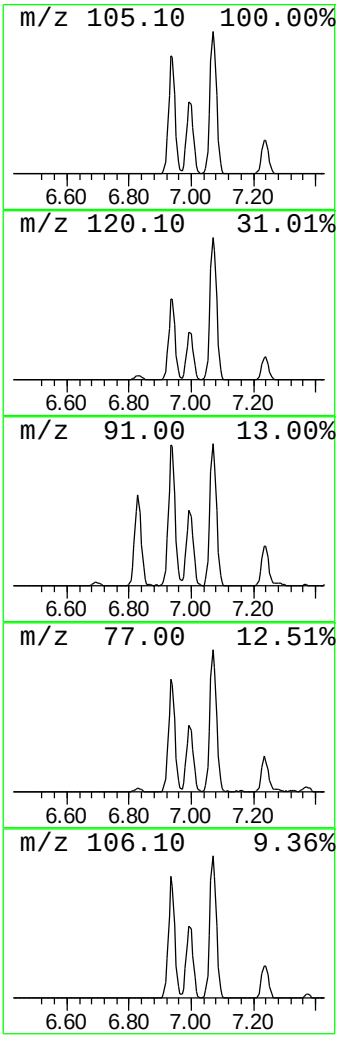
****

Peak Number 5 Benzene, 1-ethyl-3-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.93	13.62 ng/ul	1020720	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Library Search Compound Report

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Data File : BM022680.D
Acq On : 13 Sep 2019 17:11
Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

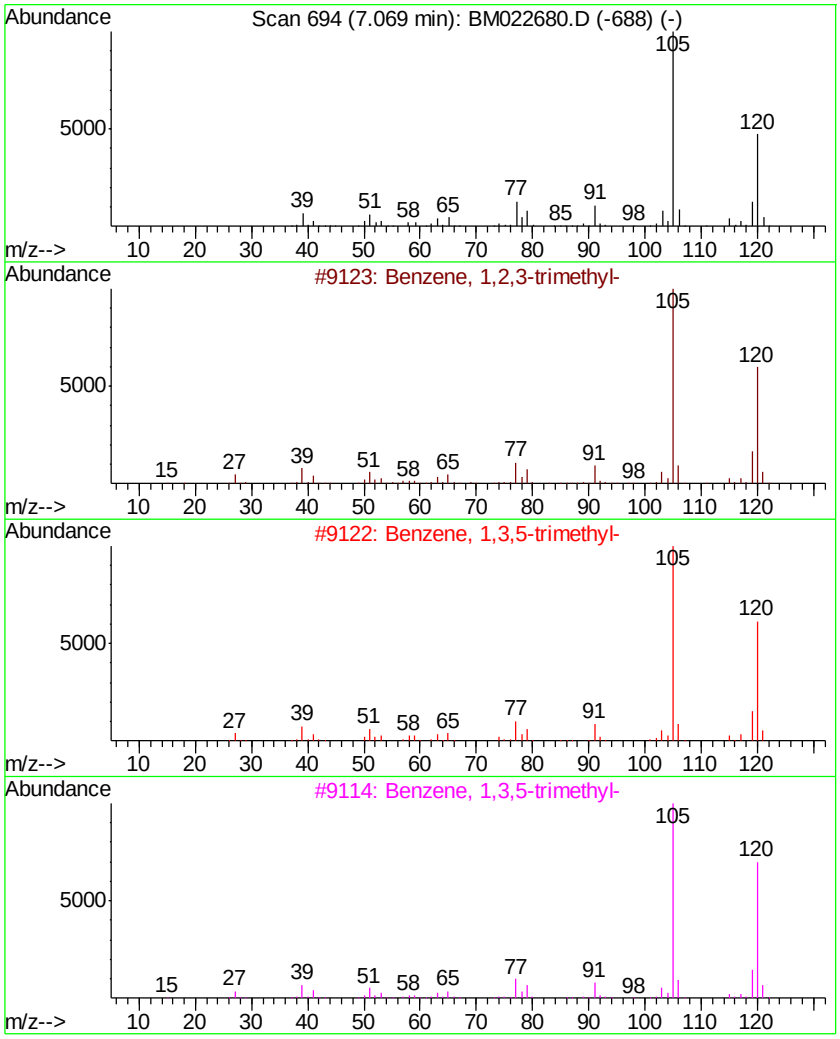
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

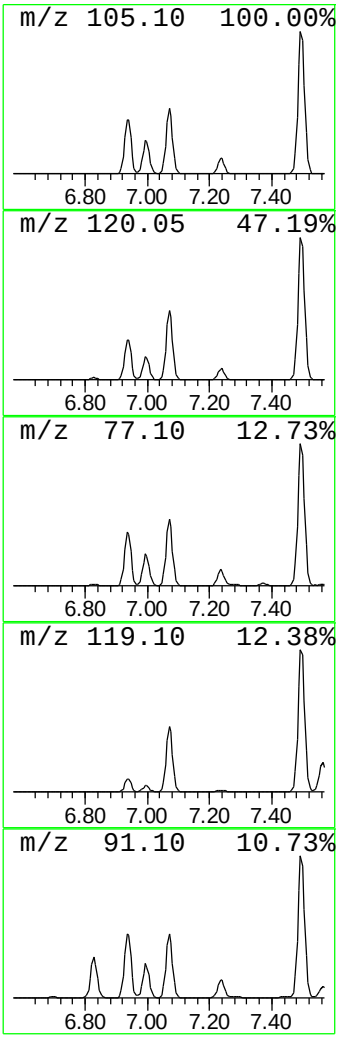
****

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.07	17.85 ng/ul	1336930	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	97
3	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	94
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94





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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

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Quant Title : SVOA CALIBRATION

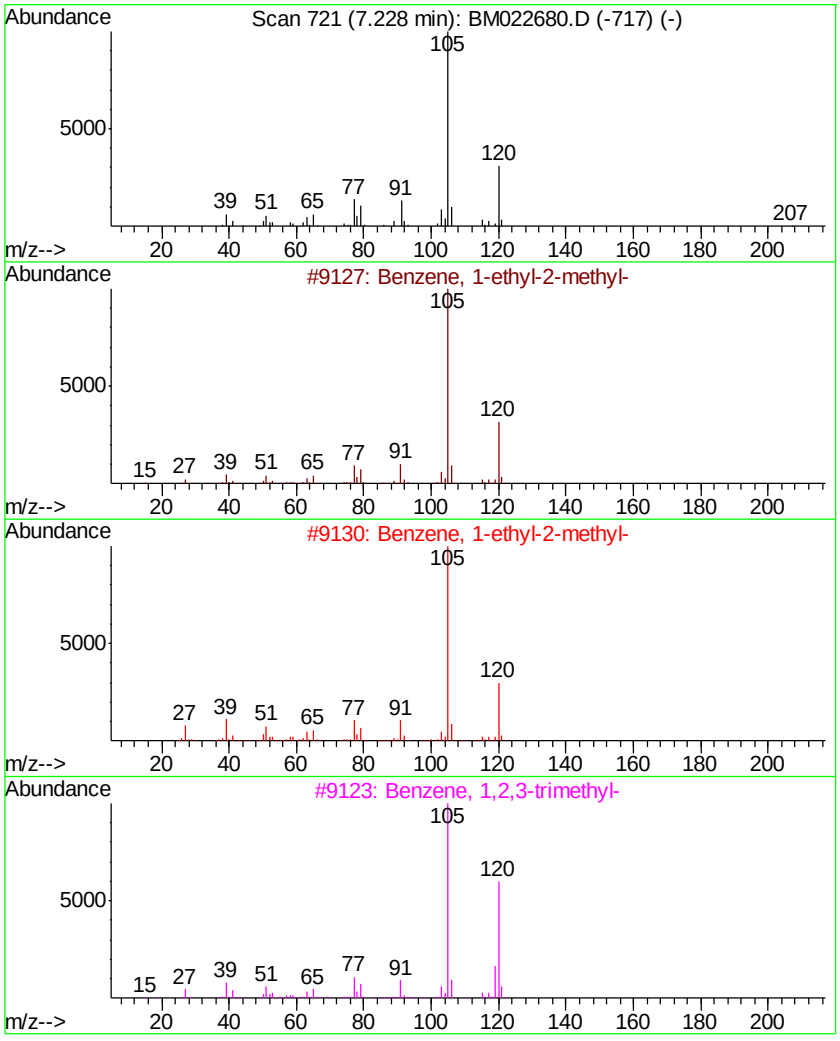
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

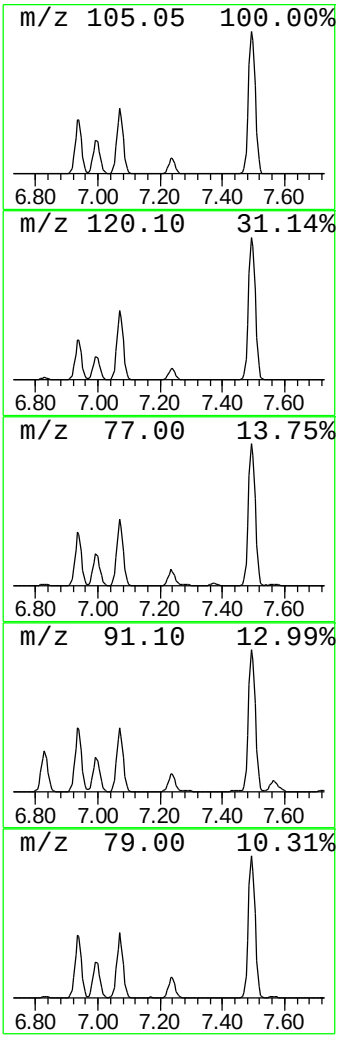
****

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.23	3.66 ng/ul	274133	1,4-Dichlorobenzene-d4	7.84

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-2-methyl-	120	C9H12	000611-14-3	94
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91





Library Search Compound Report

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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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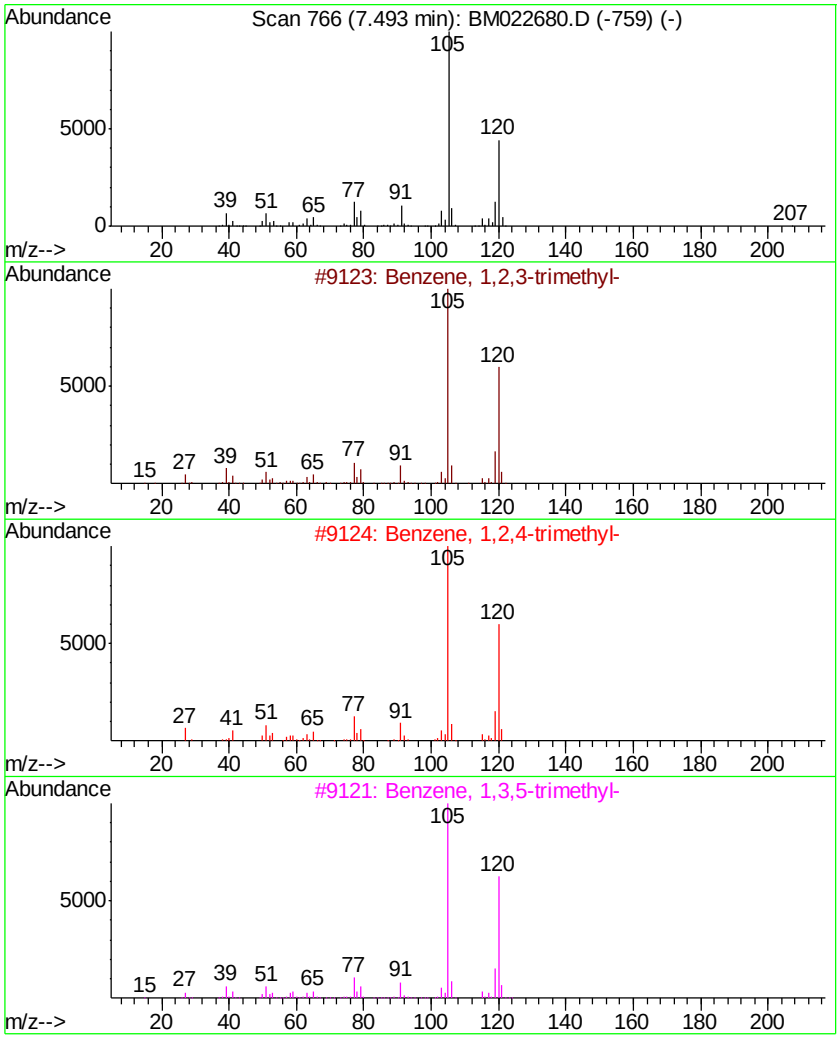
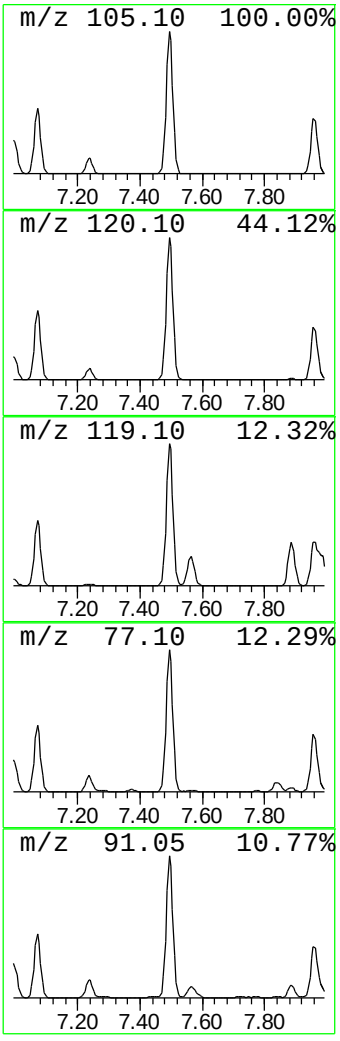
Peak Number 8 Benzene, 1,2,4-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.49	39.92 ng/ul	2990420	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	94
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





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Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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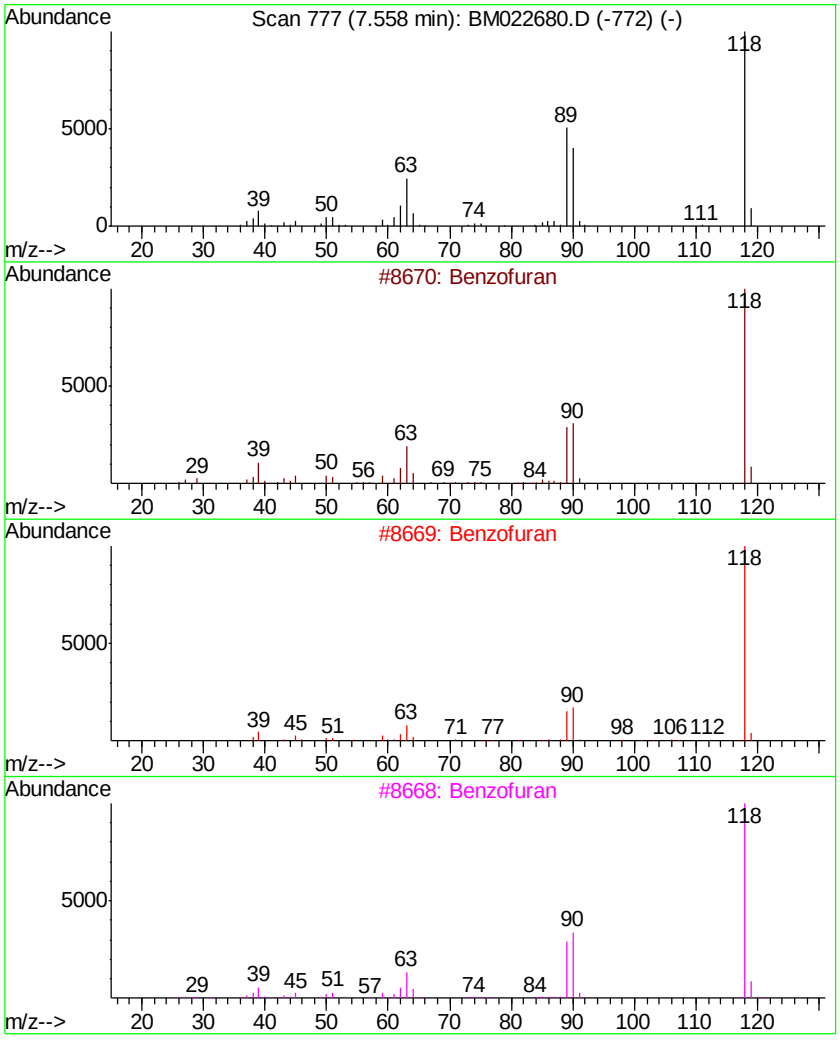
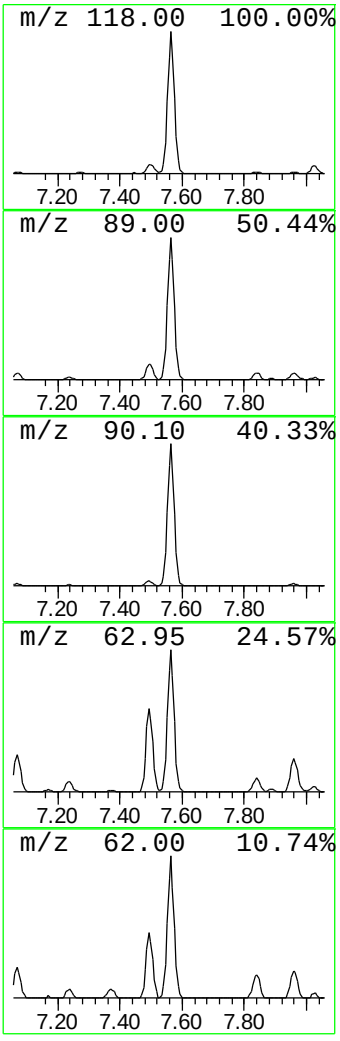
Peak Number 9 Benzofuran Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	12.37 ng/ul	927063	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran	118	C8H6O	000271-89-6	91
2	Benzofuran	118	C8H6O	000271-89-6	91
3	Benzofuran	118	C8H6O	000271-89-6	87
4	3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	83
5	3-Pyridineacetonitrile	118	C7H6N2	006443-85-2	40





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Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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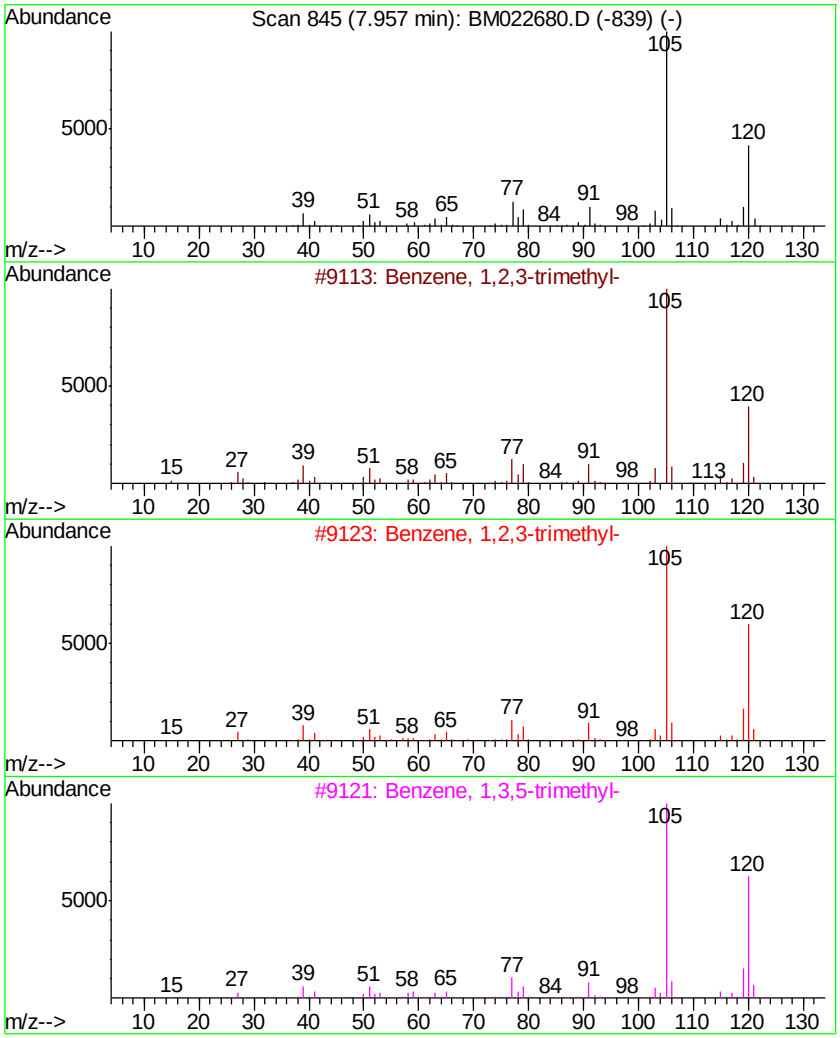
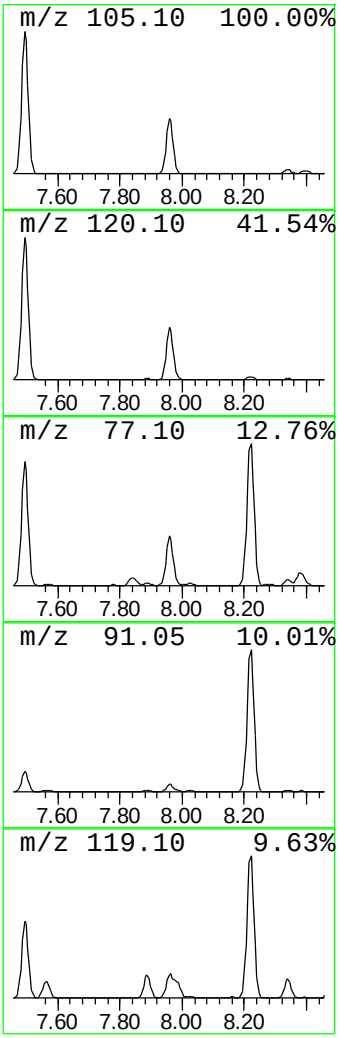
Peak Number 10 Benzene, 1,3,5-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.96	16.41 ng/ul	1229050	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	94
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





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Acq On : 13 Sep 2019 17:11
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Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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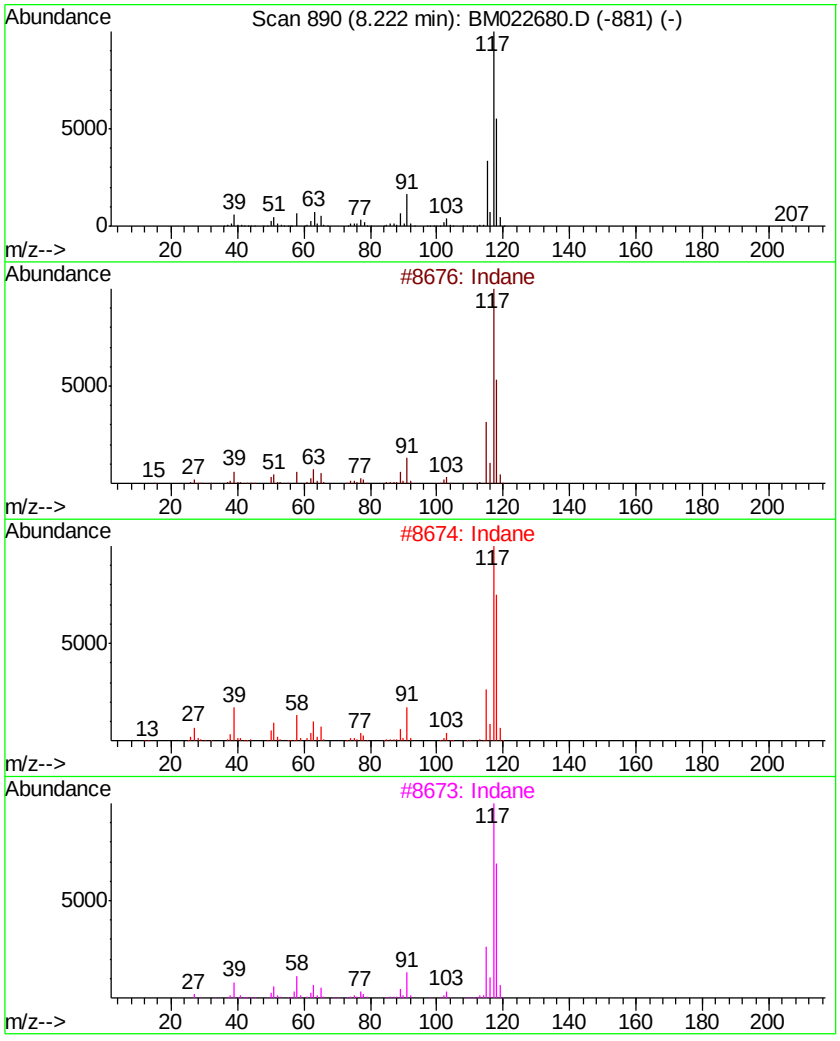
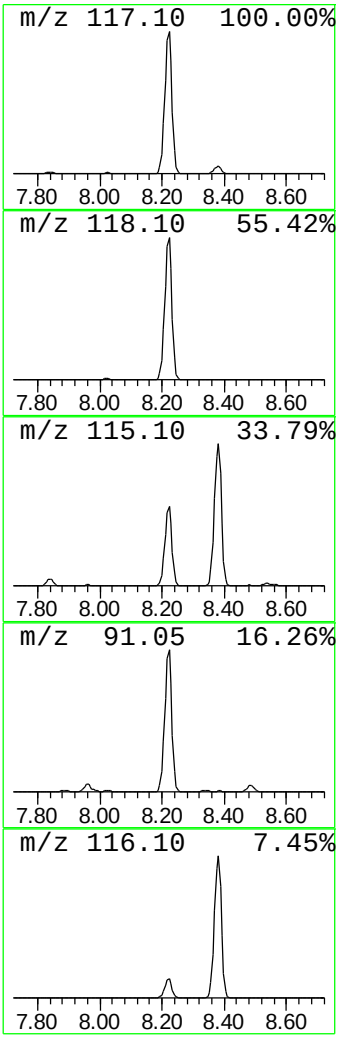
Peak Number 11 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.22	200.25 ng/ul	15002300	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	91
3	Indane	118	C9H10	000496-11-7	83
4	Deltacyclene	118	C9H10	007785-10-6	72
5	Indane	118	C9H10	000496-11-7	64





Library Search Compound Report

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Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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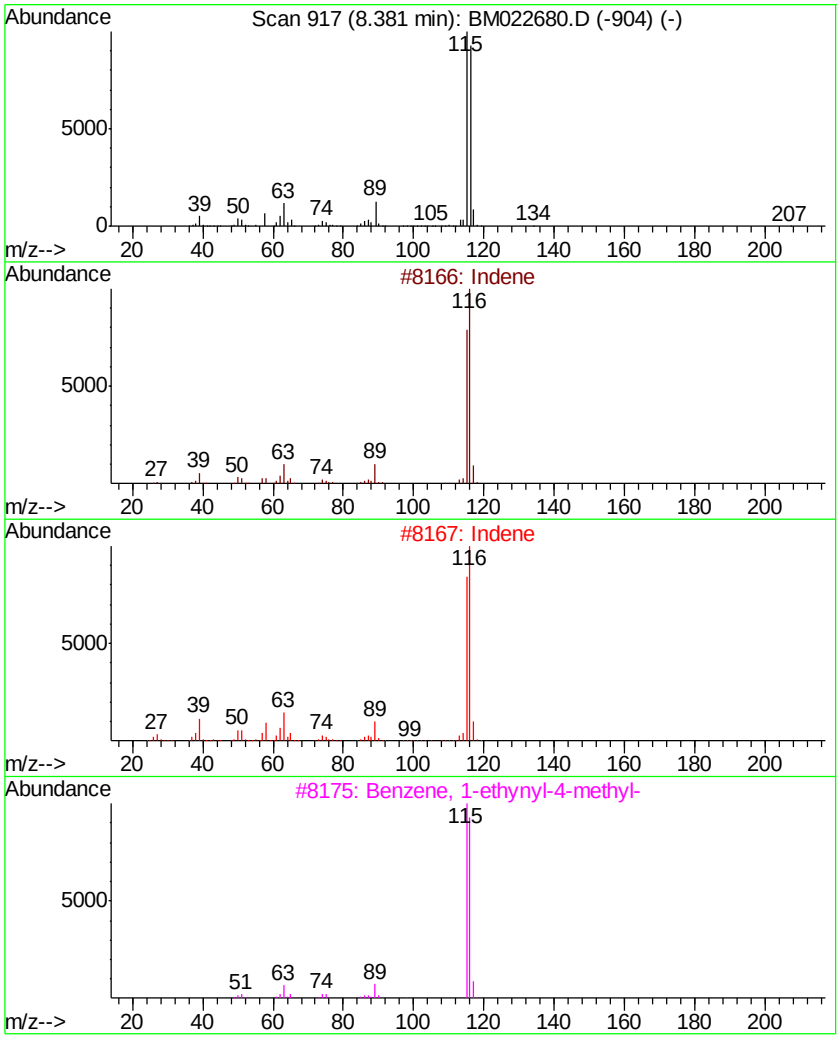
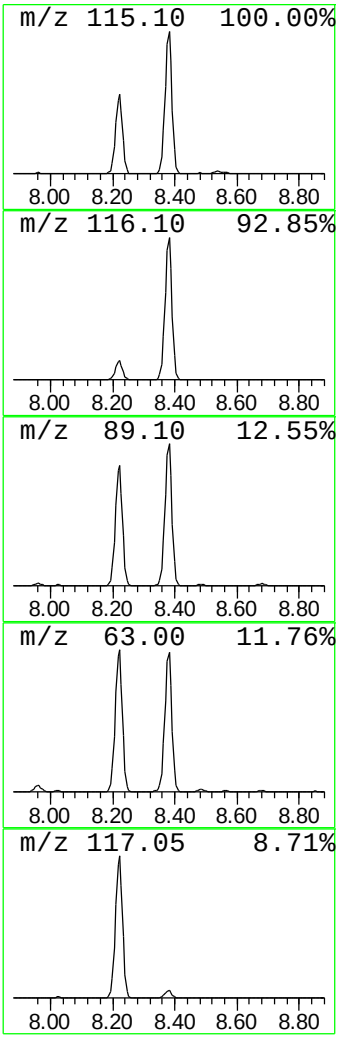
Peak Number 12 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.38	120.79 ng/ul	9048900	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	97
2	Indene	116	C9H8	000095-13-6	94
3	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91
4	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91
5	Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	90





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Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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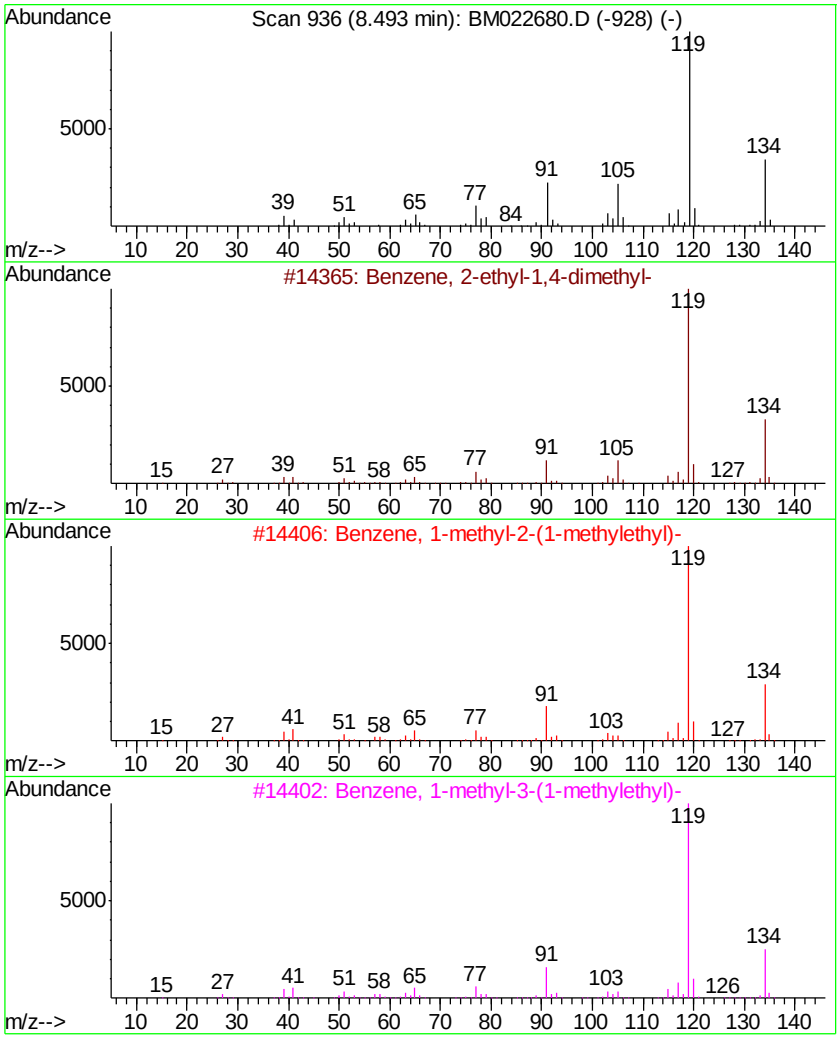
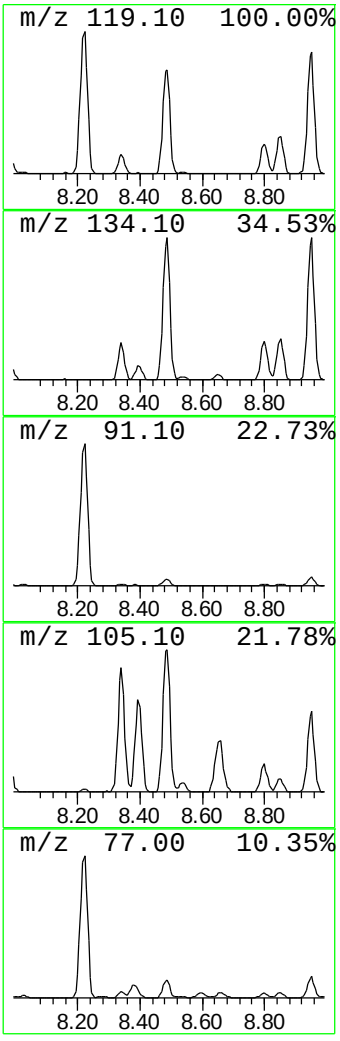
Peak Number 13 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.49	7.24 ng/ul	542207	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	95
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5	Benzene, 1-methyl-4-(1-methyleth...	134	C10H14	000099-87-6	94





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Misc :
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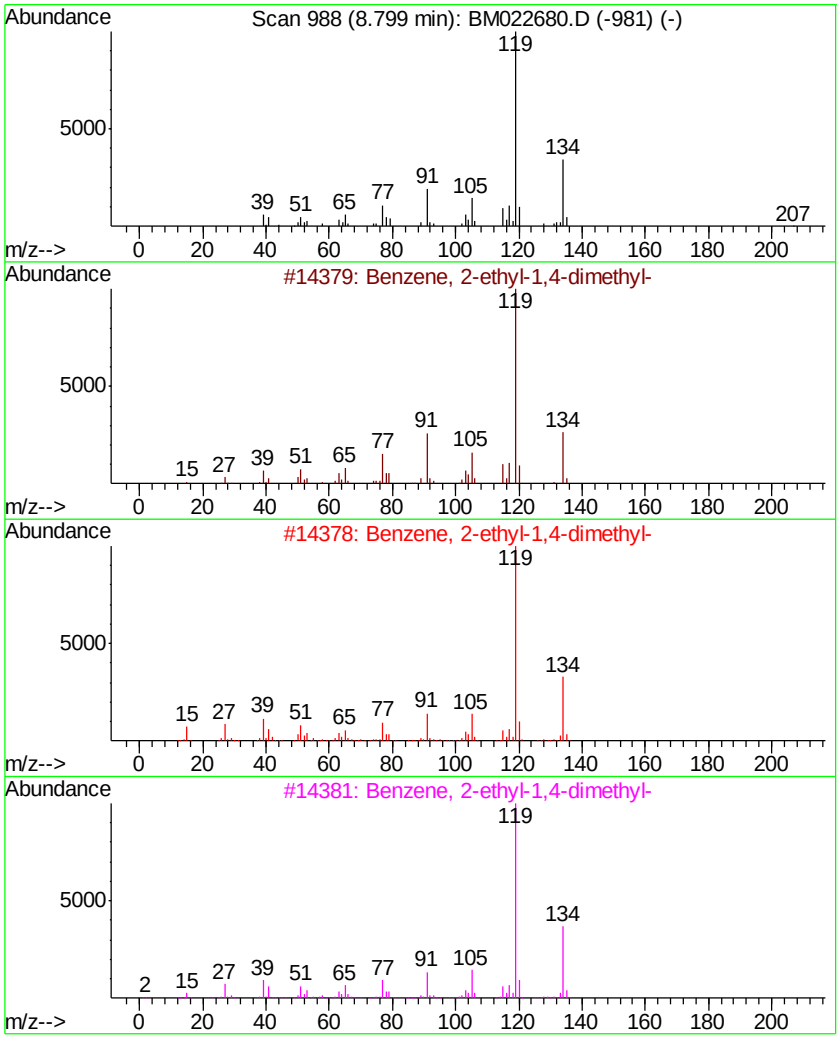
Quant Title : SVOA CALIBRATION

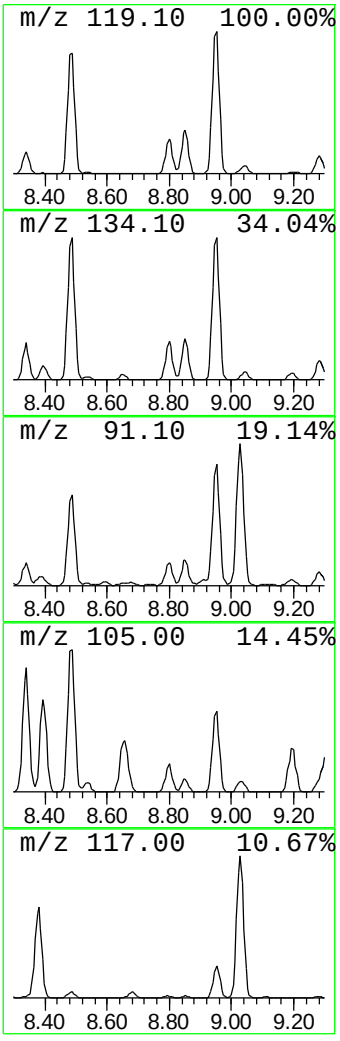
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Benzene, 1-methyl-4-(1-meth... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.80	2.04 ng/ul	152622	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4	Benzene, 1-methyl-4-(1-methyleth...	134	C10H14	000099-87-6	95
5	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	95





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Misc :
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Instrument : BNA_M
ClientSampleId : BF575DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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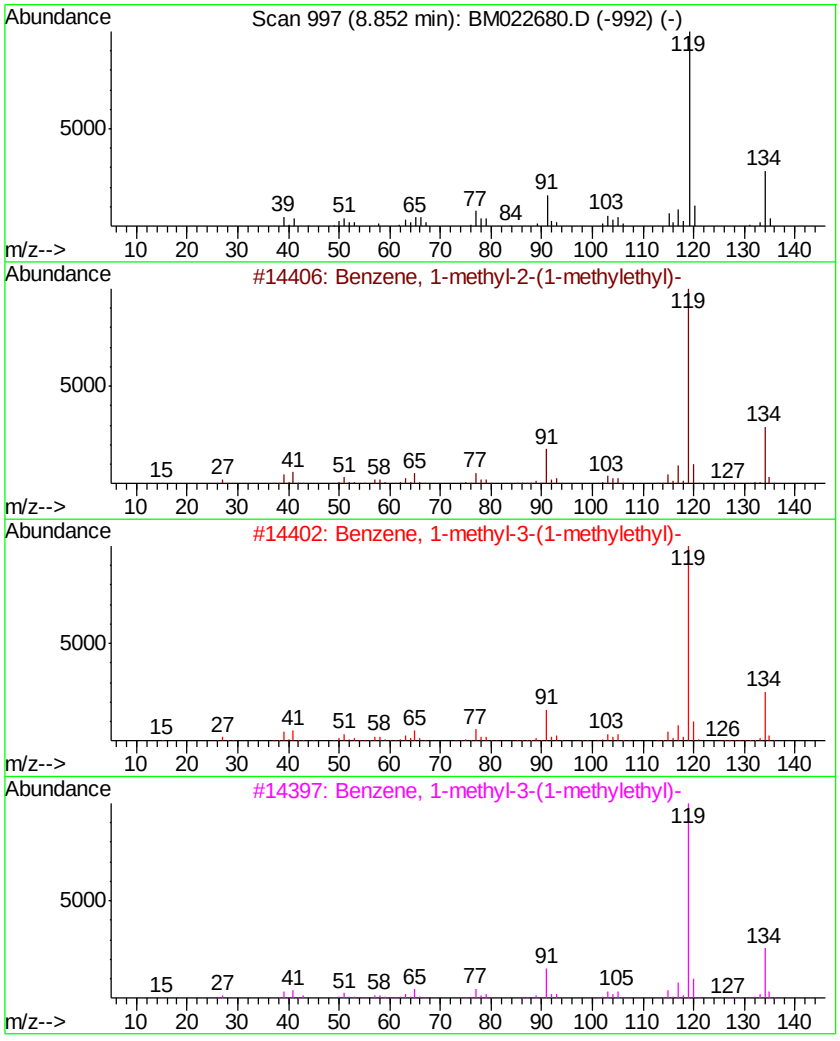
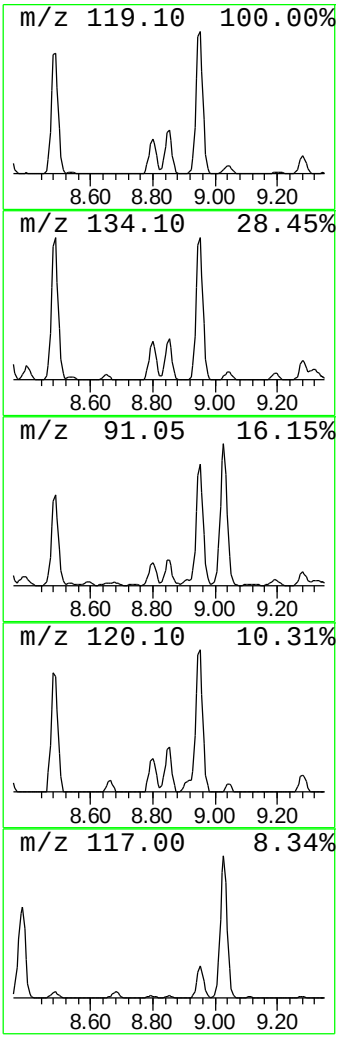
Peak Number 15 Benzene, 1-methyl-2-(1-meth... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.85	2.17 ng/ul	162493	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	97
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
4	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	97
5	Benzene, 1-methyl-4-(1-methyleth...	134	C10H14	000099-87-6	95





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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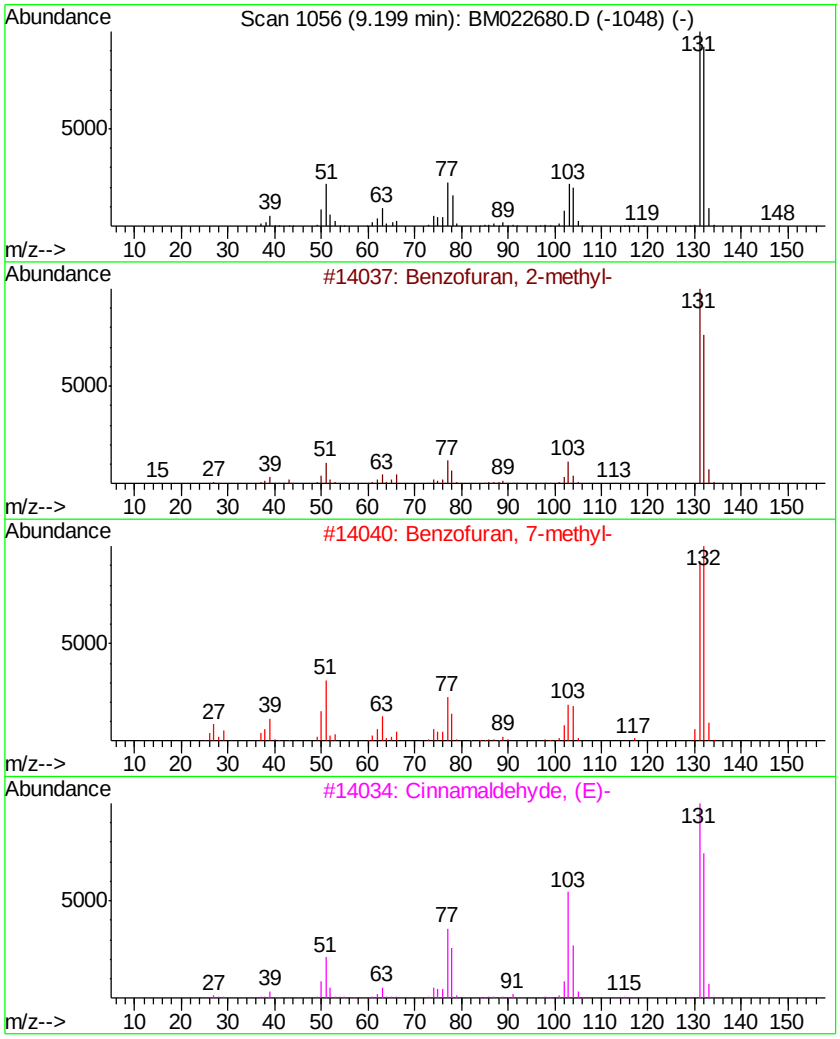
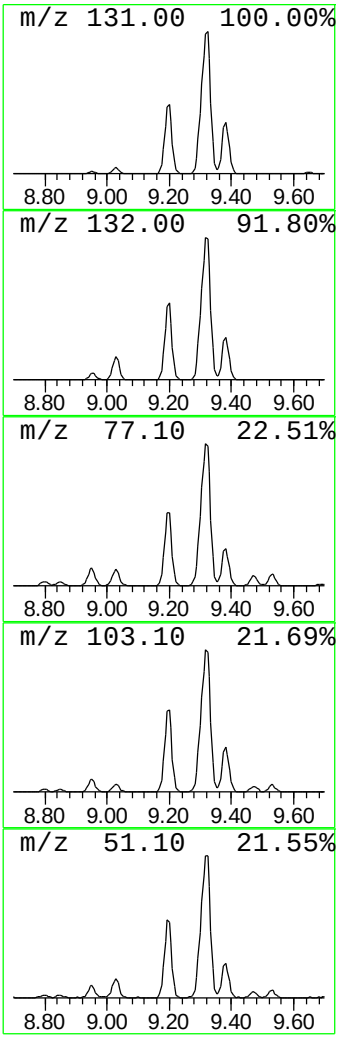
Peak Number 16 Benzofuran, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.20	23.00 ng/ul	1722910	1,4-Dichlorobenzene-d4	7.84

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	91
3	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	91
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90





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ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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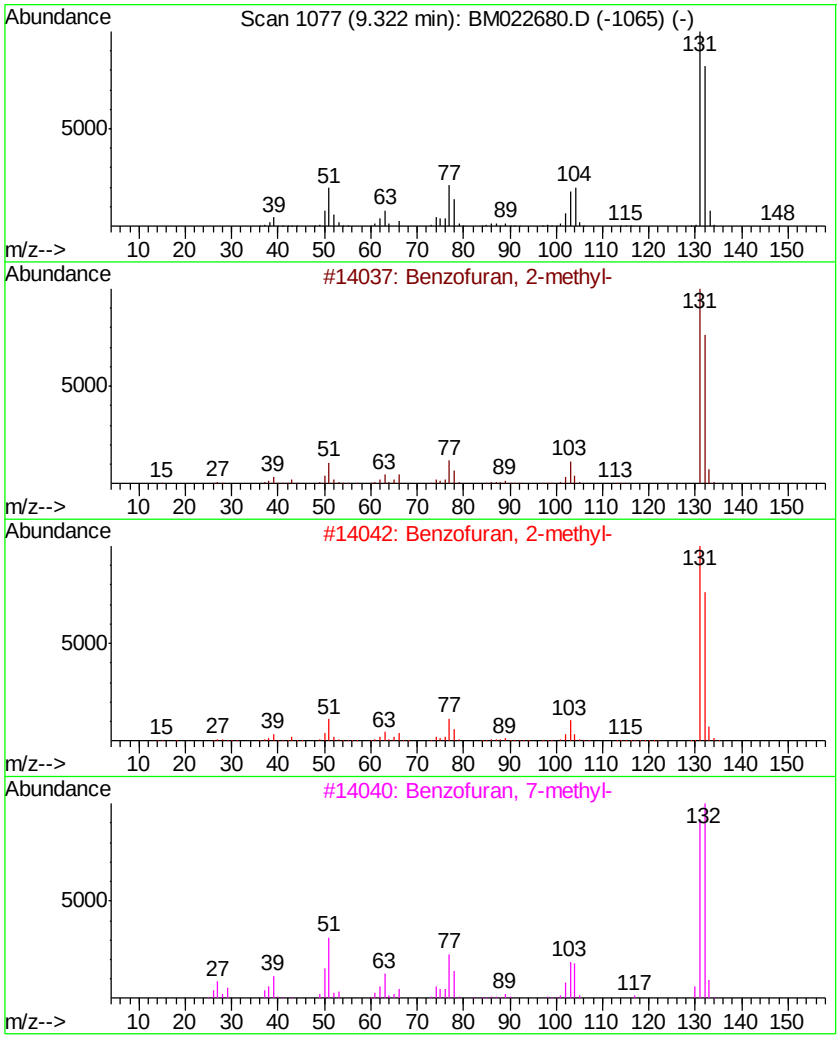
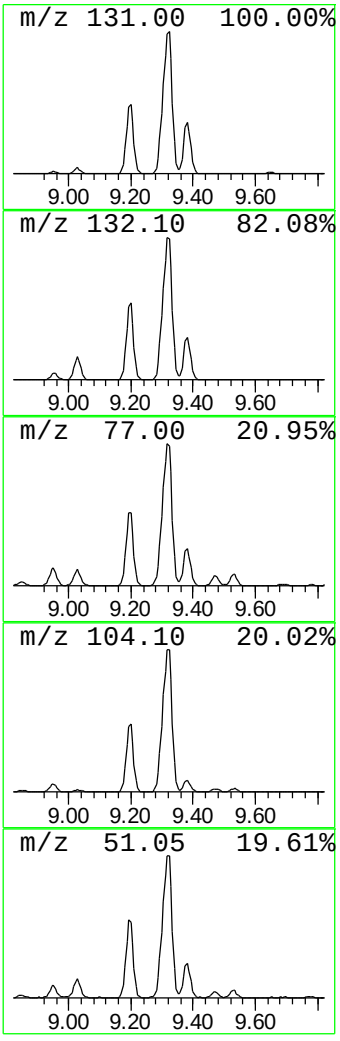
Peak Number 17 Benzofuran, 7-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.32	35.21 ng/ul	4349630	Naphthalene-d8	10.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
3	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	91
4	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	91
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91





Library Search Compound Report

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Acq On : 13 Sep 2019 17:11
Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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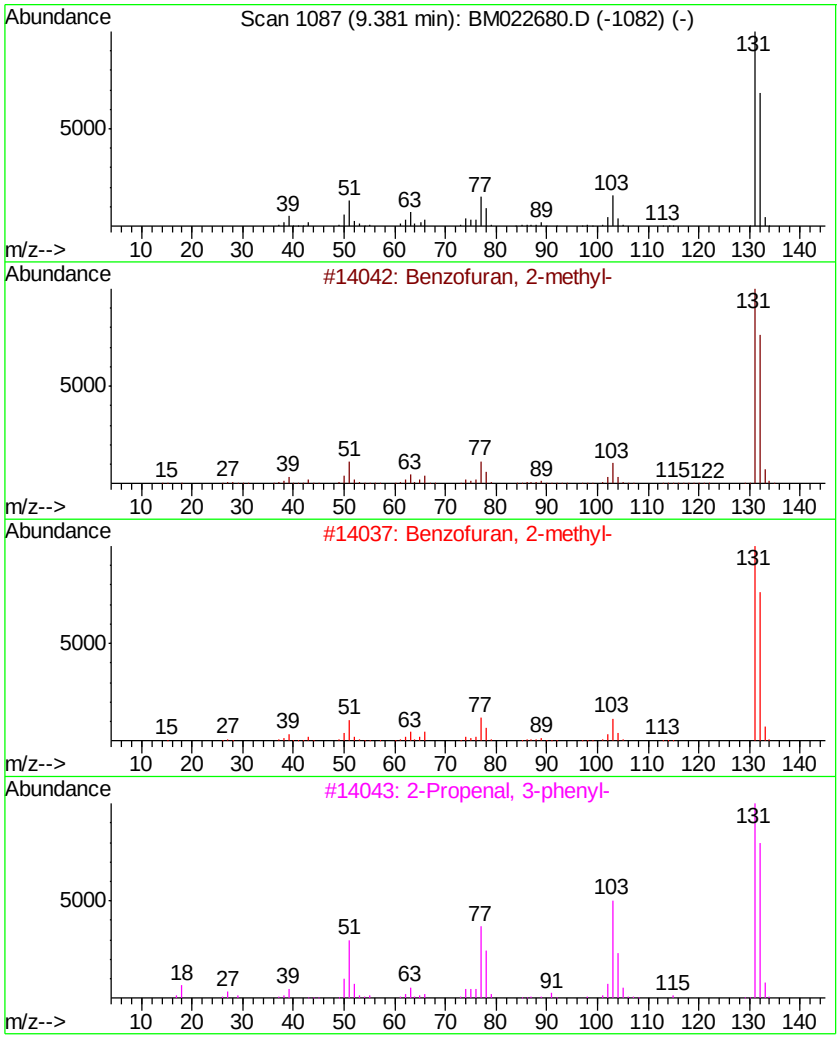
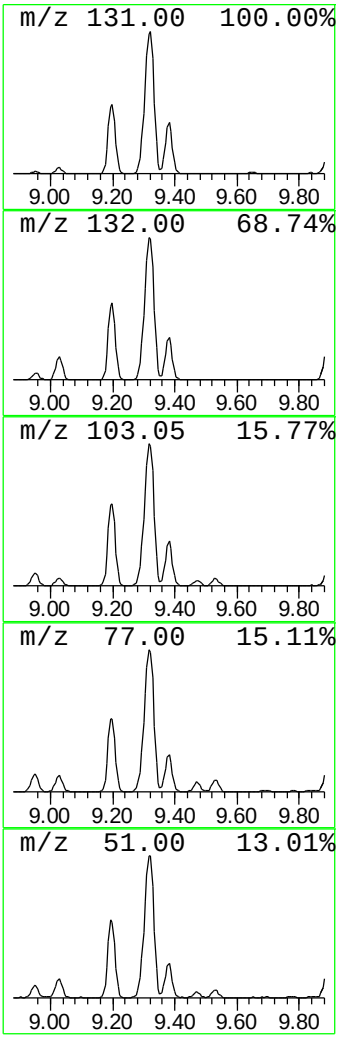
Peak Number 18 2-Propenal, 3-phenyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.38	7.81 ng/ul	964541	Naphthalene-d8	10.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	86
3	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	78
4	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	78
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	76





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Misc :
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Instrument : BNA_M
ClientSampleId : BF575DL

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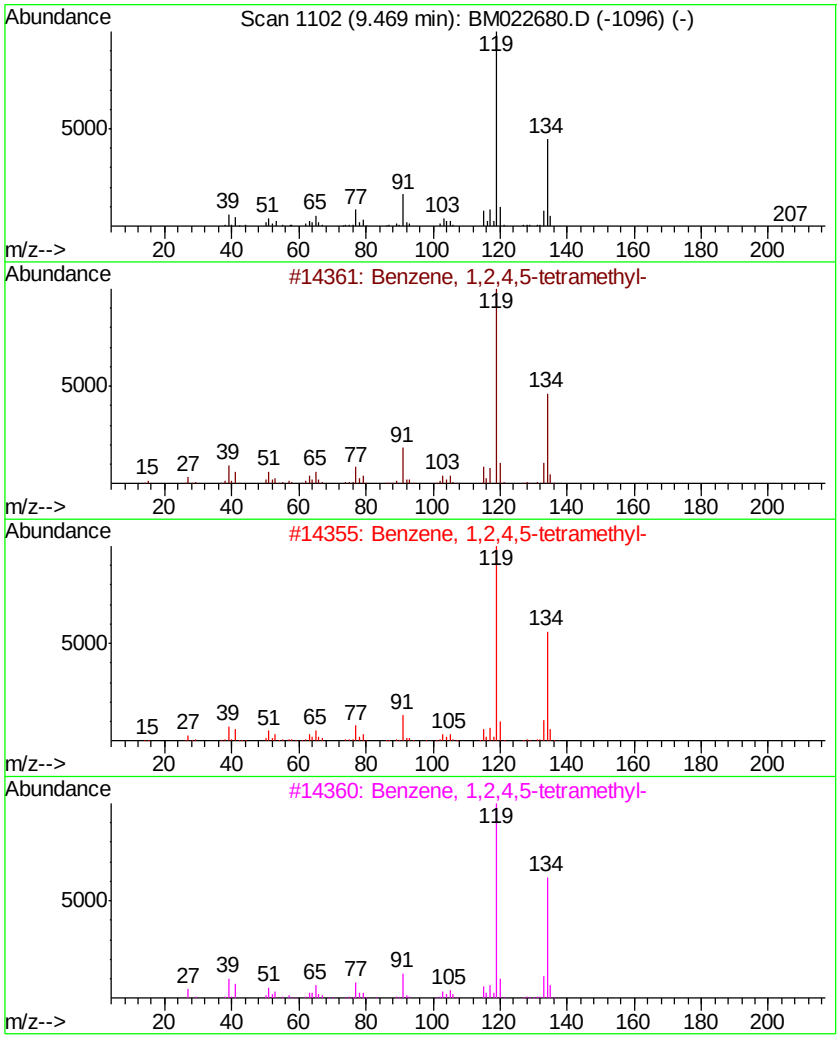
Quant Title : SVOA CALIBRATION

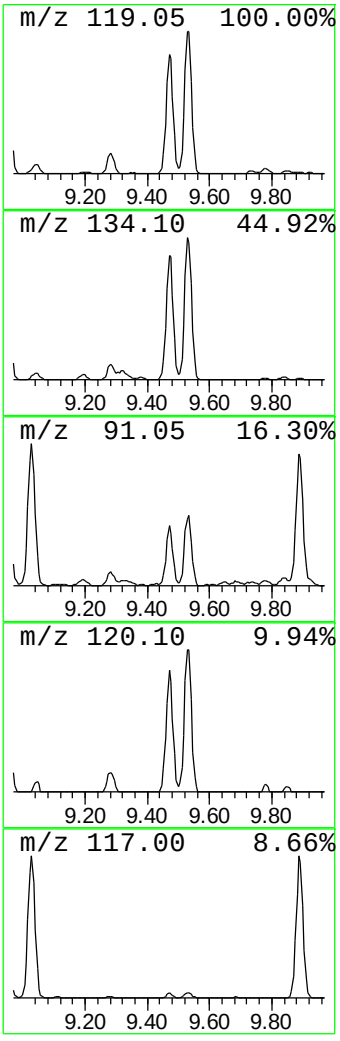
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.47	3.57 ng/ul	441117	Naphthalene-d8	10.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95





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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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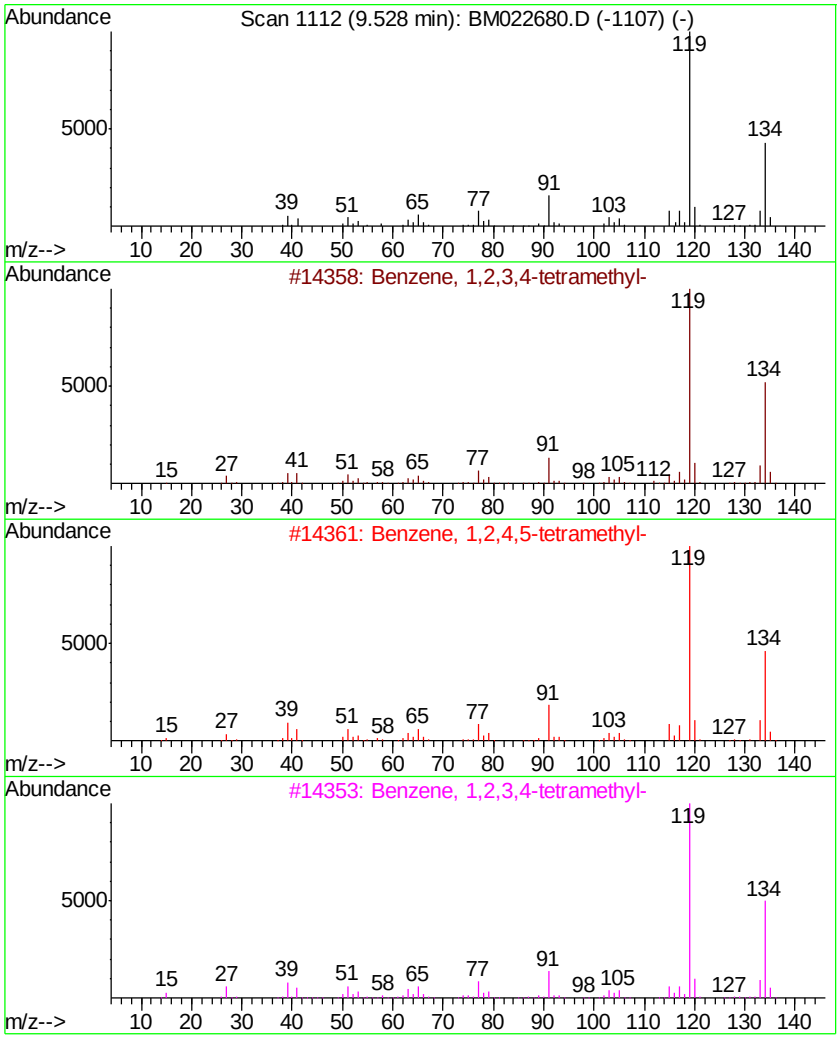
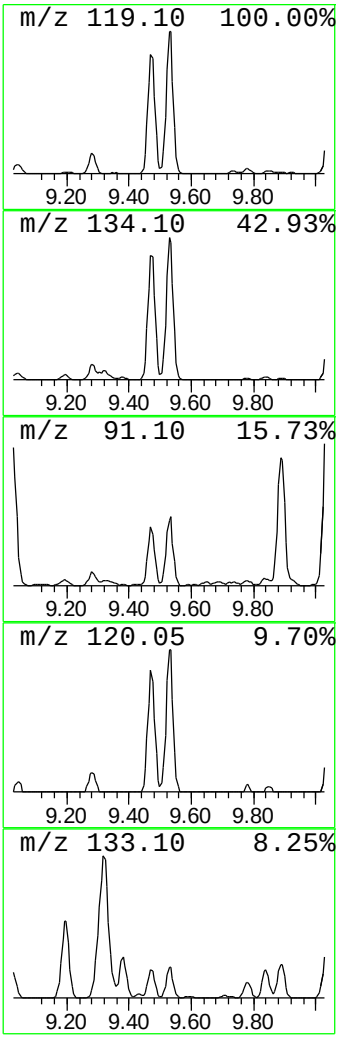
Peak Number 20 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.53	4.17 ng/ul	515623	Naphthalene-d8	10.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95





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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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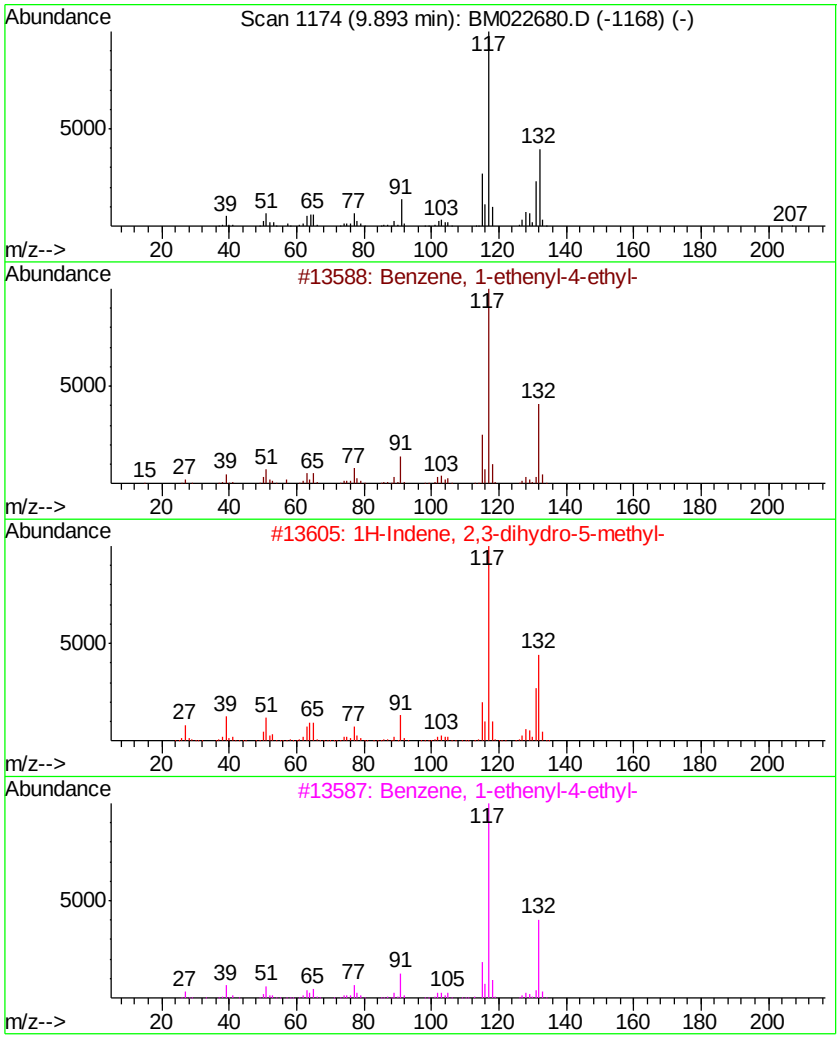
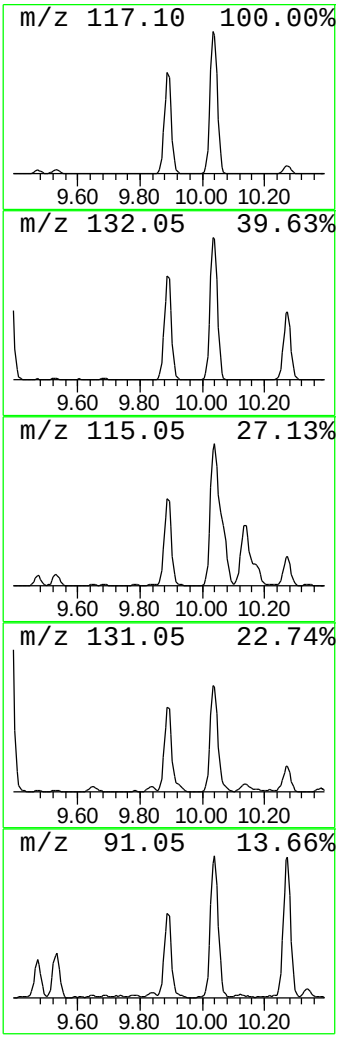
Peak Number 21 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.89	11.26 ng/ul	1391360	Naphthalene-d8	10.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90





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Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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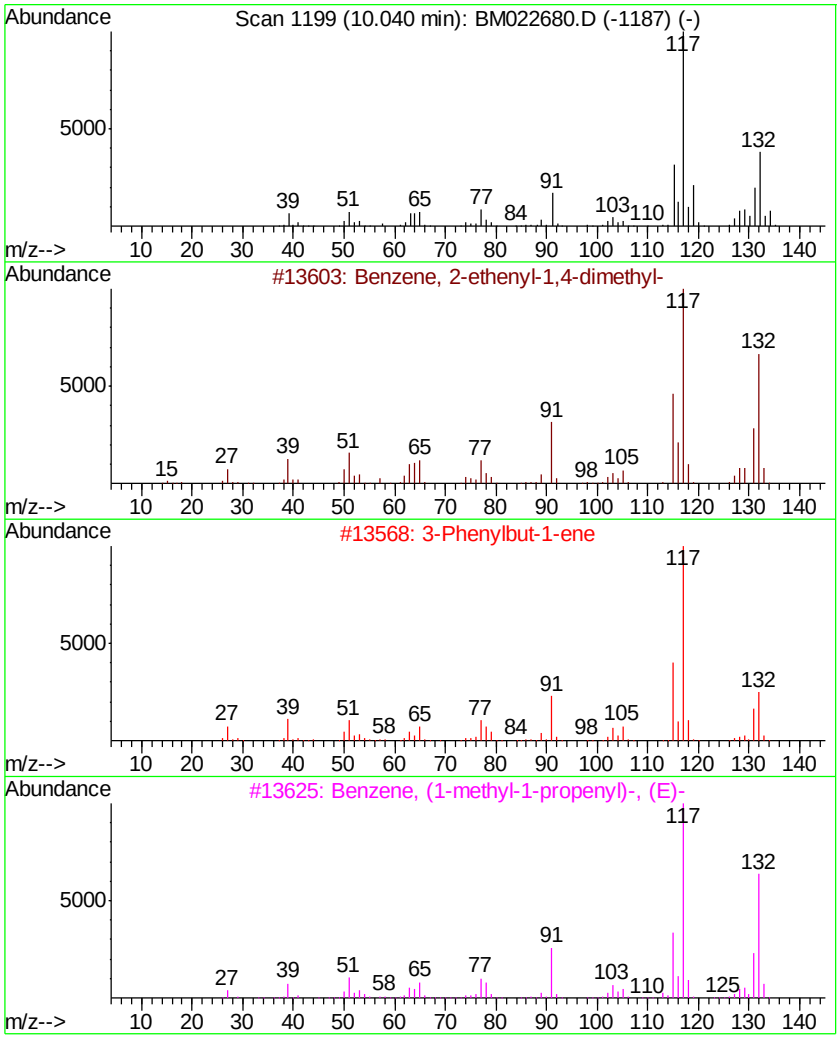
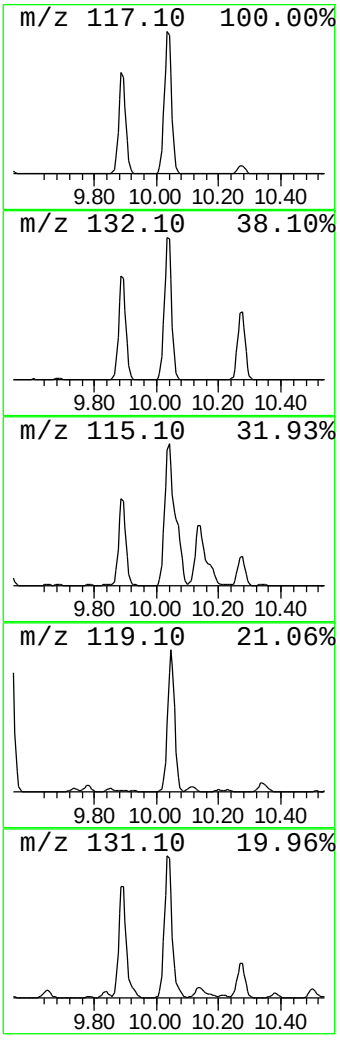
Peak Number 22 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.04	22.68 ng/ul	2801940	Naphthalene-d8	10.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83
4	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	81
5	Indan, 1-methyl-	132	C10H12	000767-58-8	81





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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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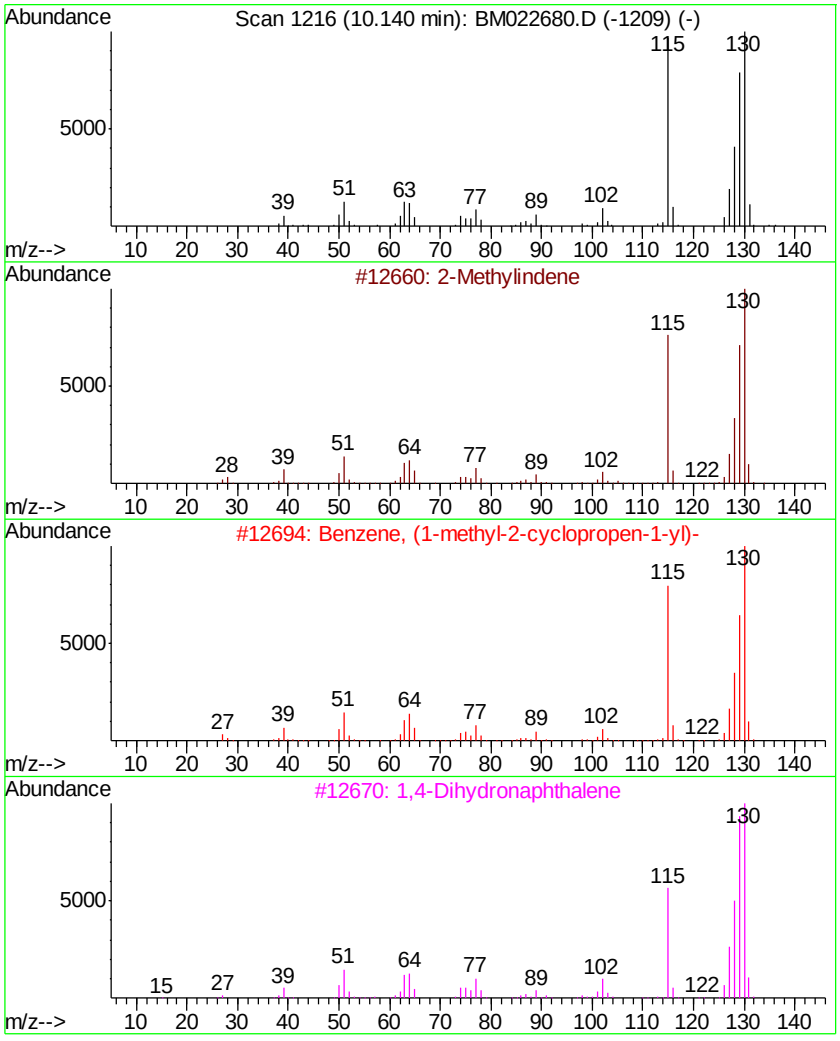
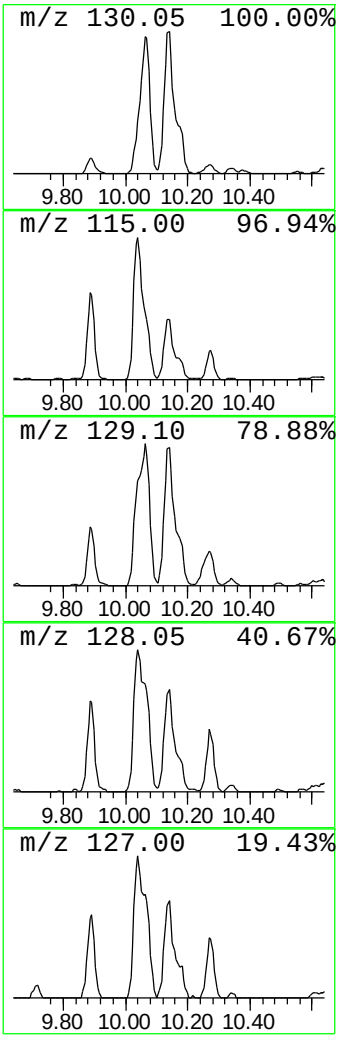
Peak Number 23 2-Methylindene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.14	5.08 ng/ul	627685	Naphthalene-d8	10.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylindene	130	C10H10	002177-47-1	97
2	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	97
3	1,4-Dihydronaphthalene	130	C10H10	000612-17-9	95
4	Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	94
5	1H-Indene, 3-methyl-	130	C10H10	000767-60-2	94





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Sample : K4706-09DL 5X
Misc :
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Instrument : BNA_M
ClientSampleId : BF575DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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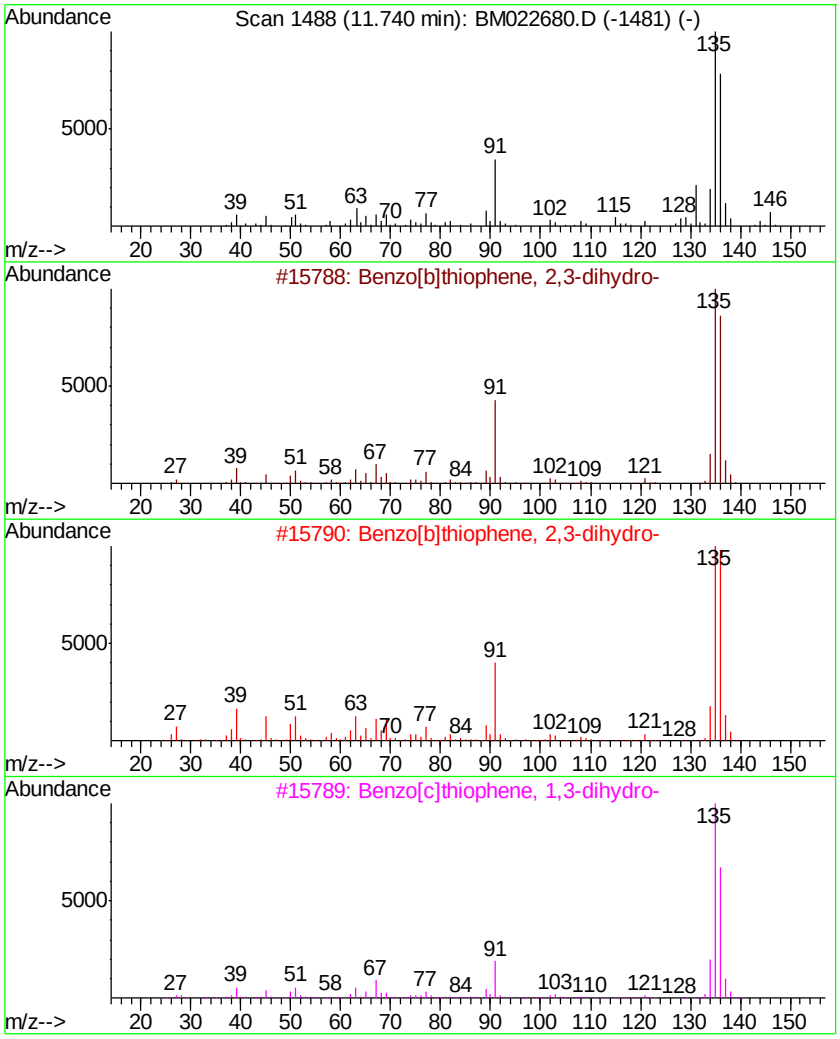
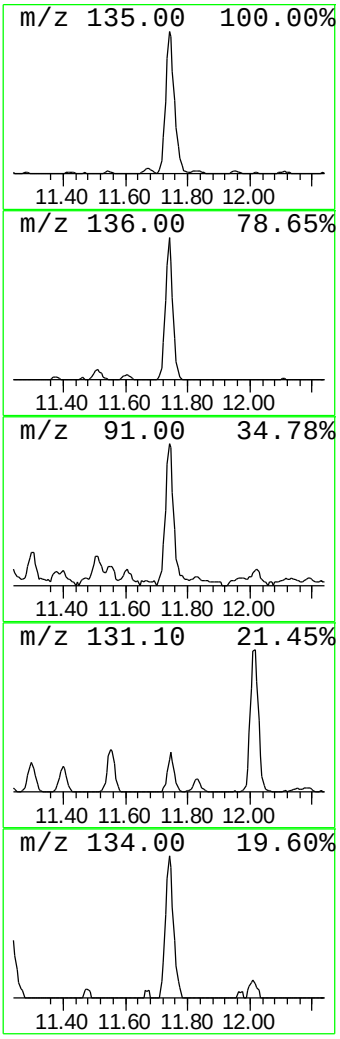
Peak Number 24 Benzo[b]thiophene, 2,3-dihy... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	2.73 ng/ul	337170	Naphthalene-d8	10.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	93
2	Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	81
3	Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	76
4	Benzene, (ethenylthio)-	136	C8H8S	001822-73-7	59
5	Benzaldehyde, 4-methoxy-	136	C8H8O2	000123-11-5	53





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Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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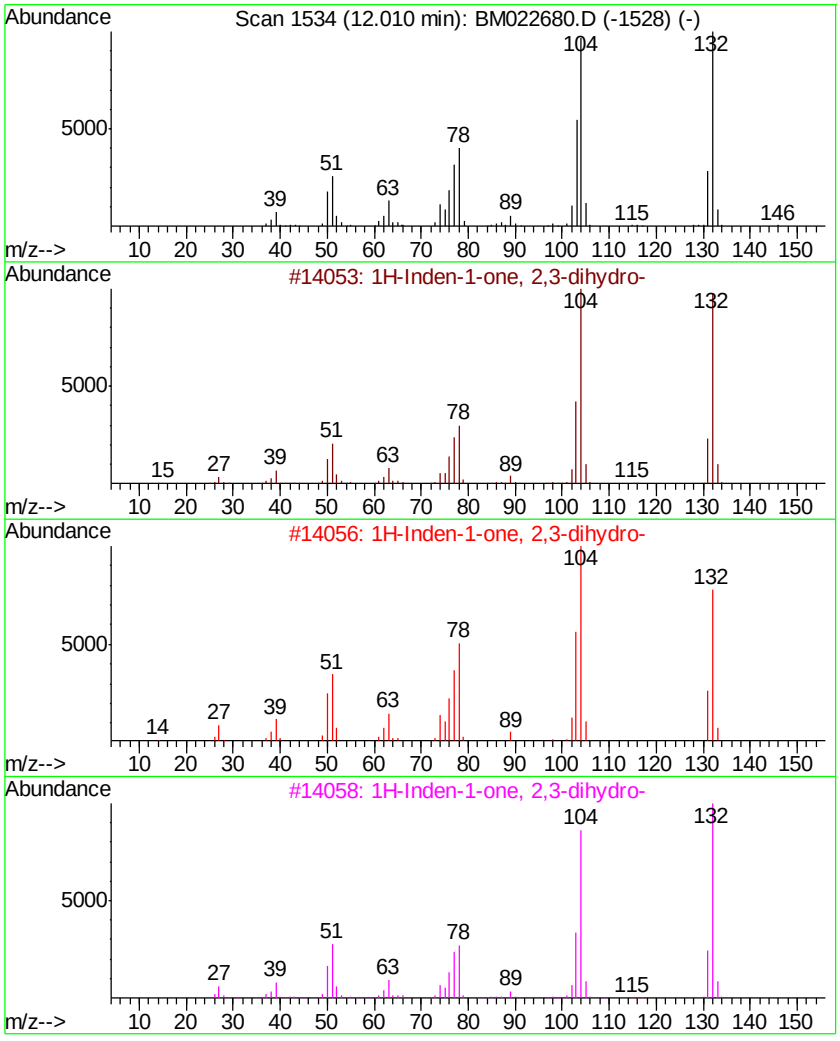
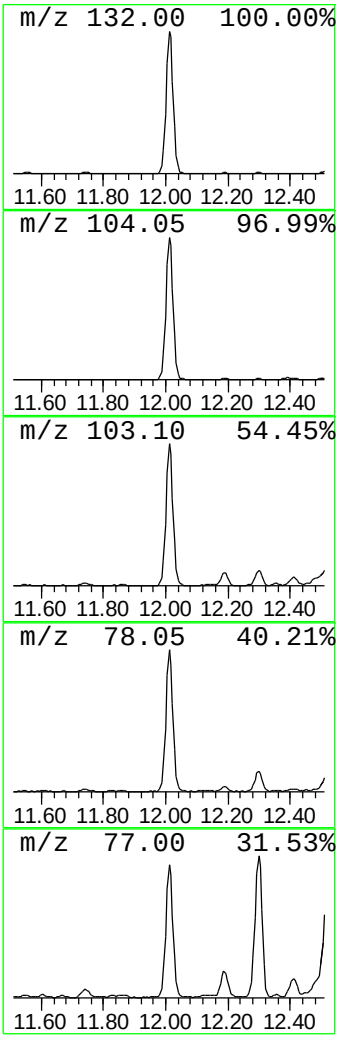
Peak Number 25 1H-Inden-1-one, 2,3-dihydro- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	9.08 ng/ul	1121330	Naphthalene-d8	10.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	97
2	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	95
3	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	93
4	Styrene	104	C8H8	000100-42-5	64
5	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	64





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Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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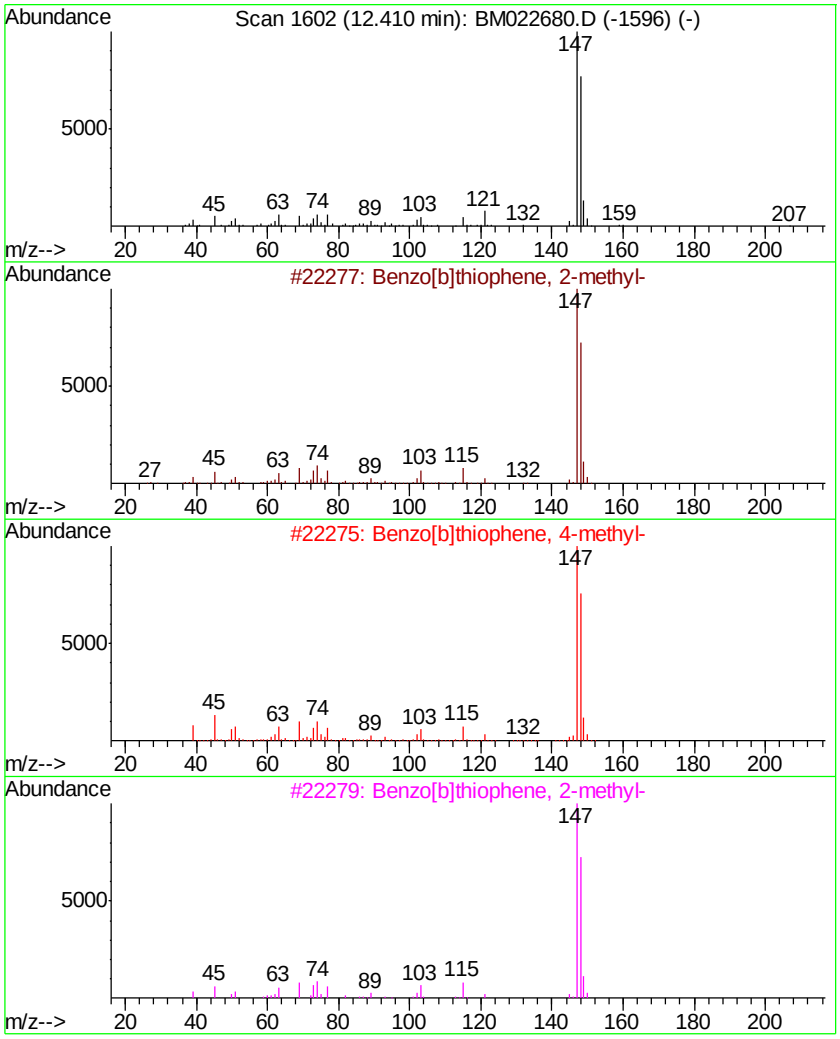
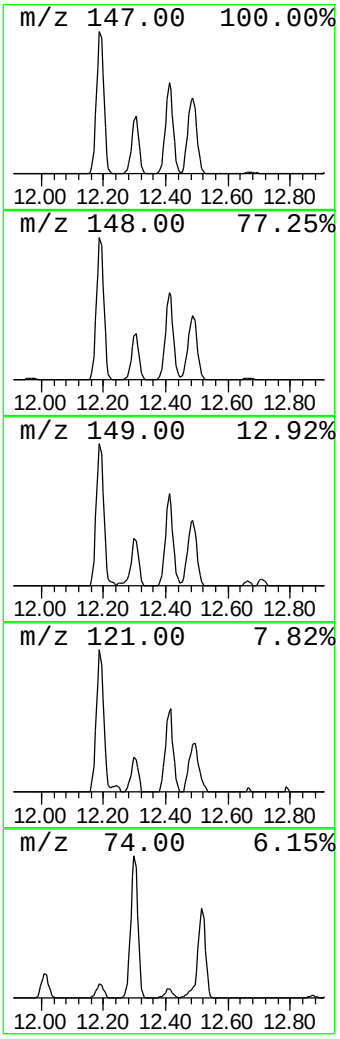
Peak Number 27 Benzo[b]thiophene, 4-methyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.41	3.86 ng/ul	476322	Naphthalene-d8	10.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
2	Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	91
3	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
4	3-Methylbenzothiophene	148	C9H8S	001455-18-1	91
5	Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	91





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Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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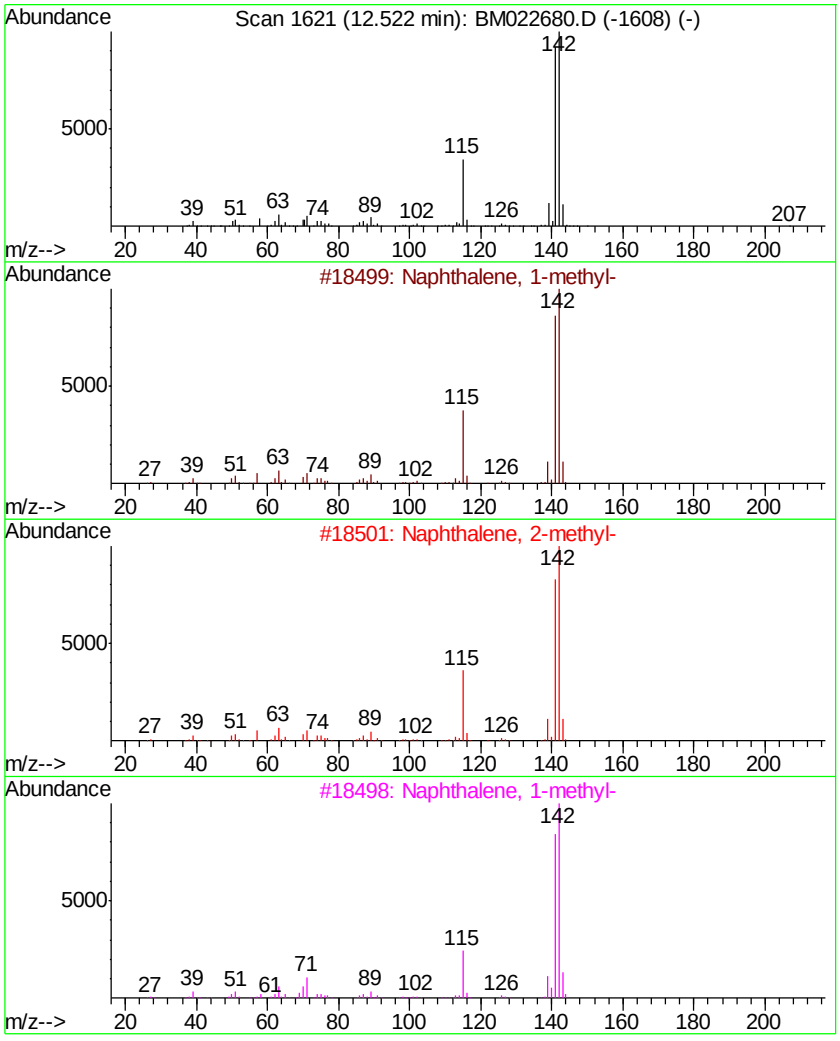
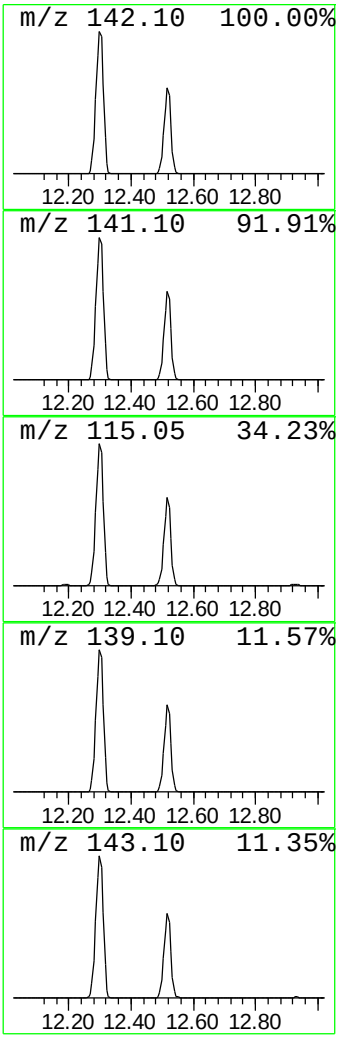
Peak Number 28 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.52	74.35 ng/ul	9183660	Naphthalene-d8	10.64

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	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91





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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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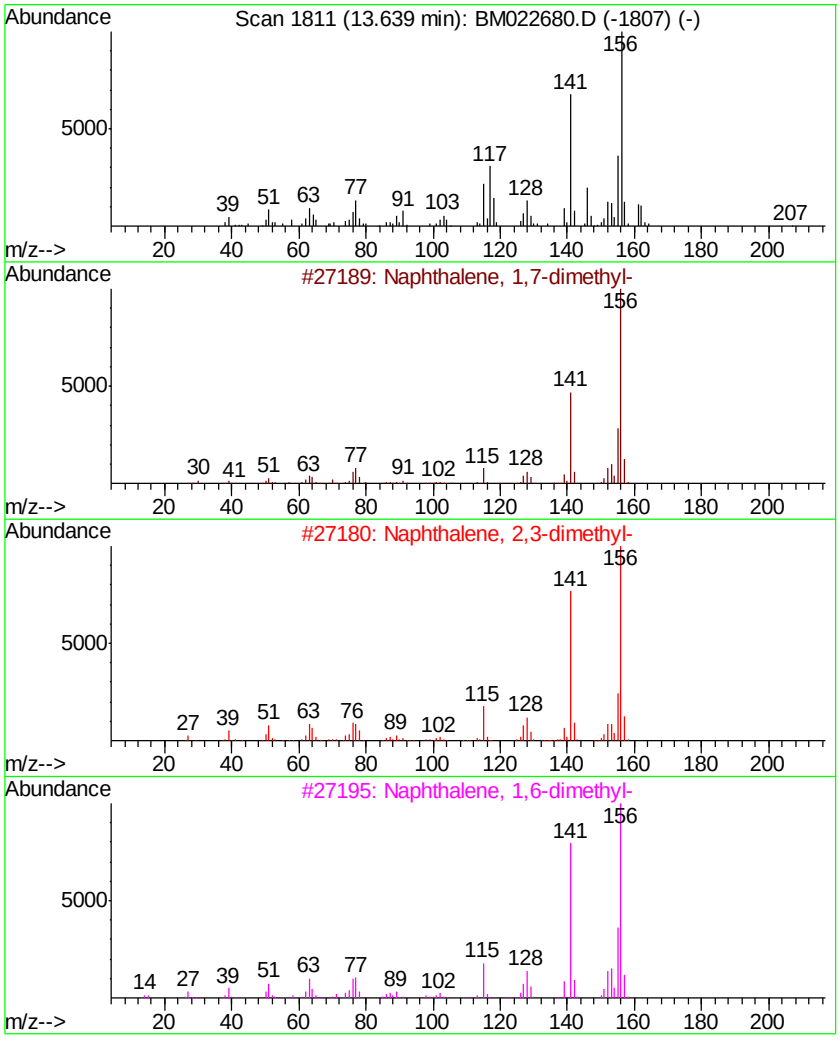
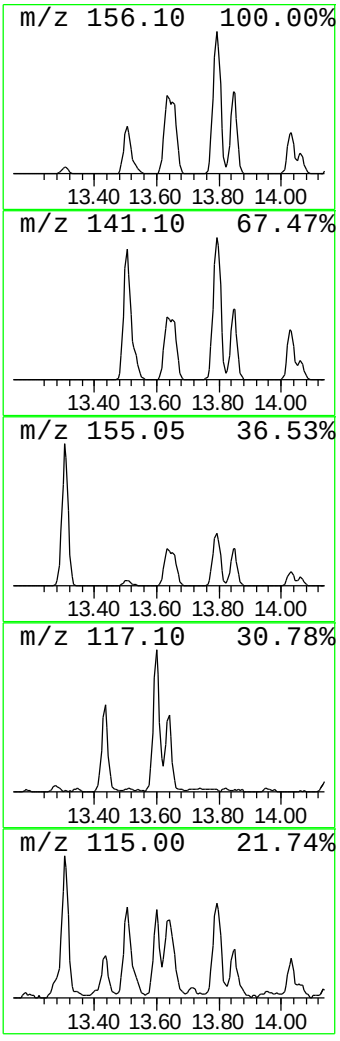
Peak Number 29 Naphthalene, 1,7-dimethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.64	3.25 ng/ul	545701	Acenaphthene-d10	14.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091319\
Data File : BM022680.D
Acq On : 13 Sep 2019 17:11
Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

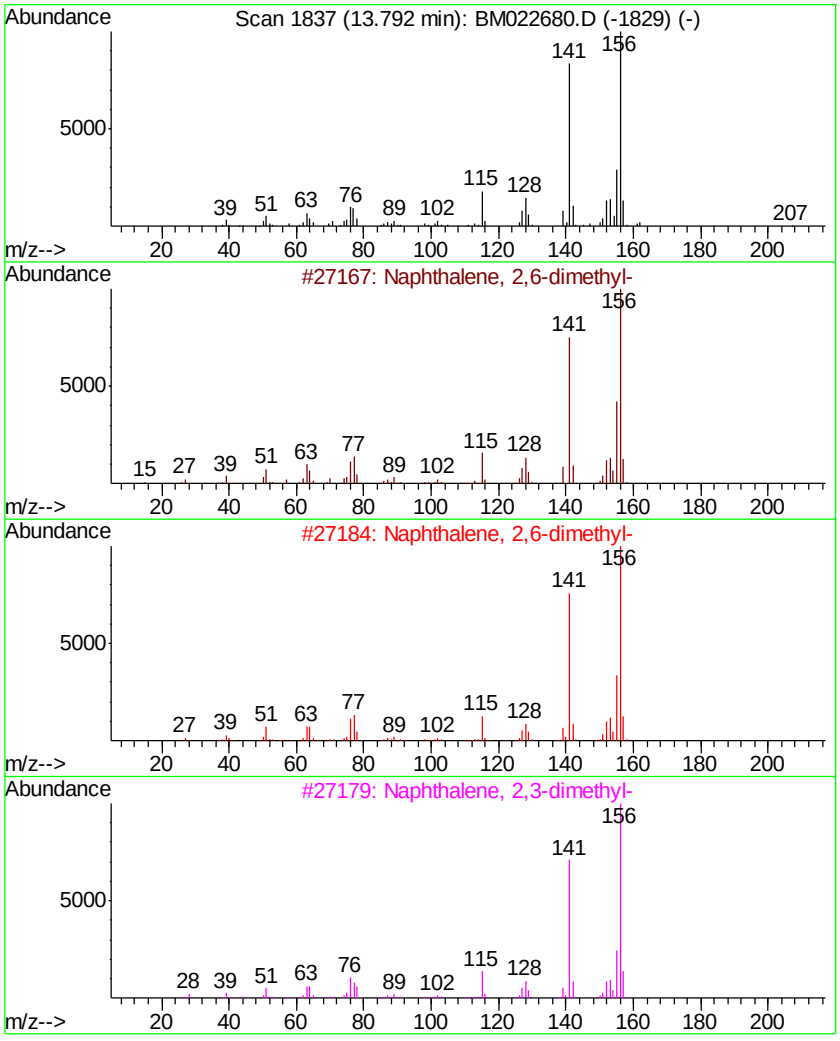
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

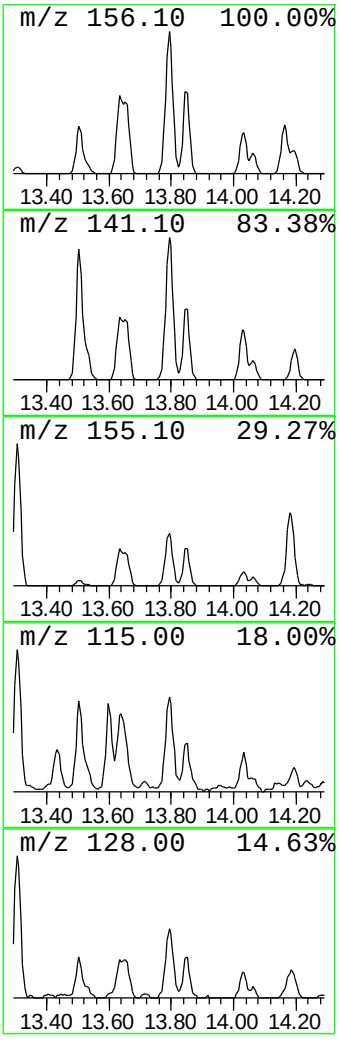
****

Peak Number 30 Naphthalene, 2,6-dimethyl- Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.79	3.14 ng/ul	527567	Acenaphthene-d10	14.47

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM091319\
Data File : BM022680.D
Acq On : 13 Sep 2019 17:11
Operator : HP/JU
Sample : K4706-09DL 5X
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BF575DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090519MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methoxy...	3.03	9.8	ng/ul	737638	1	7.84	1498340	20.0
Ethylbenzene	5.36	4.3	ng/ul	319242	1	7.84	1498340	20.0
o-Xylene	5.49	14.0	ng/ul	1048370	1	7.84	1498340	20.0
p-Xylene	5.86	9.3	ng/ul	695889	1	7.84	1498340	20.0
Benzene, 1-ethyl-...	6.93	13.6	ng/ul	1020720	1	7.84	1498340	20.0
Benzene, 1,2,3-tr...	7.07	17.9	ng/ul	1336930	1	7.84	1498340	20.0
Benzene, 1-ethyl-...	7.23	3.7	ng/ul	274133	1	7.84	1498340	20.0
Benzene, 1,2,4-tr...	7.49	39.9	ng/ul	2990420	1	7.84	1498340	20.0
Benzofuran	7.56	12.4	ng/ul	927063	1	7.84	1498340	20.0
Benzene, 1,3,5-tr...	7.96	16.4	ng/ul	1229050	1	7.84	1498340	20.0
Indane	8.22	200.3	ng/ul	15002300	1	7.84	1498340	20.0
Indene	8.38	120.8	ng/ul	9048900	1	7.84	1498340	20.0
Benzene, 2-ethyl-...	8.49	7.2	ng/ul	542207	1	7.84	1498340	20.0
Benzene, 1-methyl...	8.80	2.0	ng/ul	152622	1	7.84	1498340	20.0
Benzene, 1-methyl...	8.85	2.2	ng/ul	162493	1	7.84	1498340	20.0
Benzofuran, 2-met...	9.20	23.0	ng/ul	1722910	1	7.84	1498340	20.0
Benzofuran, 7-met...	9.32	35.2	ng/ul	4349630	2	10.64	2470540	20.0
2-Propenal, 3-phe...	9.38	7.8	ng/ul	964541	2	10.64	2470540	20.0
Benzene, 1,2,4,5-...	9.47	3.6	ng/ul	441117	2	10.64	2470540	20.0
Benzene, 1,2,3,4-...	9.53	4.2	ng/ul	515623	2	10.64	2470540	20.0
Benzene, 1-etheny...	9.89	11.3	ng/ul	1391360	2	10.64	2470540	20.0
Benzene, 2-etheny...	10.04	22.7	ng/ul	2801940	2	10.64	2470540	20.0
2-Methylindene	10.14	5.1	ng/ul	627685	2	10.64	2470540	20.0
Benzo[b]thiophene...	11.74	2.7	ng/ul	337170	2	10.64	2470540	20.0
1H-Inden-1-one, 2...	12.01	9.1	ng/ul	1121330	2	10.64	2470540	20.0
Benzo[b]thiophene...	12.41	3.9	ng/ul	476322	2	10.64	2470540	20.0
Naphthalene, 1-me...	12.52	74.3	ng/ul	9183660	2	10.64	2470540	20.0
Naphthalene, 1,7-...	13.64	3.3	ng/ul	545701	3	14.47	3359100	20.0
Naphthalene, 2,6-...	13.79	3.1	ng/ul	527567	3	14.47	3359100	20.0