LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
Data File : BN010391.D
Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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_N\METHODS\SOM-EPA-BN040120MA.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	5.311	424	429	440	rBV	298913	589512	0.37%	0.139%
2	5.646	480	486	488	rBV	170746	267606	0.17%	0.063%
3	5.669	488	490	501	rVB	190234	269072	0.17%	0.063%
4	6.334	594	603	612	rBV	1058080	1663625	1.06%	0.391%
5	6.628	649	653	660	rVV	239637	383488	0.24%	0.090%
6	6.728	663	670	675	rVV	395219	598815	0.38%	0.141%
7	6.781	675	679	687	rVV2	193550	354463	0.22%	0.083%
8	6.857	687	692	700	rVB	418159	656982	0.42%	0.154%
9	7.069	723	728	737	rVV	366172	670834	0.43%	0.158%
10	7.275	756	763	769	rBV2	1343057	2120376	1.35%	0.498%
11	7.340	769	774	786	rVB	676393	1263047	0.80%	0.297%
12	7.610	813	820	827	rBV	1638439	2606662	1.65%	0.613%
13	7.757	834	845	850	rBV2	1165267	2258034	1.43%	0.531%
14	7.840	854	859	865	rVV	517317	802872	0.51%	0.189%
15	7.981	877	883	892	rVB	3183215	4979147	3.16%	1.170%
16	8.140	899	910	921	rBV	8317446	13022358	8.26%	3.060%
17	8.740	1002	1012	1016	rBV3	307898	804002	0.51%	0.189%
18	8.787	1016	1020	1024	rVV	292232	541888	0.34%	0.127%
19	8.840	1024	1029	1041	rVB	837911	1436605	0.91%	0.338%
20	8.951	1041	1048	1056	rBV	477557	756418	0.48%	0.178%
21	9.069	1056	1068	1074	rBV	1021425	2142053	1.36%	0.503%
22	9.134	1074	1079	1085	rVV	346214	534360	0.34%	0.126%
23	9.463	1127	1135	1140	rBV3	124529	271384	0.17%	0.064%
24	9.634	1155	1164	1171	rVB2	350852	670785	0.43%	0.158%
25	9.757	1171	1185	1188	rBV	1706442	3412276	2.17%	0.802%
26	9.798	1188	1192	1201	rVV4	1999551	4248286	2.70%	0.998%
27	9.881	1201	1206	1210	rVV	586251	1125635	0.71%	0.264%
28	9.922	1210	1213	1219	rVV3	380162	648261	0.41%	0.152%
29	9.998	1219	1226	1235	rVB4	343549	893656	0.57%	0.210%
30	10.257	1261	1270	1276	rBV7	182269	470102	0.30%	0.110%
31	10.381	1284	1291	1292	rBV	1820997	2849580	1.81%	0.670%
32	10.451	1292	1303	1309	rVV2	56330599	157561174	100.00%	37.023%
33	10.546	1309	1319	1329	rVB	3747034	6349278	4.03%	1.492%
34	11.198	1424	1430	1439	rVV	2088010	3203412	2.03%	0.753%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
Data File : BN010391.D
Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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_N\METHODS\SOM-EPA-BN040120MA.M
Title : SVOA CALIBRATION

35	11.928	1548	1554	1561	rVB	511826	825785	0.52%	0.194%
36	12.040	1561	1573	1581	rBV	20209860	33783089	21.44%	7.938%
37	12.151	1586	1592	1598	rVV	621795	1166510	0.74%	0.274%
38	12.263	1598	1611	1621	rVB	10157353	17007112	10.79%	3.996%
39	12.687	1671	1683	1690	rBV2	164474	356488	0.23%	0.084%
40	13.063	1734	1747	1755	rBV	3805346	5423020	3.44%	1.274%
41	13.263	1775	1781	1792	rVB	799726	1414964	0.90%	0.332%
42	13.398	1793	1804	1805	rBV2	717452	1212616	0.77%	0.285%
43	13.557	1823	1831	1836	rVV	1148111	2032837	1.29%	0.478%
44	13.610	1836	1840	1843	rVV	782694	1141109	0.72%	0.268%
45	13.651	1843	1847	1855	rVB	573426	864050	0.55%	0.203%
46	13.792	1865	1871	1874	rBV	465702	663221	0.42%	0.156%
47	13.922	1887	1893	1896	rVV	622623	963626	0.61%	0.226%
48	13.957	1896	1899	1910	rVB2	408687	634553	0.40%	0.149%
49	14.234	1938	1946	1952	rBV2	3845193	6287598	3.99%	1.477%
50	14.304	1952	1958	1963	rVV	14926902	21711027	13.78%	5.102%
51	14.363	1963	1968	1974	rVB	2715699	3963124	2.52%	0.931%
52	14.516	1989	1994	1997	rBV2	334637	505594	0.32%	0.119%
53	14.551	1997	2000	2005	rVV	279299	366376	0.23%	0.086%
54	14.639	2005	2015	2023	rVV2	8687314	13060812	8.29%	3.069%
55	14.716	2023	2028	2038	rVB	355064	558217	0.35%	0.131%
56	15.233	2111	2116	2120	rVV	901596	1316228	0.84%	0.309%
57	15.286	2120	2125	2131	rVV2	6878239	9669533	6.14%	2.272%
58	15.416	2143	2147	2150	rVV3	305229	509437	0.32%	0.120%
59	15.451	2150	2153	2158	rVV	477260	698232	0.44%	0.164%
60	15.504	2158	2162	2164	rVV2	175651	284521	0.18%	0.067%
61	15.533	2164	2167	2172	rVV	245314	422679	0.27%	0.099%
62	15.586	2172	2176	2184	rVV	389060	579878	0.37%	0.136%
63	15.728	2195	2200	2208	rVV2	457043	1002077	0.64%	0.235%
64	15.957	2233	2239	2249	rBV	1084196	1583680	1.01%	0.372%
65	16.086	2253	2261	2265	rVV	370203	585623	0.37%	0.138%
66	16.175	2272	2276	2284	rVV2	351349	576931	0.37%	0.136%
67	16.251	2284	2289	2294	rVV2	304633	622231	0.39%	0.146%
68	16.604	2343	2349	2351	rBV3	170958	291683	0.19%	0.069%
69	16.798	2378	2382	2393	rVB	576637	833768	0.53%	0.196%
70	16.980	2408	2413	2417	rVV	5189952	7611362	4.83%	1.788%
71	17.022	2417	2420	2425	rVV	7282691	9675791	6.14%	2.274%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
Data File : BN010391.D
Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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_N\METHODS\SOM-EPA-BN040120MA.M
Title : SVOA CALIBRATION

72	17.080	2425	2430	2433	rVV	1977695	2909415	1.85%	0.684%
73	17.110	2433	2435	2442	rVV	544671	786575	0.50%	0.185%
74	17.386	2471	2482	2488	rBV	8422399	12415163	7.88%	2.917%
75	17.480	2493	2498	2503	rVV2	403894	728467	0.46%	0.171%
76	17.539	2503	2508	2514	rVV2	332408	532480	0.34%	0.125%
77	17.816	2547	2555	2559	rVV2	171238	319491	0.20%	0.075%
78	17.869	2559	2564	2566	rVV2	172874	343214	0.22%	0.081%
79	17.898	2566	2569	2577	rVV3	798079	1269539	0.81%	0.298%
80	17.975	2577	2582	2586	rVV2	255498	363255	0.23%	0.085%
81	18.051	2587	2595	2600	rVV3	464324	959030	0.61%	0.225%
82	18.157	2606	2613	2617	rVV2	295741	706346	0.45%	0.166%
83	18.198	2617	2620	2624	rVV	309092	483973	0.31%	0.114%
84	18.298	2631	2637	2642	rVV3	300059	576423	0.37%	0.135%
85	18.357	2643	2647	2651	rVV2	250533	394343	0.25%	0.093%
86	19.045	2759	2764	2770	rVB	787661	1093804	0.69%	0.257%
87	19.380	2813	2821	2824	rBV	1402129	2101842	1.33%	0.494%
88	19.410	2824	2826	2838	rVB2	527877	778935	0.49%	0.183%
89	19.786	2882	2890	2895	rBV2	295883	675467	0.43%	0.159%
90	19.845	2895	2900	2908	rVV	489358	735222	0.47%	0.173%
91	20.463	2998	3005	3008	rBV2	180139	412693	0.26%	0.097%
92	20.504	3008	3012	3018	rVB2	239342	439433	0.28%	0.103%
93	21.180	3122	3127	3138	rVB	7933516	9689106	6.15%	2.277%
94	22.251	3307	3309	3314	rVB3	246042	282387	0.18%	0.066%
95	22.745	3389	3393	3400	rVB2	452036	633581	0.40%	0.149%
96	23.215	3468	3473	3480	rBV	1118995	1891384	1.20%	0.444%
97	23.292	3482	3486	3490	rVV2	559927	885958	0.56%	0.208%
98	23.357	3490	3497	3506	rVB2	5763011	10590013	6.72%	2.488%
99	23.915	3588	3592	3600	rVB	524327	930705	0.59%	0.219%
100	24.627	3709	3713	3721	rVB4	481510	1017075	0.65%	0.239%

Sum of corrected areas: 425576749

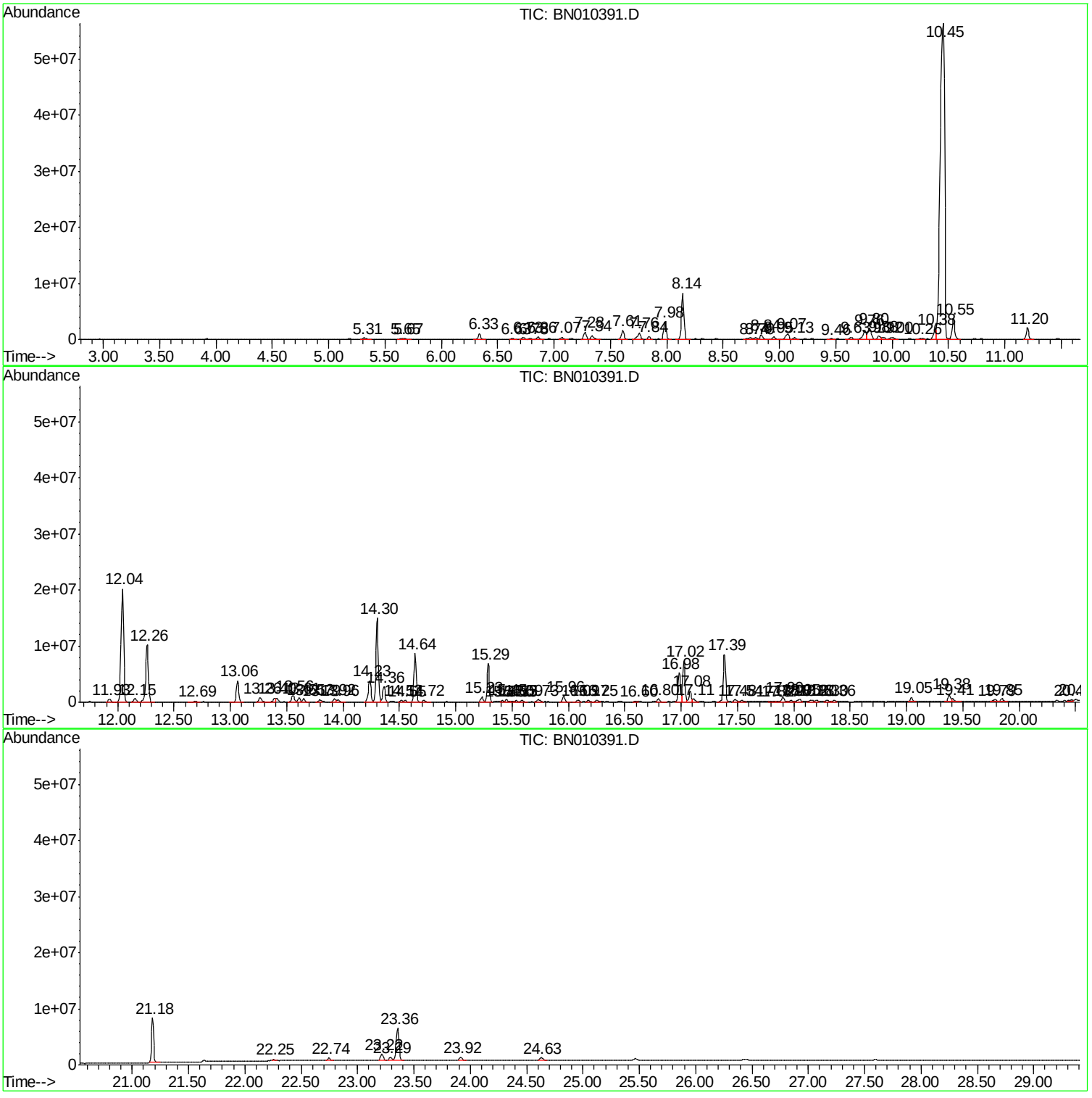
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
Data File : BN010391.D
Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
Data File : BN010391.D
Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

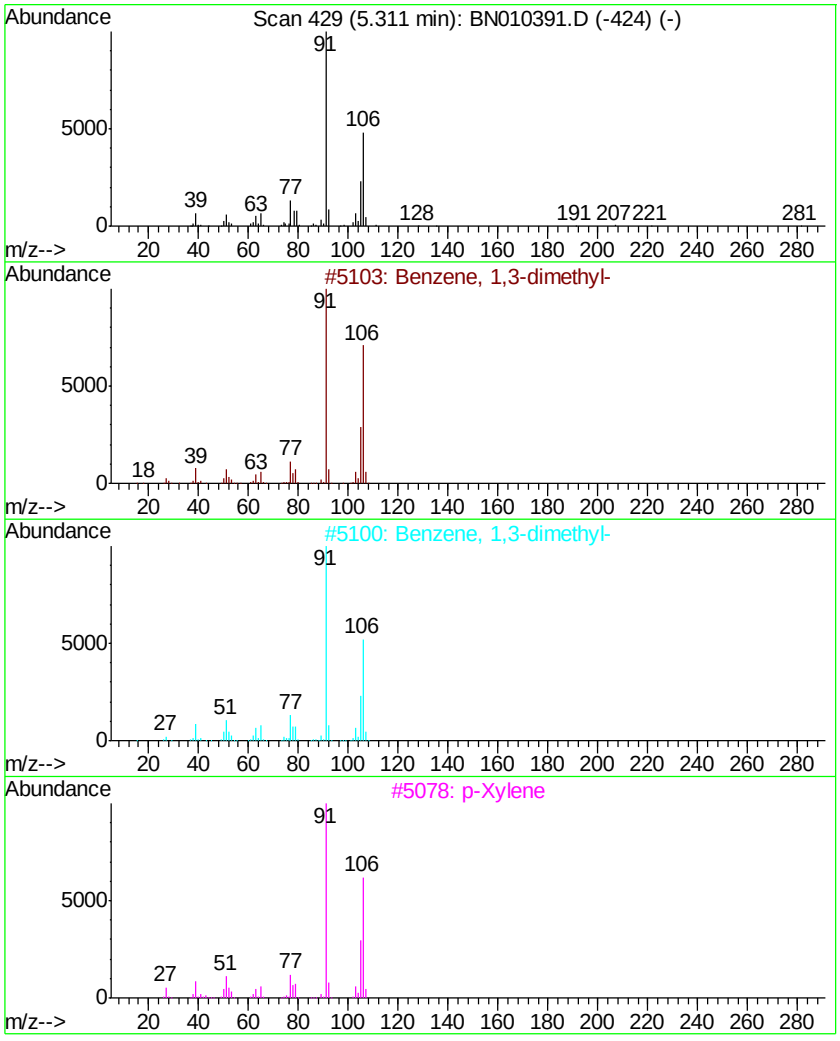
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

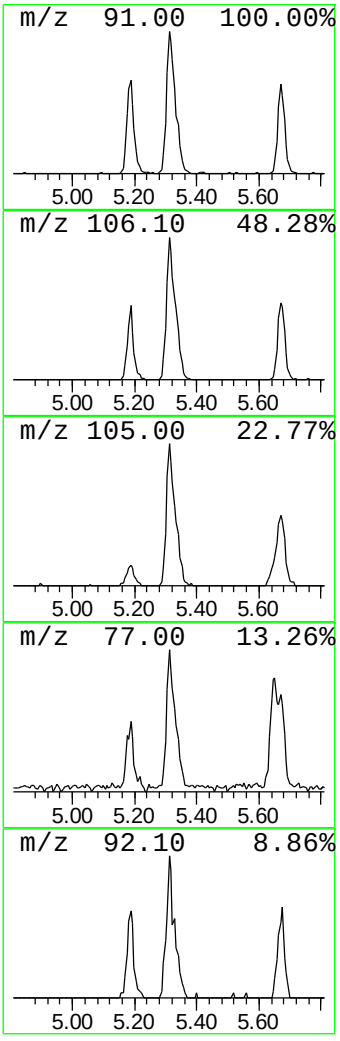
****

Peak Number 1 Benzene, 1,3-dimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.31	4.52 ng/ul	589512	1,4-Dichlorobenzene-d4	7.61

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





Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
Data File : BN010391.D
Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

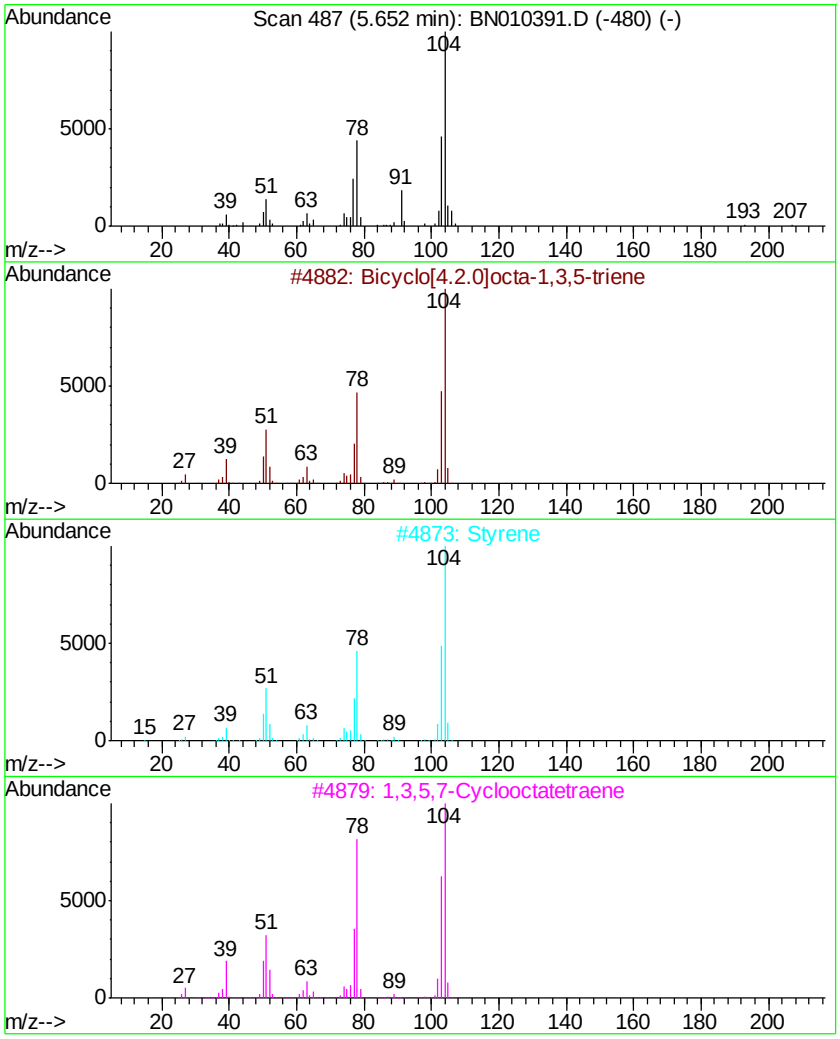
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

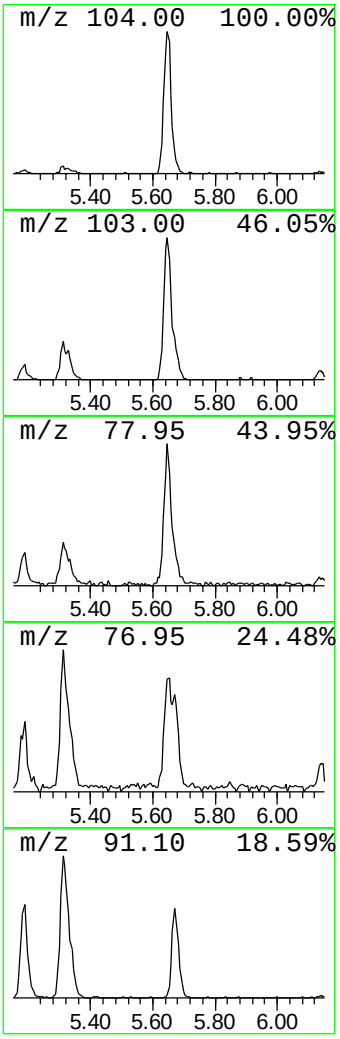
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.65	2.05 ng/ul	267606	1,4-Dichlorobenzene-d4	7.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.2.0]octa-1,3,5-triene	104	C8H8	000694-87-1	93
2	Styrene	104	C8H8	000100-42-5	93
3	1,3,5,7-Cyclooctatetraene	104	C8H8	000629-20-9	91
4	1,3,5,7-Cyclooctatetraene	104	C8H8	000629-20-9	91
5	1,3,5,7-Cyclooctatetraene	104	C8H8	000629-20-9	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
Data File : BN010391.D
Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

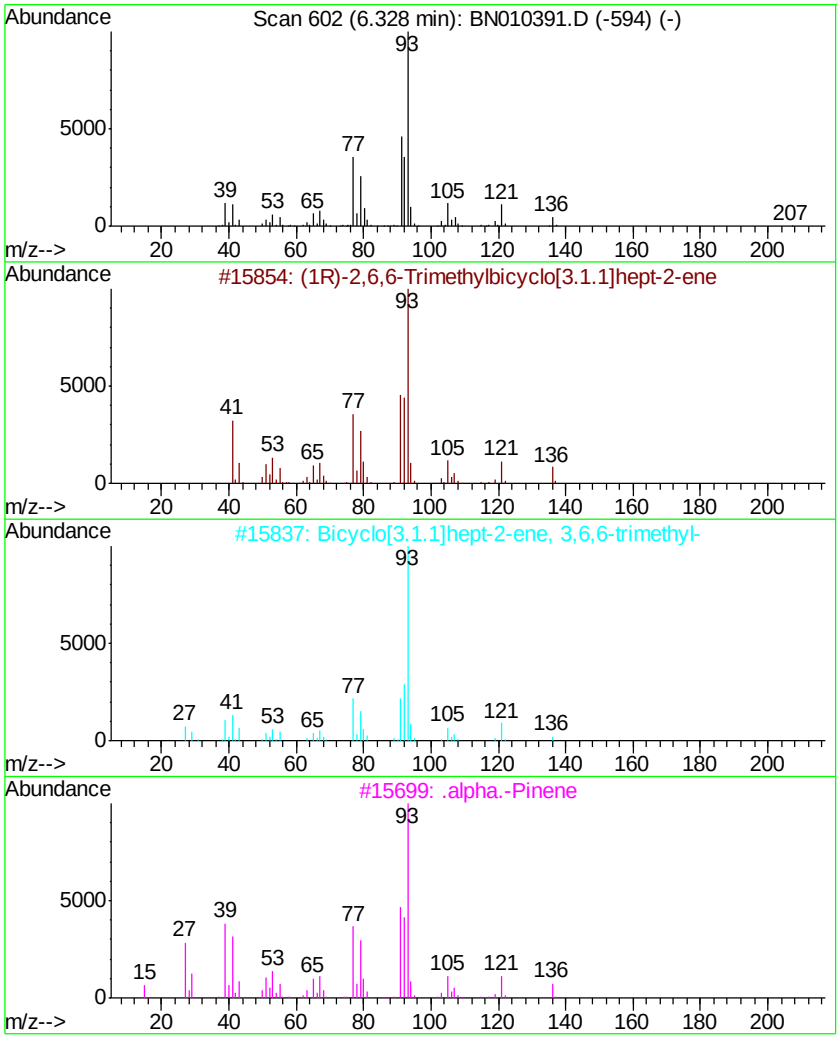
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

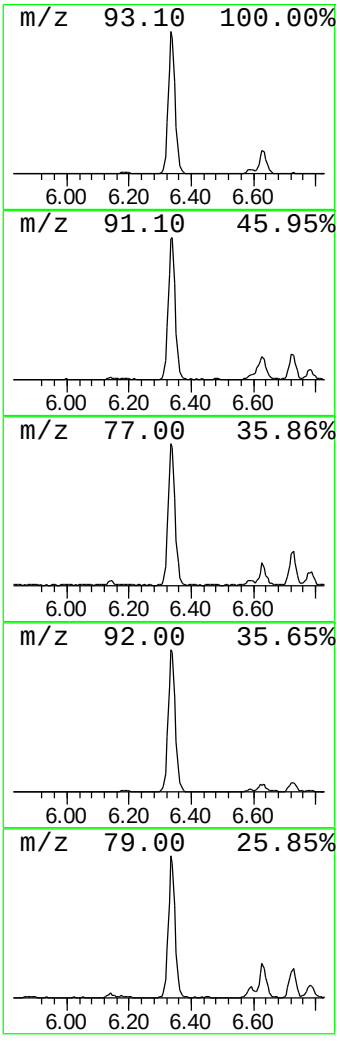
****

Peak Number 4 (1R)-2,6,6-Trimethylbicyclo... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.33	12.76 ng/ul	1663630	1,4-Dichlorobenzene-d4	7.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1R)-2,6,6-Trimethylbicyclo[3.1....	136	C10H16	007785-70-8	95
2	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	94
3	.alpha.-Pinene	136	C10H16	000080-56-8	94
4	.gamma.-Terpinene	136	C10H16	000099-85-4	91
5	(+)-3-Carene	136	C10H16	000498-15-7	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
Data File : BN010391.D
Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

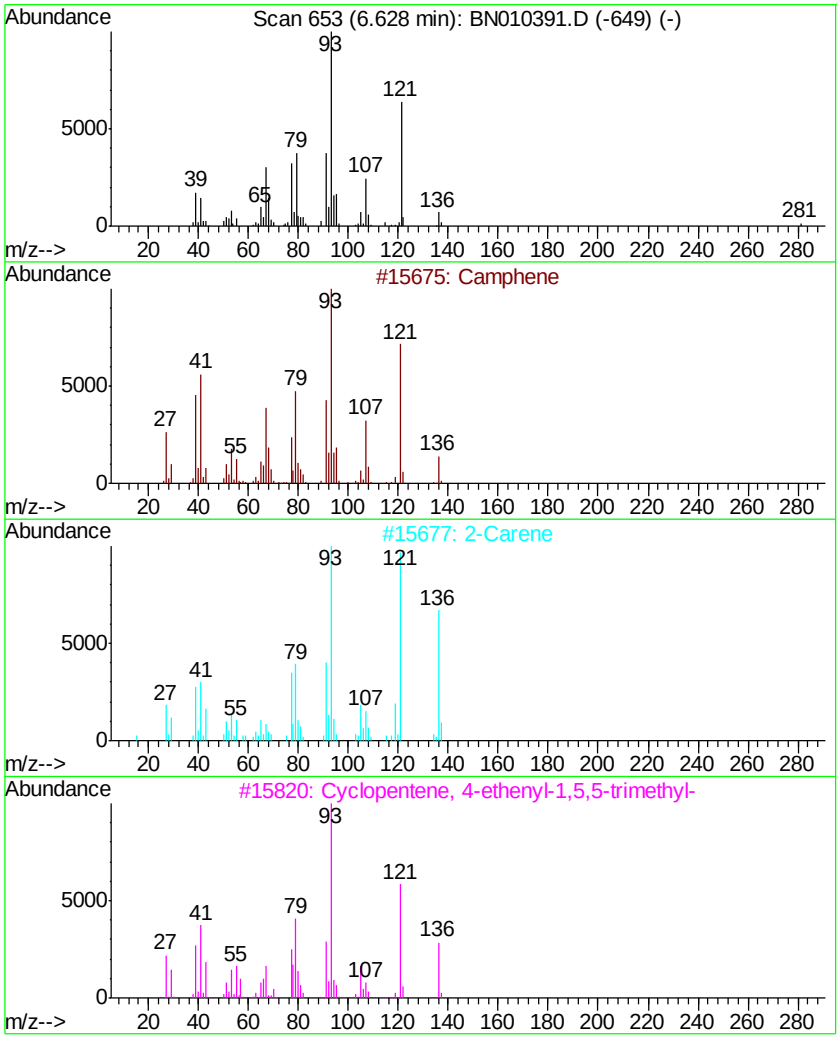
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

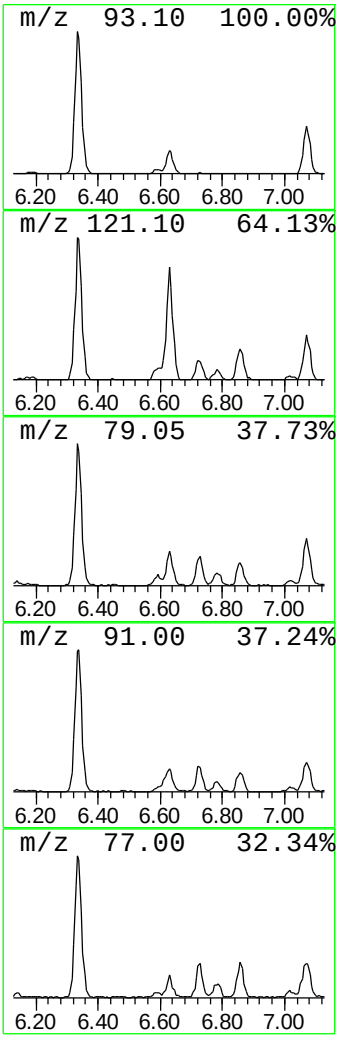
****

Peak Number 5 Camphene Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	2.94 ng/ul	383488	1,4-Dichlorobenzene-d4	7.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Camphene	136	C10H16	000079-92-5	80
2	2-Carene	136	C10H16	000554-61-0	72
3	Cyclopentene, 4-ethenyl-1,5,5-tr...	136	C10H16	001727-69-1	64
4	.beta.-Phellandrene	136	C10H16	000555-10-2	53
5	Bicyclo[2.2.1]heptane, 2-chloro-...	172	C10H17Cl	000559-45-5	53





Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040120\
Data File : BN010391.D
Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

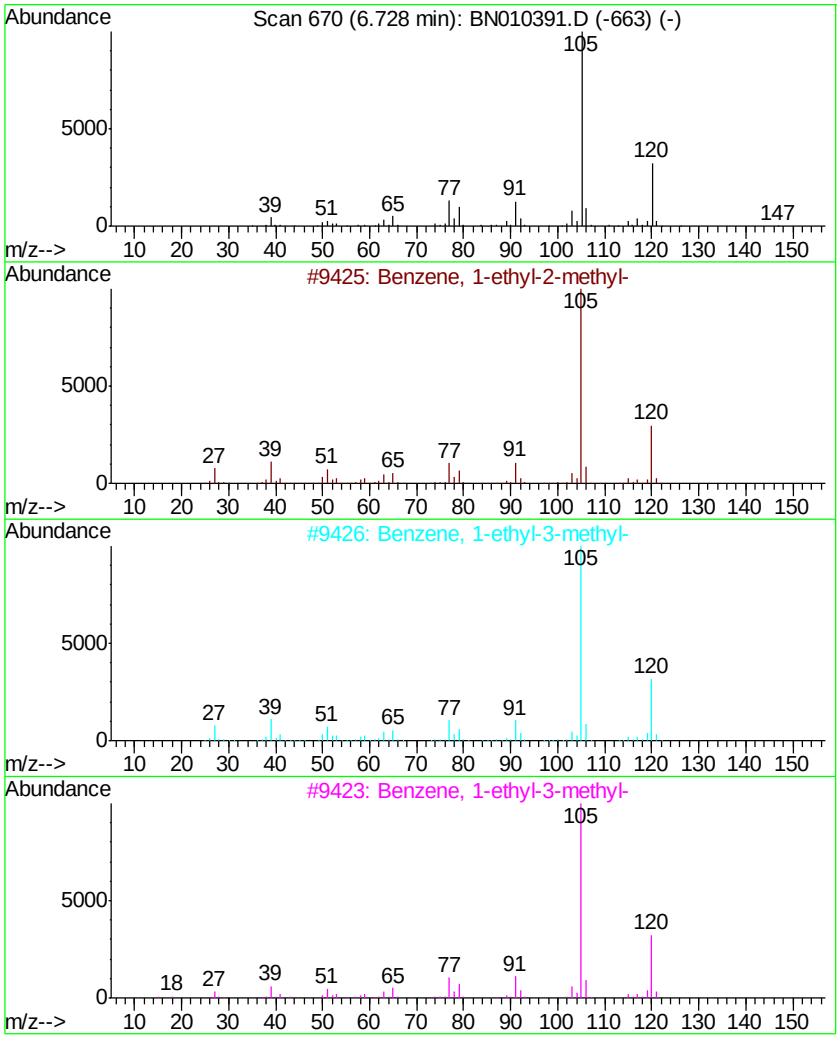
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

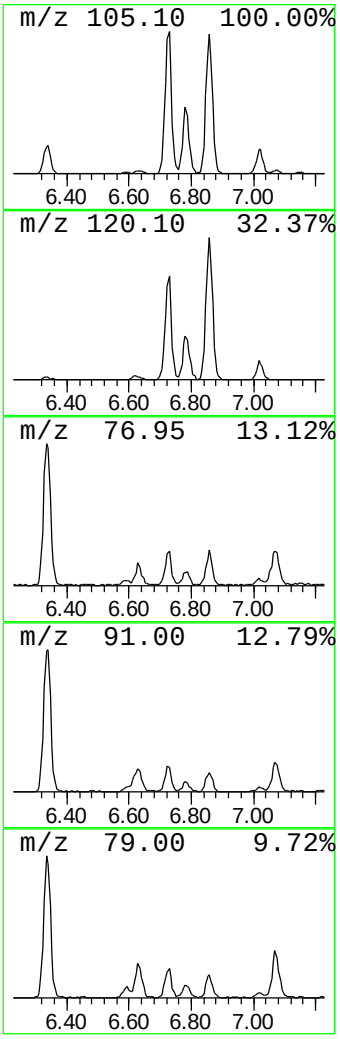
****

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	4.59 ng/ul	598815	1,4-Dichlorobenzene-d4	7.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





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Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

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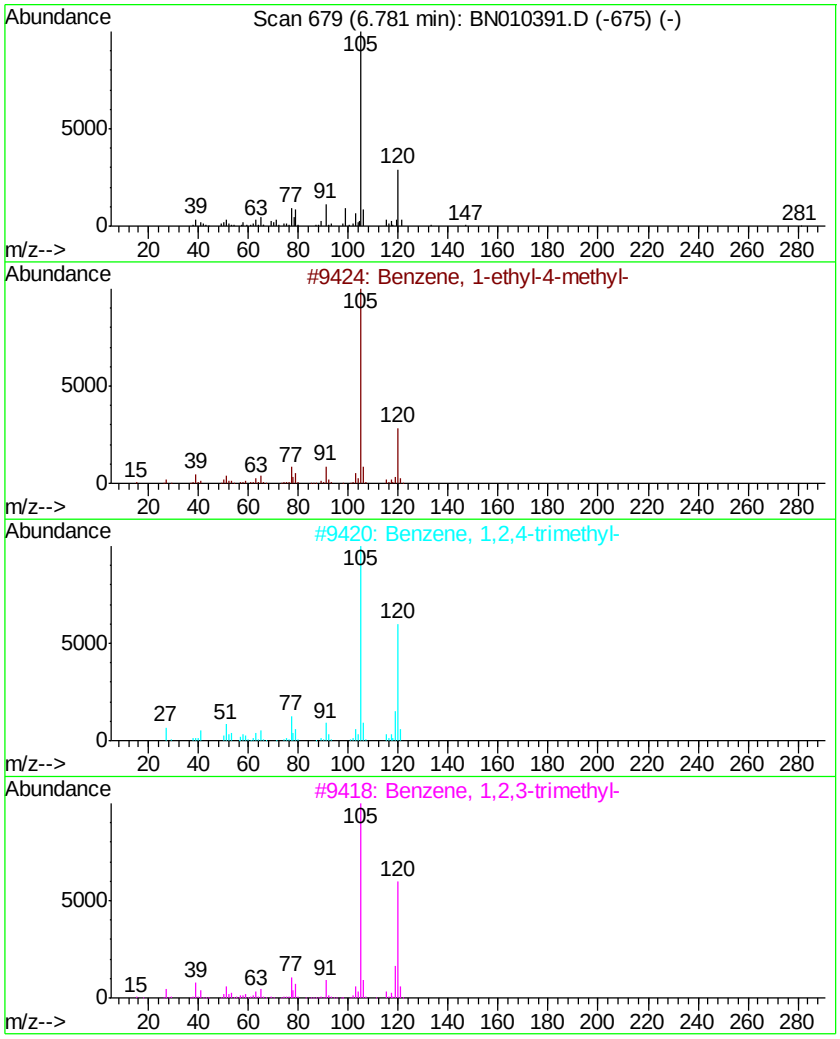
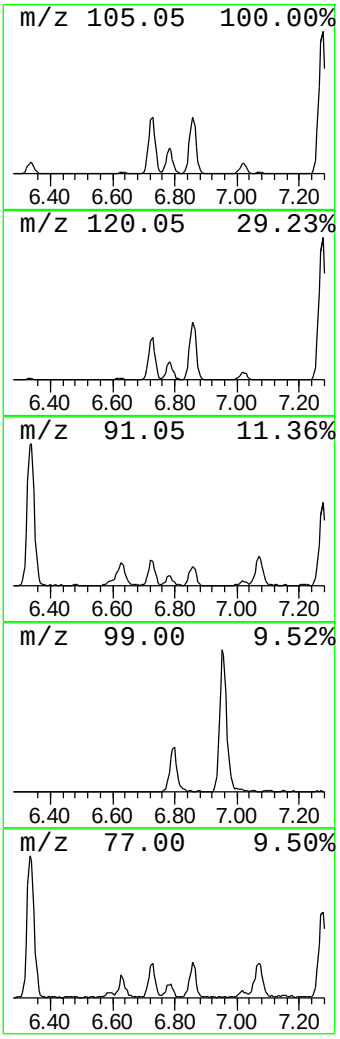
Peak Number 7 Benzene, 1-ethyl-4-methyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.78	2.72 ng/ul	354463	1,4-Dichlorobenzene-d4	7.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	87
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	87
4	Mesitylene	120	C9H12	000108-67-8	87
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87





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Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

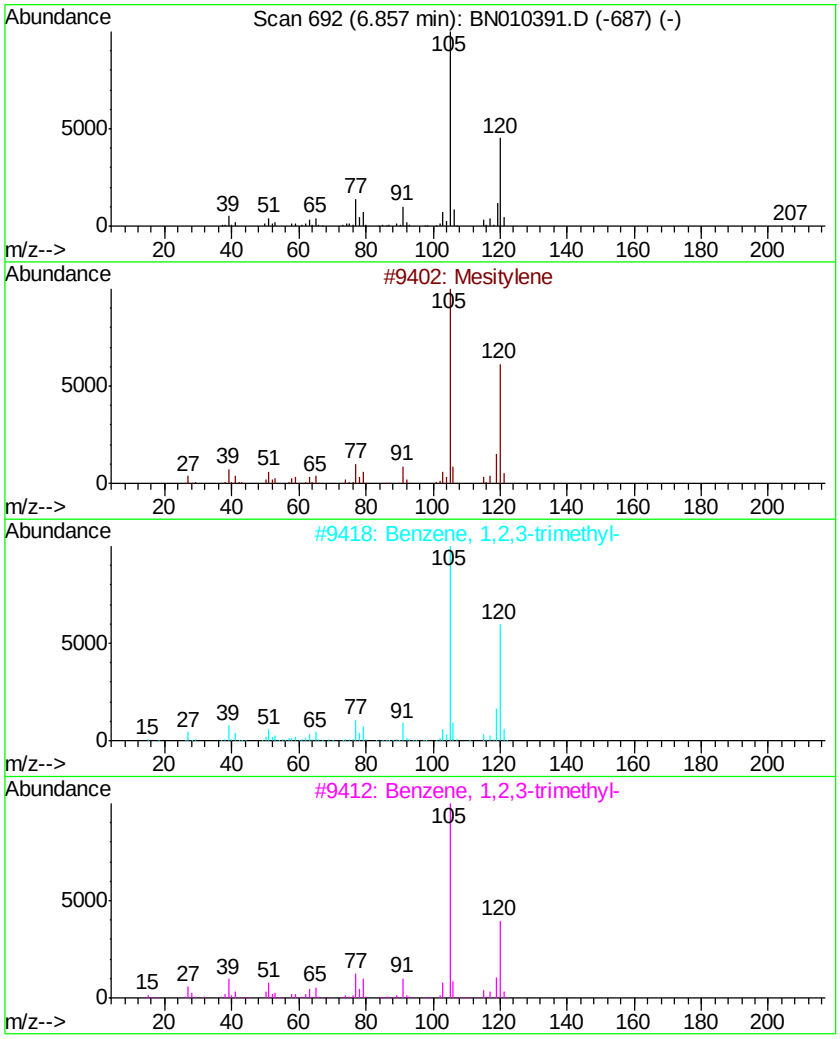
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

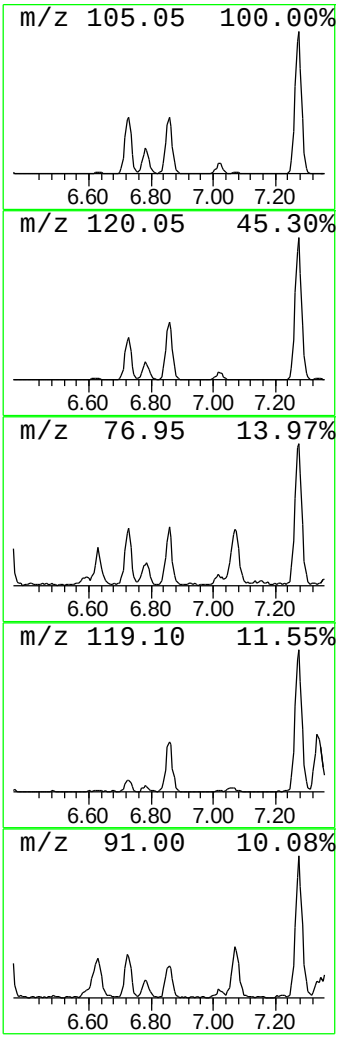
****

Peak Number 8 Mesitylene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.86	5.04 ng/ul	656982	1,4-Dichlorobenzene-d4	7.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93





Library Search Compound Report

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Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

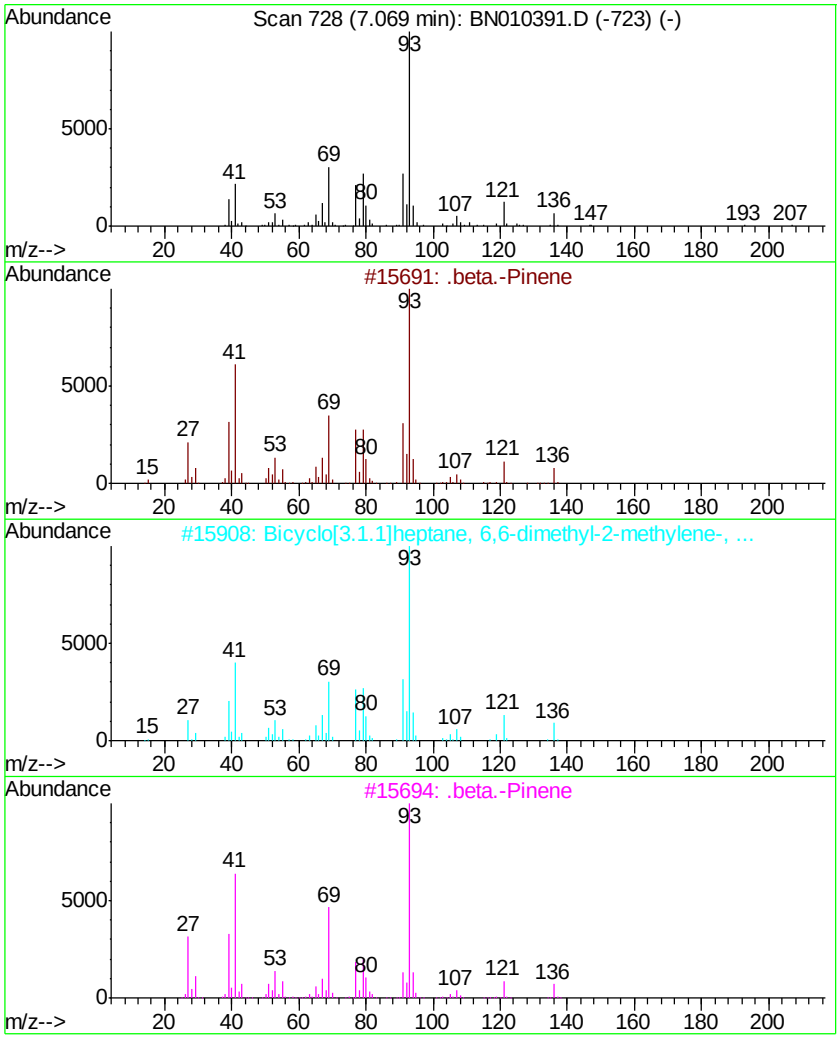
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

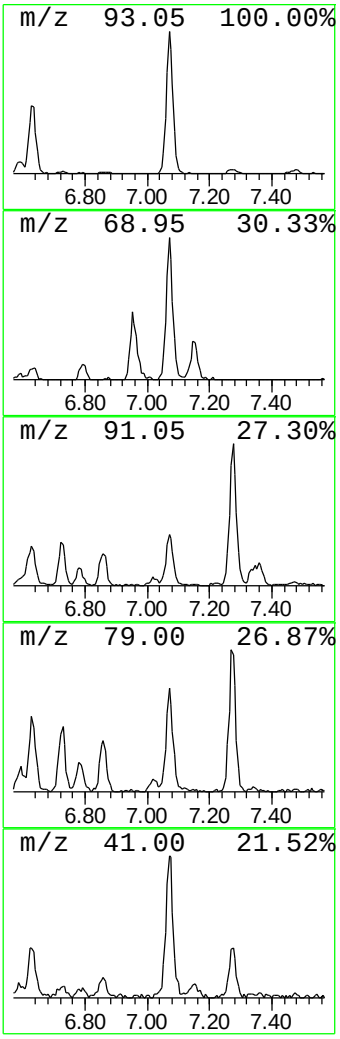
****

Peak Number 9 .beta.-Pinene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.07	5.15 ng/ul	670834	1,4-Dichlorobenzene-d4	7.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-Pinene	136	C10H16	000127-91-3	94
2	Bicyclo[3.1.1]heptane, 6,6-dimet...	136	C10H16	018172-67-3	91
3	.beta.-Pinene	136	C10H16	000127-91-3	91
4	.beta.-Pinene	136	C10H16	000127-91-3	91
5	Tricyclo[2.2.1.0(2,6)]heptane, 1...	136	C10H16	000508-32-7	91





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Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

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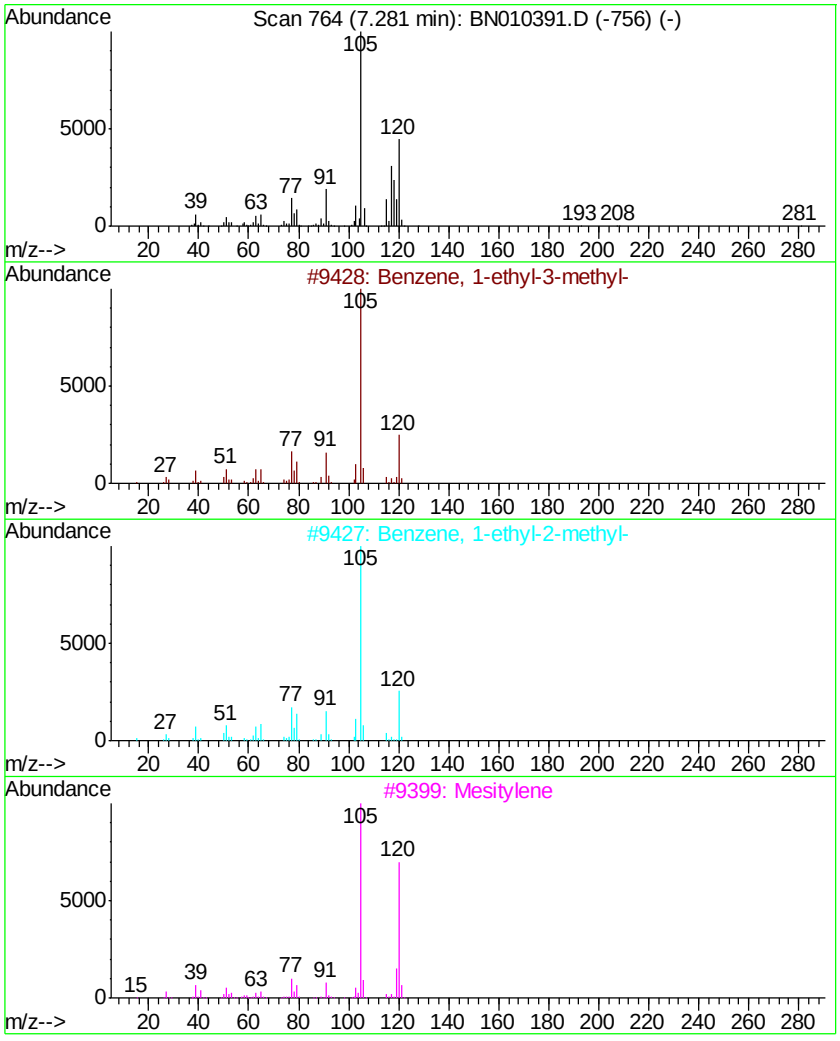
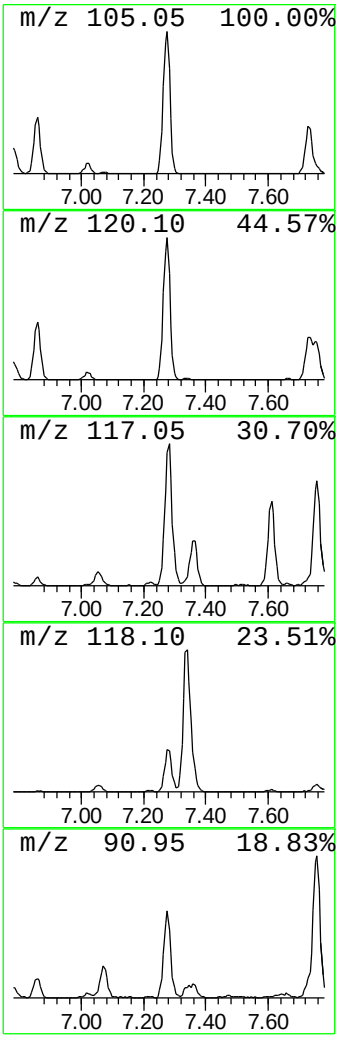
Peak Number 10 Benzene, 1-ethyl-3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.28	16.27 ng/ul	2120380	1,4-Dichlorobenzene-d4	7.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
3	Mesitylene	120	C9H12	000108-67-8	70
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	70
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	70





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Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

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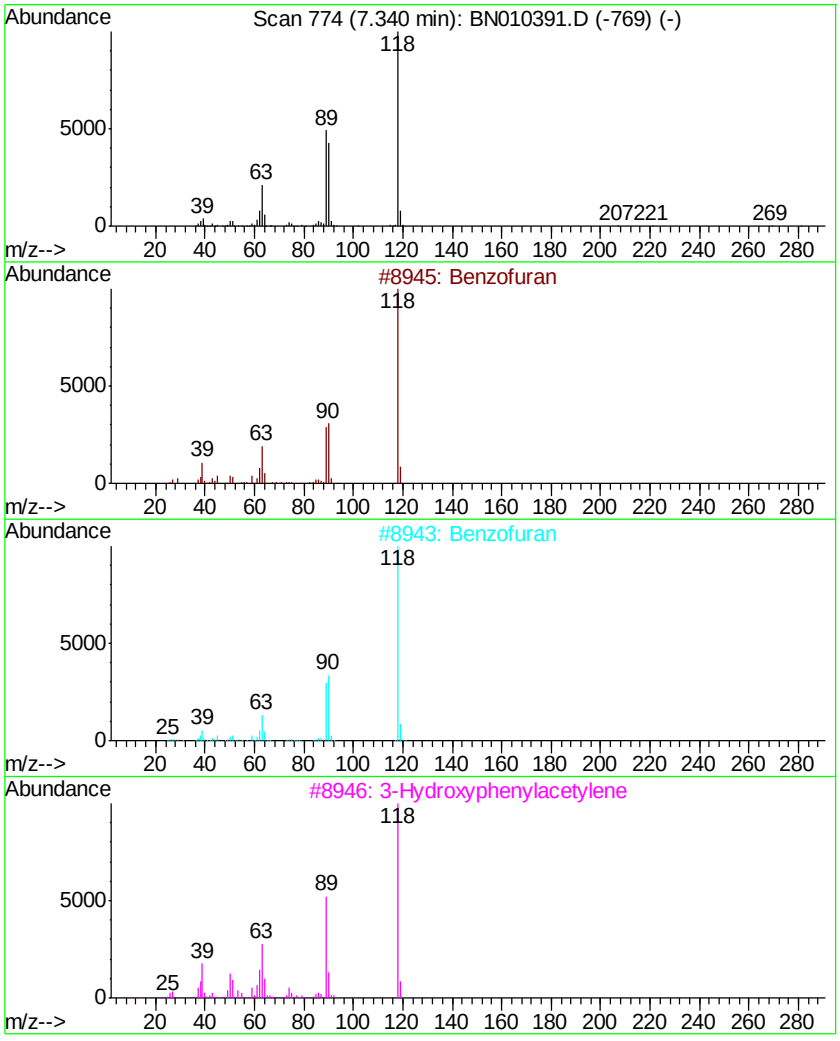
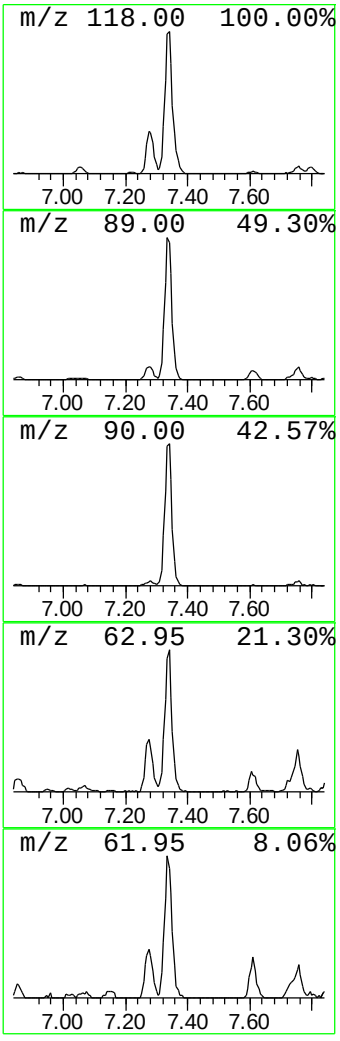
Peak Number 11 Benzofuran Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.34	9.69 ng/ul	1263050	1,4-Dichlorobenzene-d4	7.61

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	3-Hydroxyphenylacetylene	118	C8H6O	010401-11-3	64
4	2H-Cyclopenta[d]pyridazine	118	C7H6N2	000270-64-4	47
5	3-Pyridineacetonitrile	118	C7H6N2	006443-85-2	43





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Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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Quant Title : SVOA CALIBRATION

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

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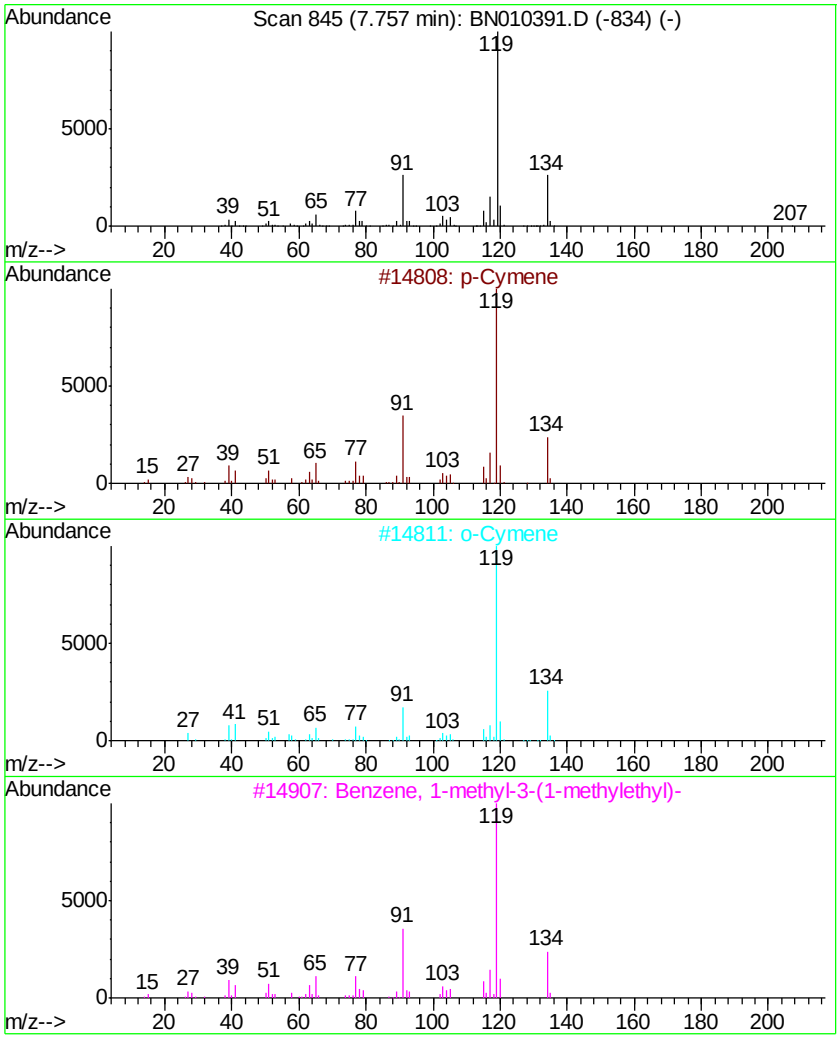
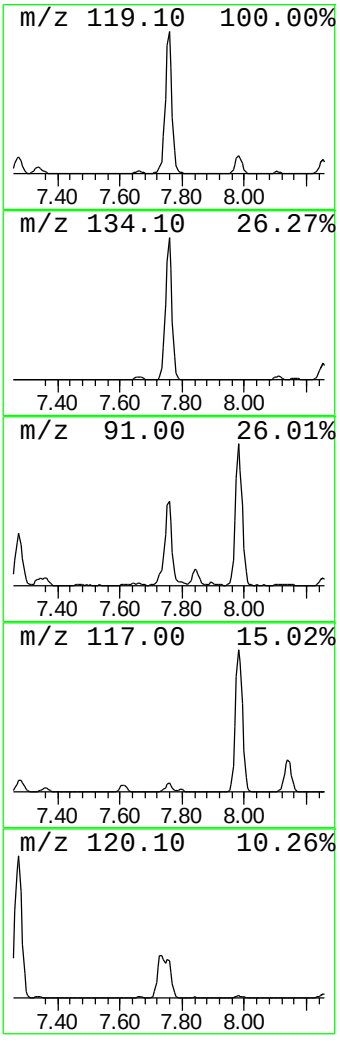
Peak Number 12 p-Cymene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.76	17.33 ng/ul	2258030	1,4-Dichlorobenzene-d4	7.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	95
2	o-Cymene	134	C10H14	000527-84-4	94
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
4	o-Cymene	134	C10H14	000527-84-4	94
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91





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Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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Quant Title : SVOA CALIBRATION

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

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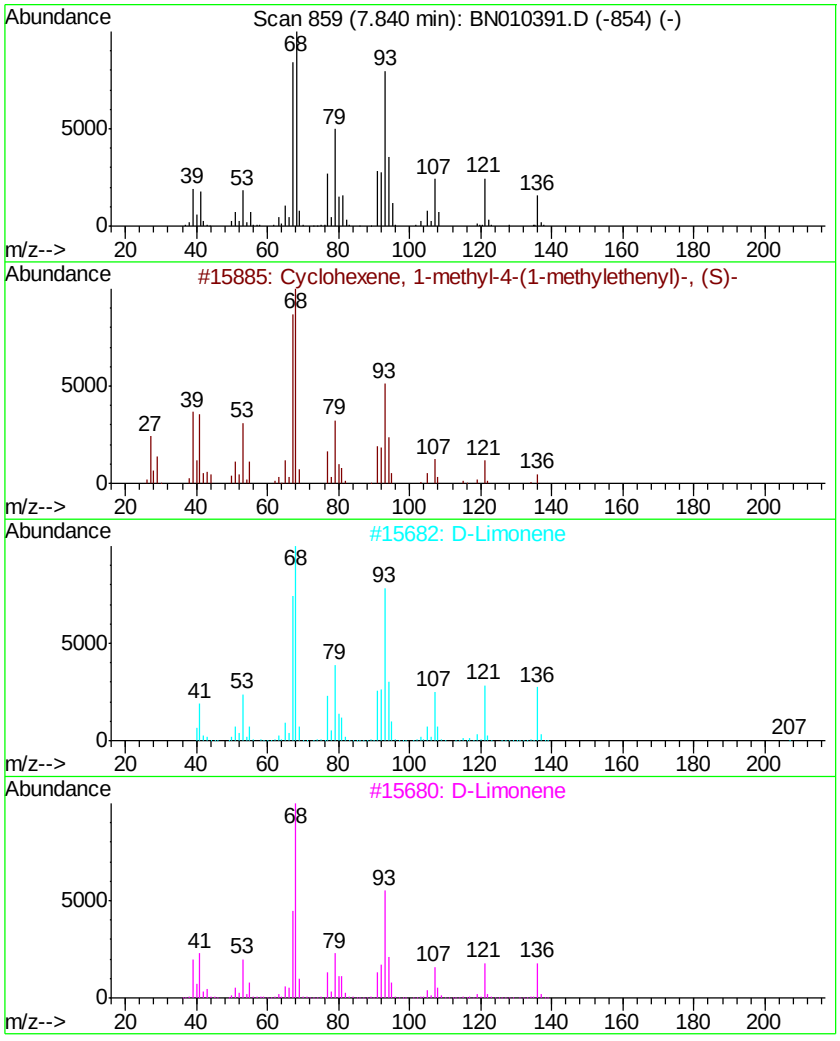
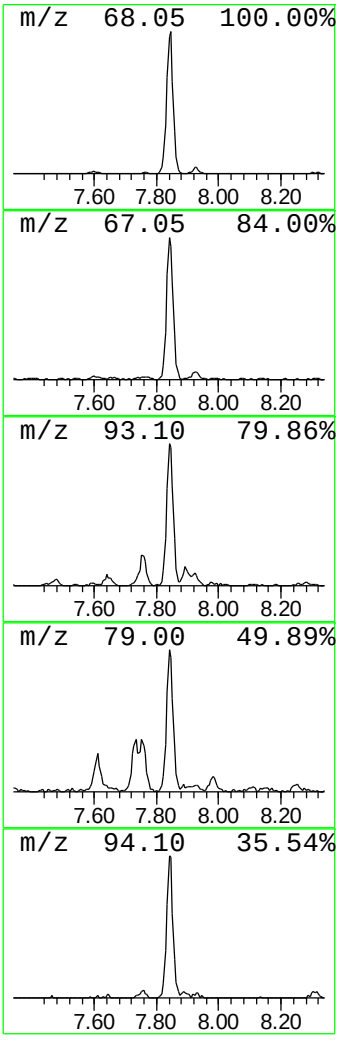
Peak Number 13 Cyclohexene, 1-methyl-4-(1-... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.84	6.16 ng/ul	802872	1,4-Dichlorobenzene-d4	7.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexene, 1-methyl-4-(1-methy...	136	C10H16	005989-54-8	97
2	D-Limonene	136	C10H16	005989-27-5	96
3	D-Limonene	136	C10H16	005989-27-5	94
4	D-Limonene	136	C10H16	005989-27-5	94
5	D-Limonene	136	C10H16	005989-27-5	93





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Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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Quant Title : SVOA CALIBRATION

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

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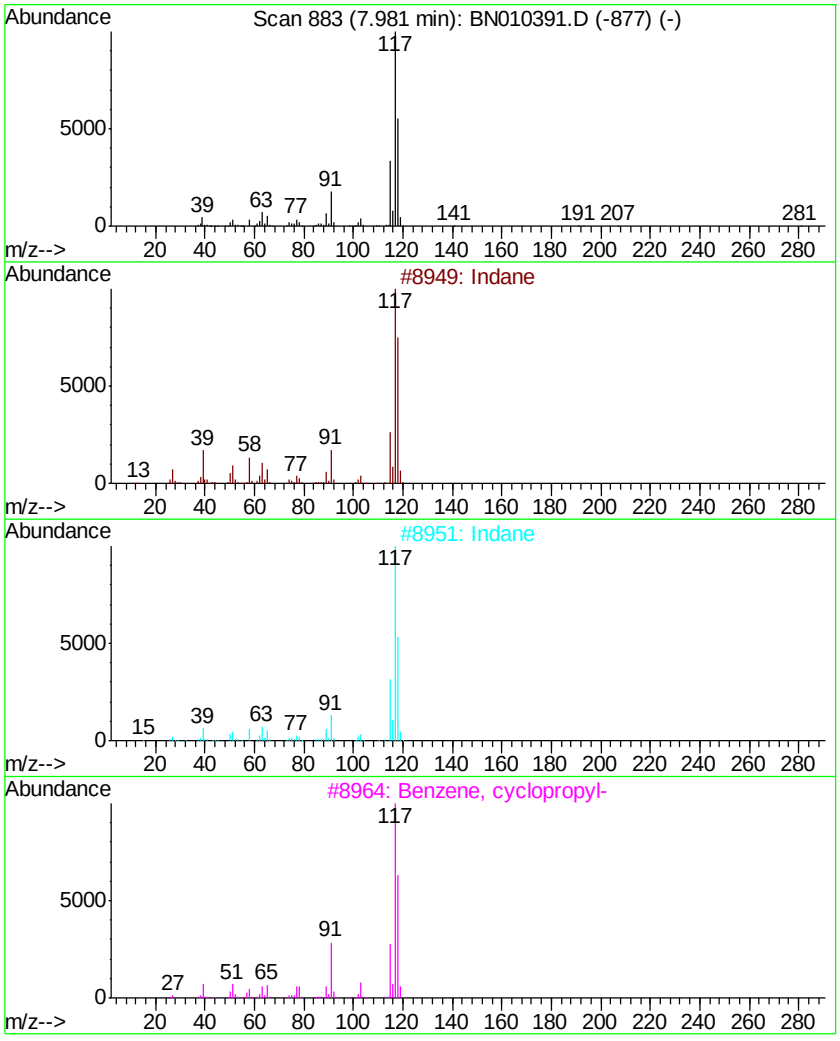
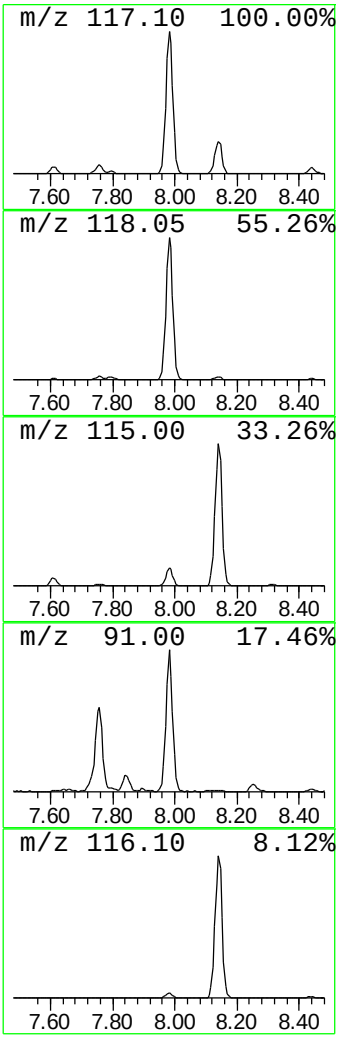
Peak Number 14 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.98	38.20 ng/ul	4979150	1,4-Dichlorobenzene-d4	7.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Indane	118	C9H10	000496-11-7	87
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
4	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	64
5	Indane	118	C9H10	000496-11-7	60





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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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Quant Title : SVOA CALIBRATION

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

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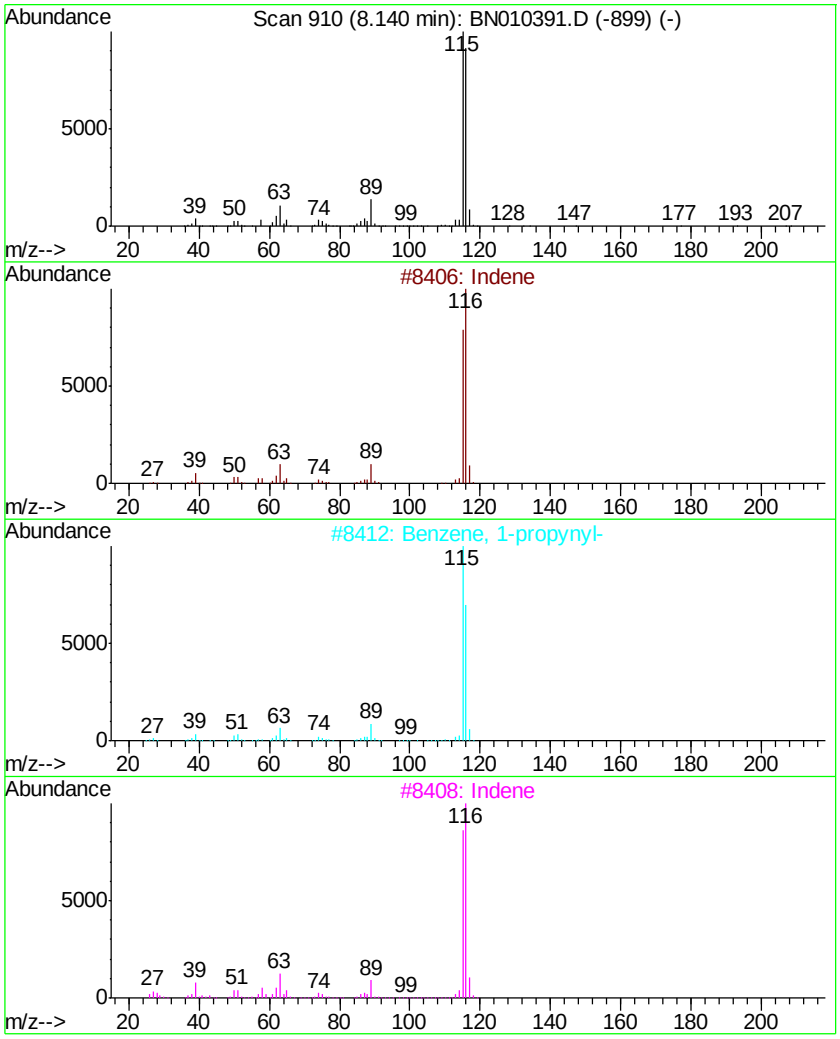
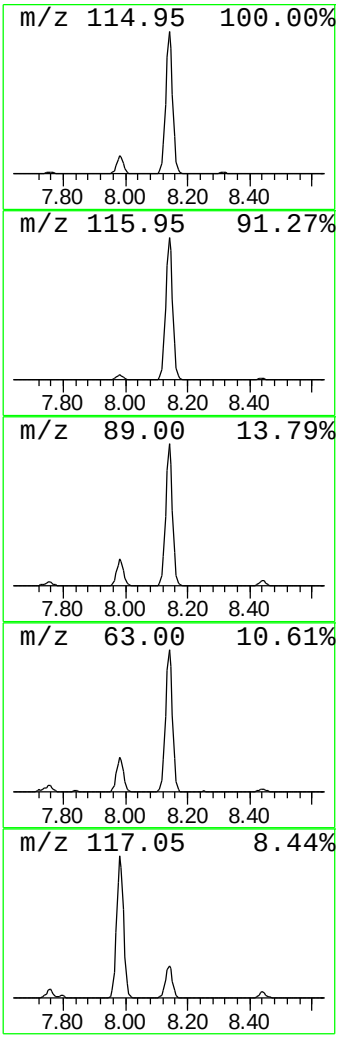
Peak Number 15 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.14	99.92 ng/ul	13022400	1,4-Dichlorobenzene-d4	7.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	97
2	Benzene, 1-propynyl-	116	C9H8	000673-32-5	95
3	Indene	116	C9H8	000095-13-6	95
4	3-Methylphenylacetylene	116	C9H8	000766-82-5	94
5	Benzene, 1-propynyl-	116	C9H8	000673-32-5	91





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Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

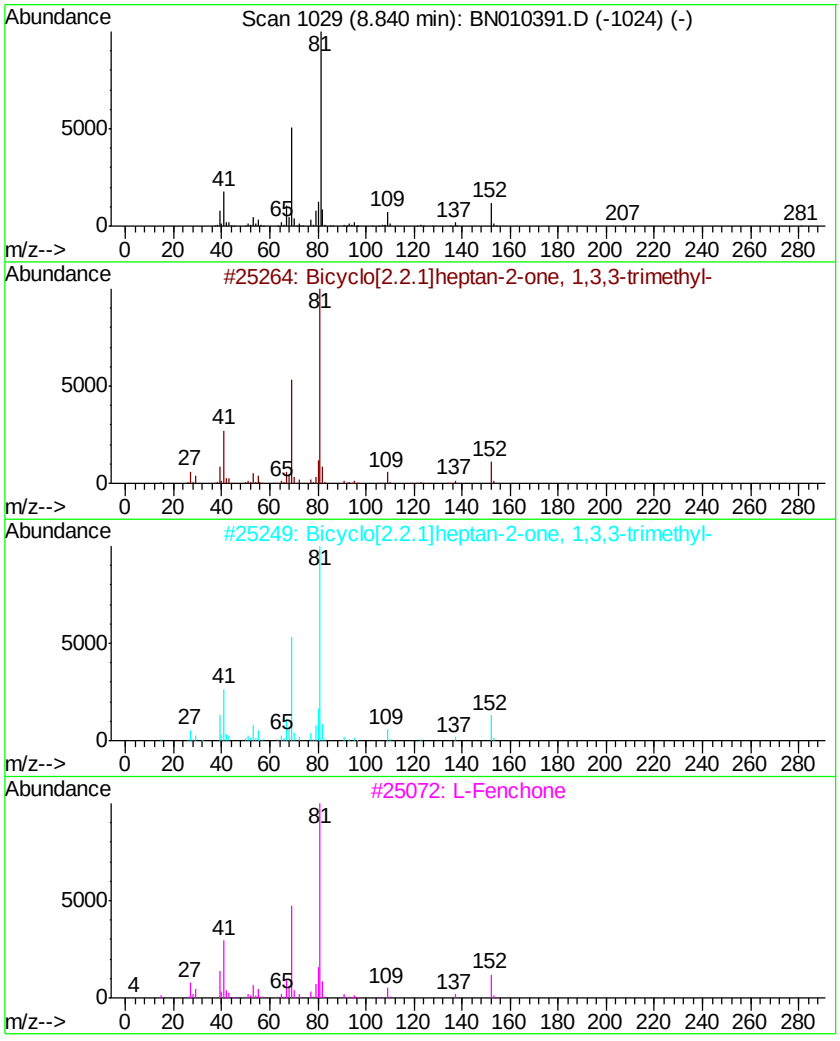
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

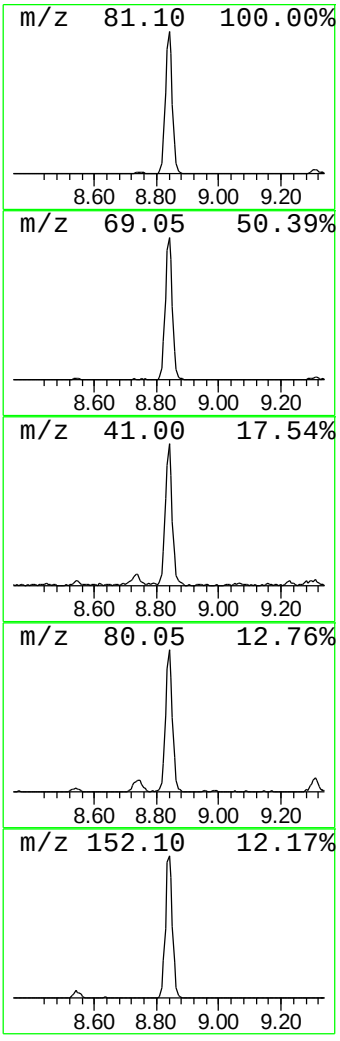
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.84	11.02 ng/ul	1436610	1,4-Dichlorobenzene-d4	7.61

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	94
2	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	94
3	L-Fenchone	152	C10H16O	007787-20-4	94
4	L-Fenchone	152	C10H16O	007787-20-4	91
5	Bicyclo[2.2.1]heptan-2-one, 1,3,...	152	C10H16O	001195-79-5	91





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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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Quant Title : SVOA CALIBRATION

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

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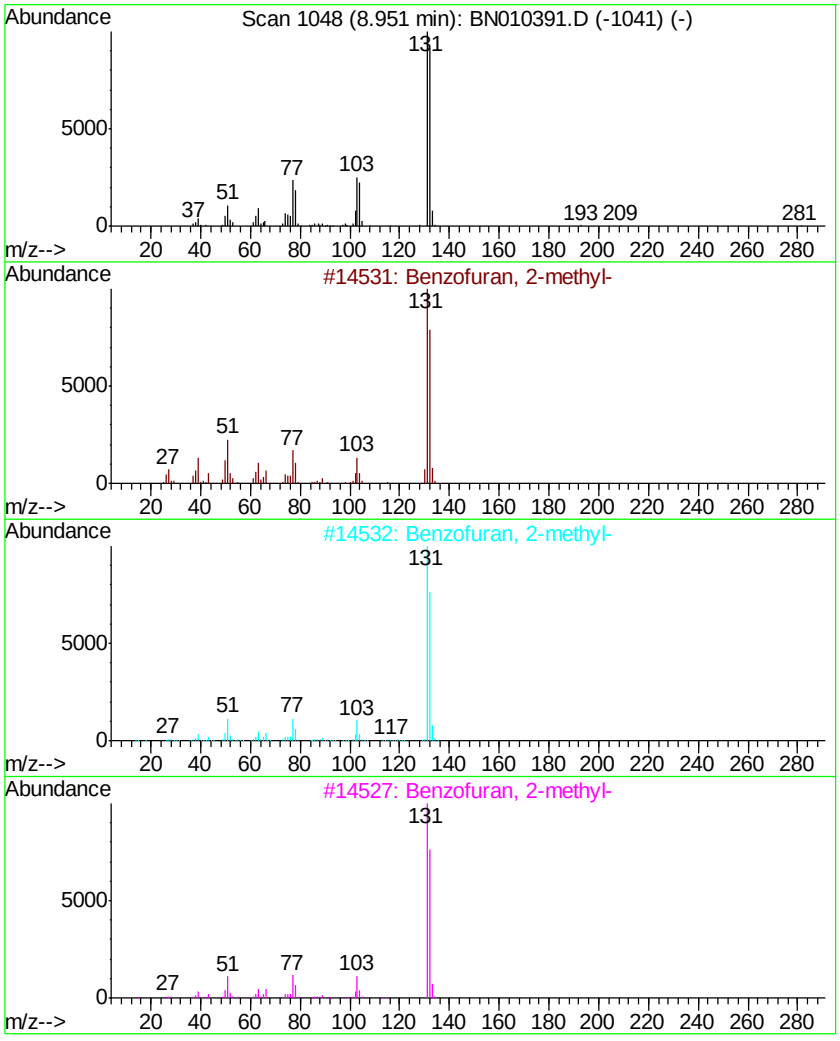
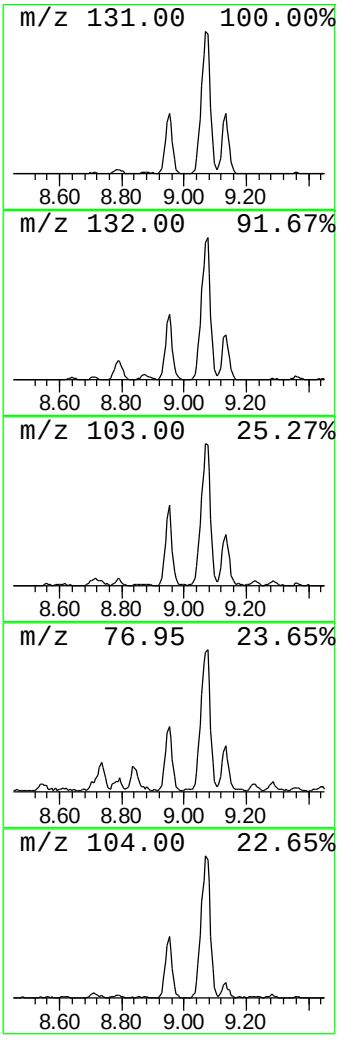
Peak Number 17 Benzofuran, 2-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.95	5.80 ng/ul	756418	1,4-Dichlorobenzene-d4	7.61

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, 2-methyl-	132	C9H8O	004265-25-2	91
4	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	90
5	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	87





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Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

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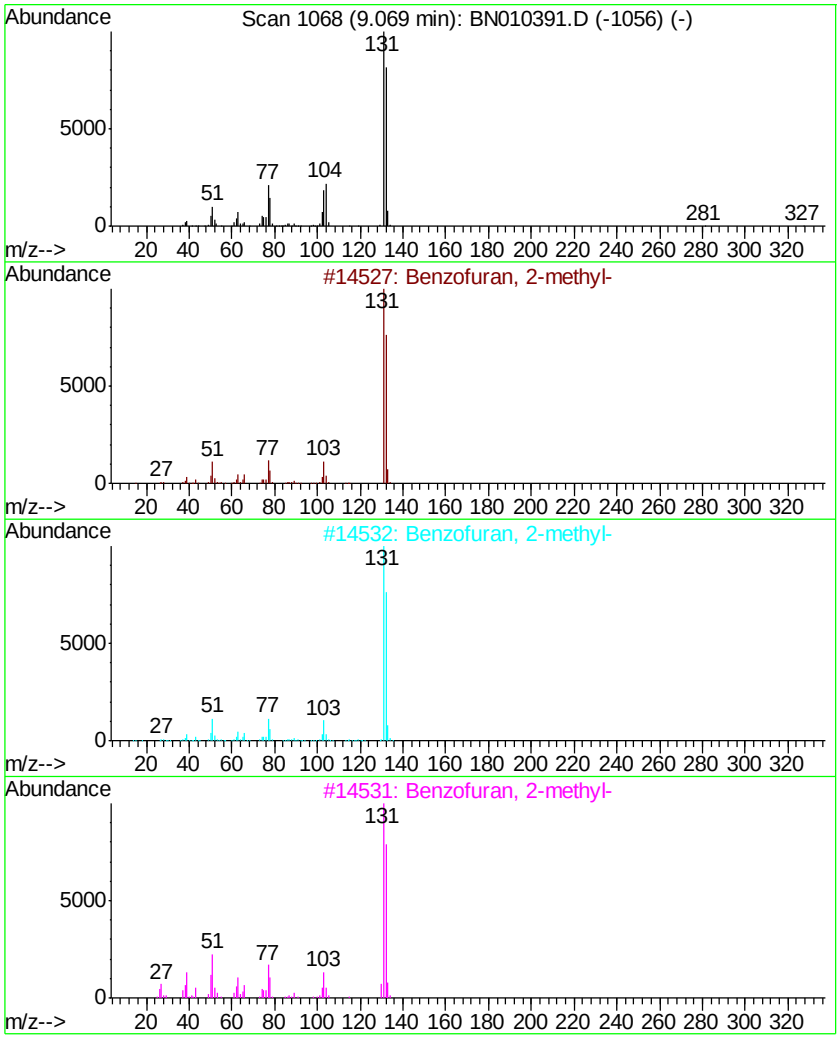
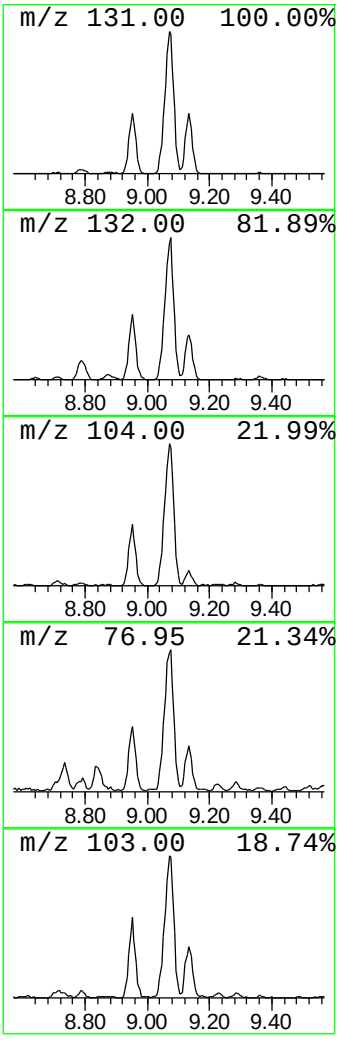
Peak Number 18 4-Aminobenzyl cyanide Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.07	15.03 ng/ul	2142050	Naphthalene-d8	10.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
4	4-Aminobenzyl cyanide	132	C8H8N2	003544-25-0	87
5	4-Aminobenzyl cyanide	132	C8H8N2	003544-25-0	86





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Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

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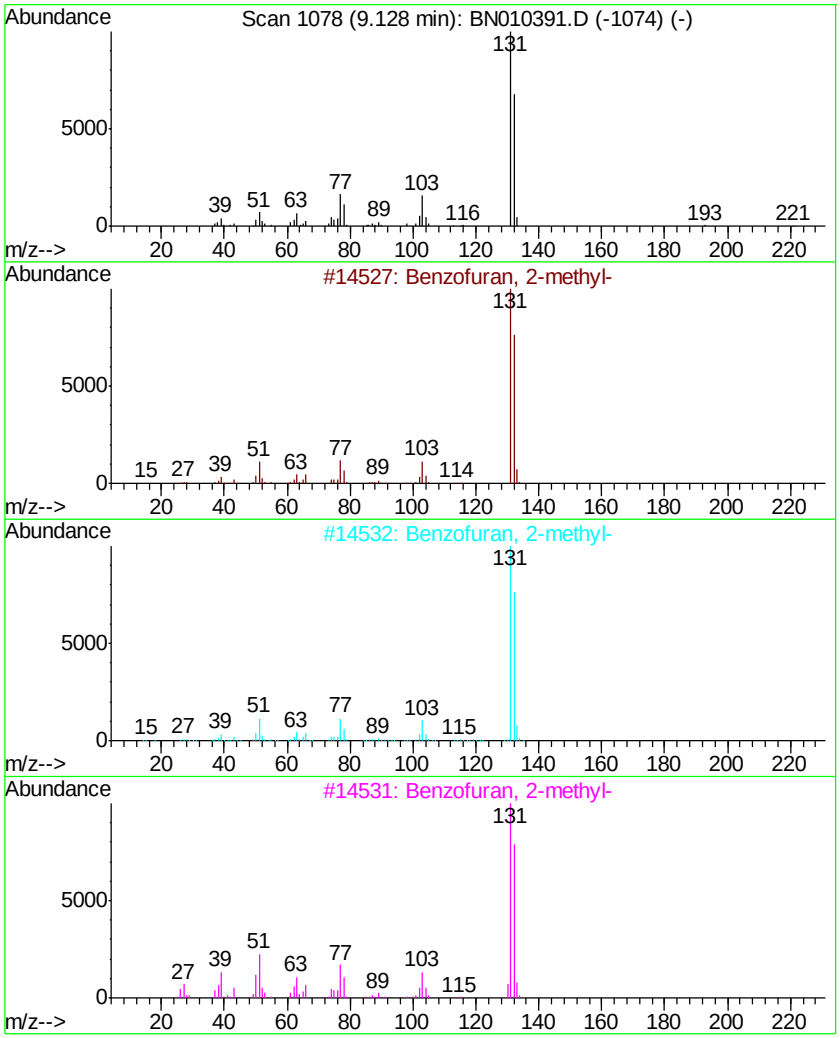
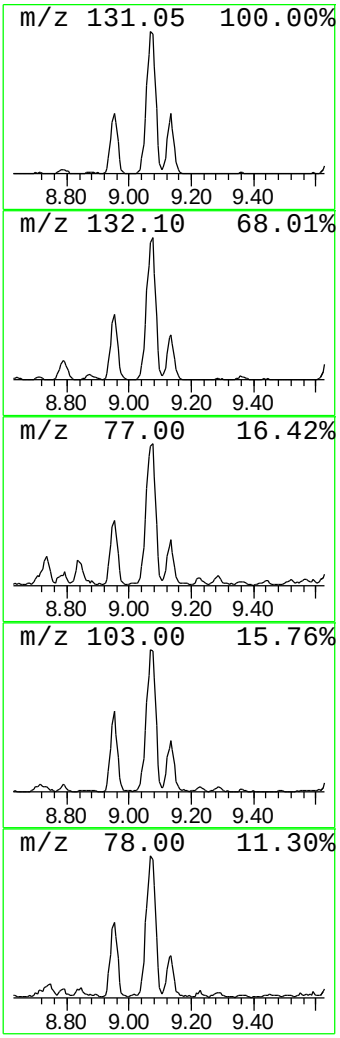
Peak Number 19 Benzofuran, 7-methyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.13	3.75 ng/ul	534360	Naphthalene-d8	10.38

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	91
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	64
4	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	64
5	Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	59





Library Search Compound Report

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Data File : BN010391.D
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Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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Quant Title : SVOA CALIBRATION

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

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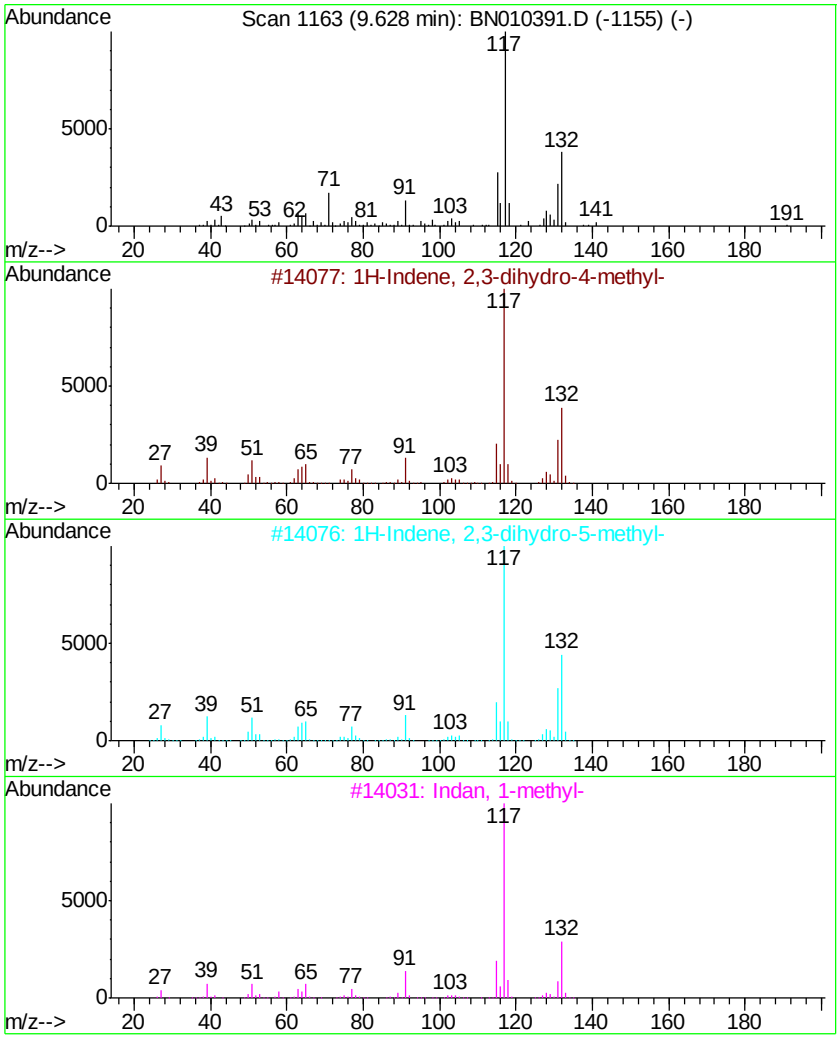
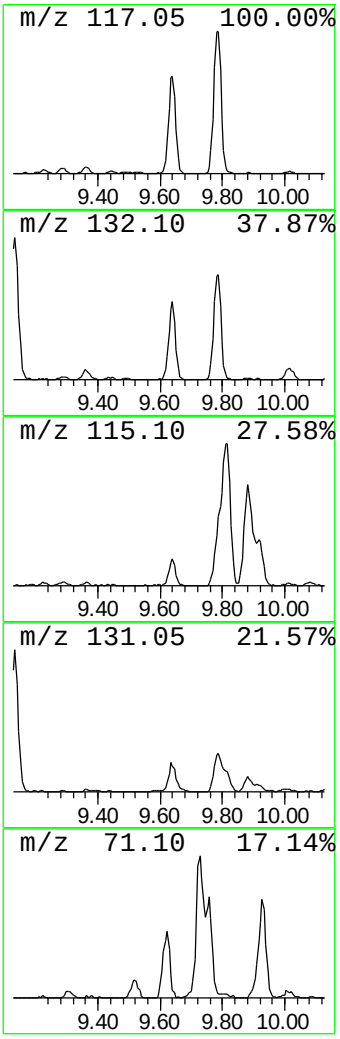
Peak Number 20 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.63	4.71 ng/ul	670785	Naphthalene-d8	10.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
3	Indan, 1-methyl-	132	C10H12	000767-58-8	87
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5	Indan, 1-methyl-	132	C10H12	000767-58-8	83





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Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

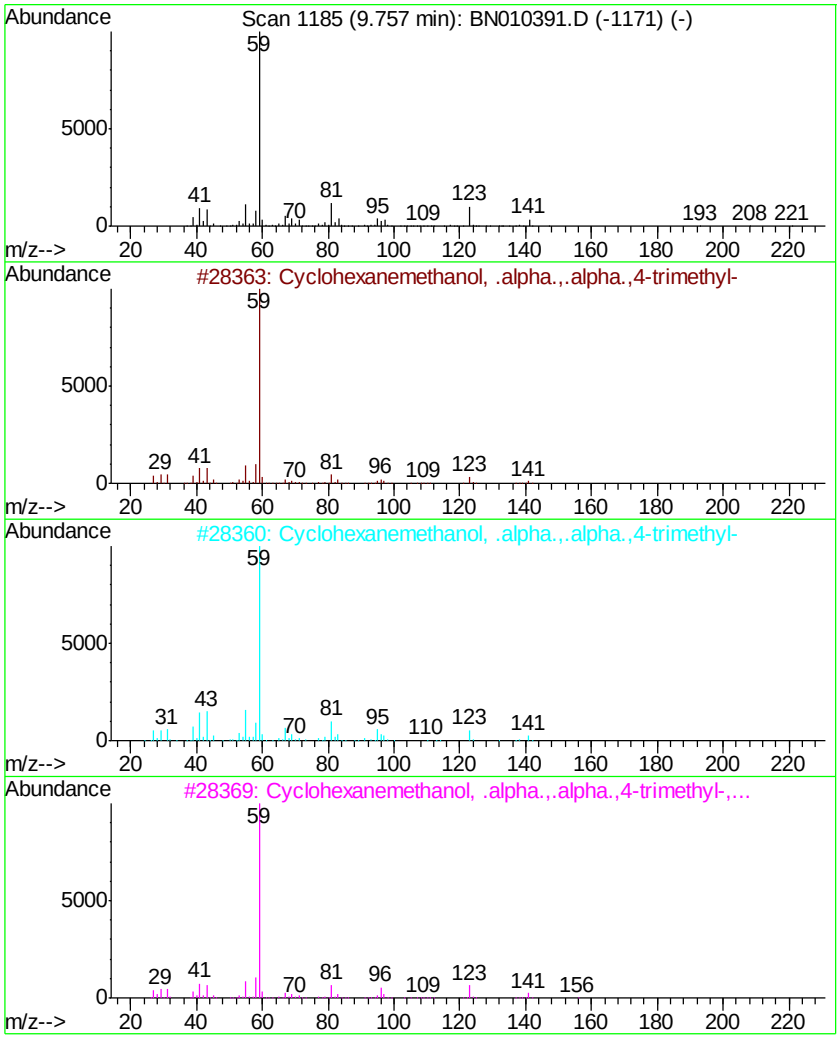
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

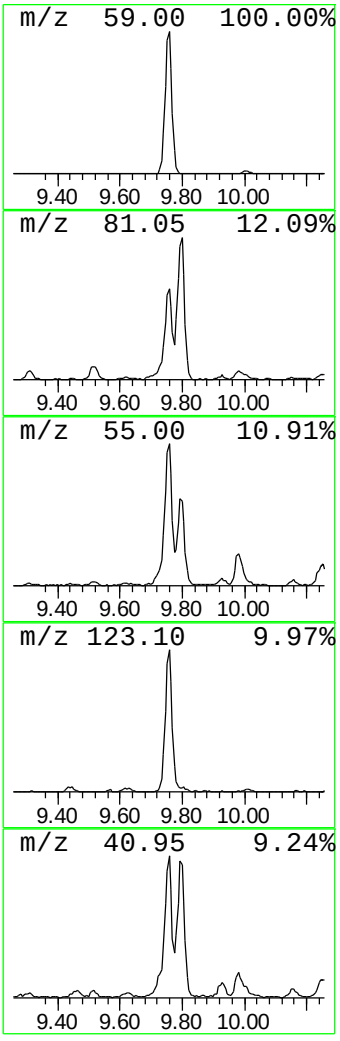
****

Peak Number 21 Cyclohexanemethanol, .alpha... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.76	23.95 ng/ul	3412280	Naphthalene-d8	10.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	74
2	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	000498-81-7	64
3	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	005114-00-1	64
4	Cyclohexanemethanol, .alpha.,.al...	156	C10H20O	007322-63-6	56
5	2,4-Dimethyl-4-penten-2-ol	114	C7H14O	019781-53-4	53





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Data File : BN010391.D
Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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Quant Title : SVOA CALIBRATION

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

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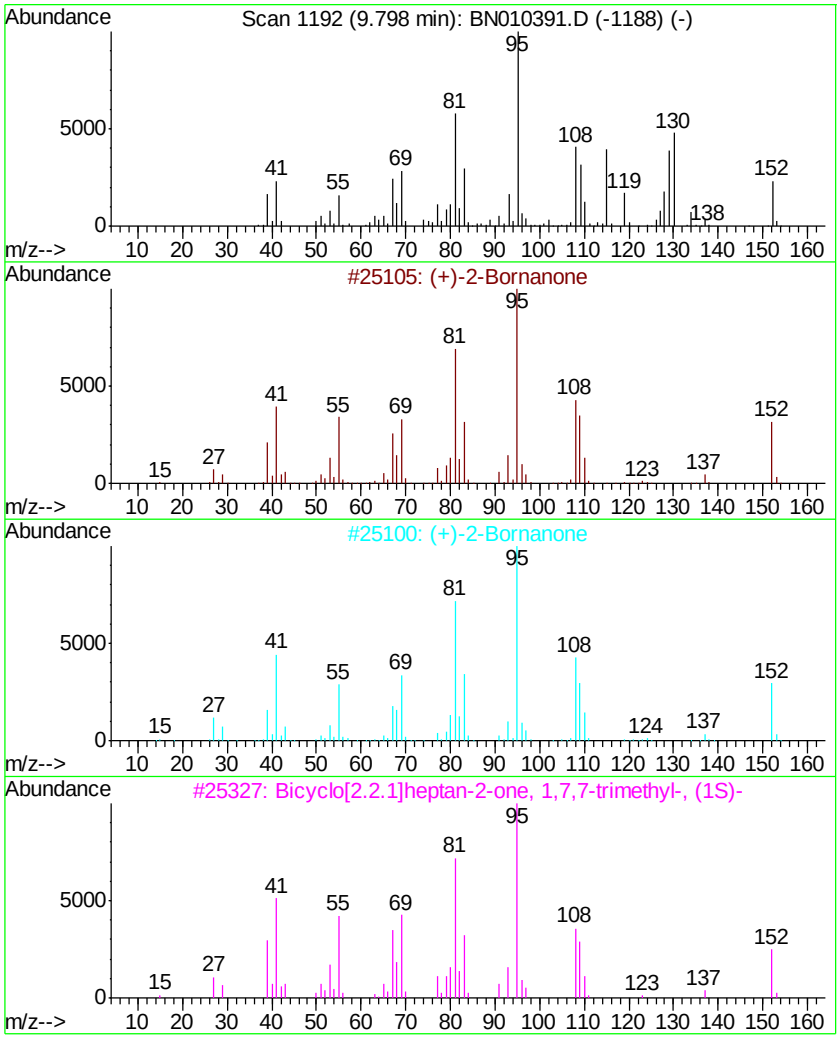
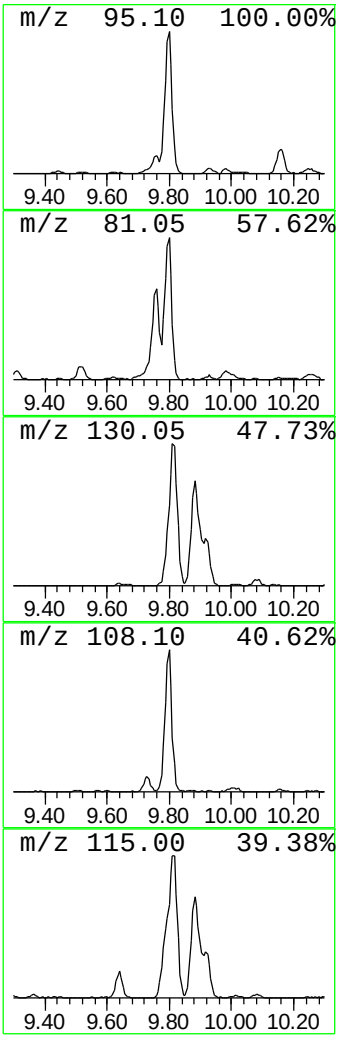
Peak Number 22 (+)-2-Bornanone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.80	29.82 ng/ul	4248290	Naphthalene-d8	10.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-2-Bornanone	152	C10H16O	000464-49-3	95
2	(+)-2-Bornanone	152	C10H16O	000464-49-3	95
3	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	94
4	Camphor	152	C10H16O	000076-22-2	90
5	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	90





Library Search Compound Report

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Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

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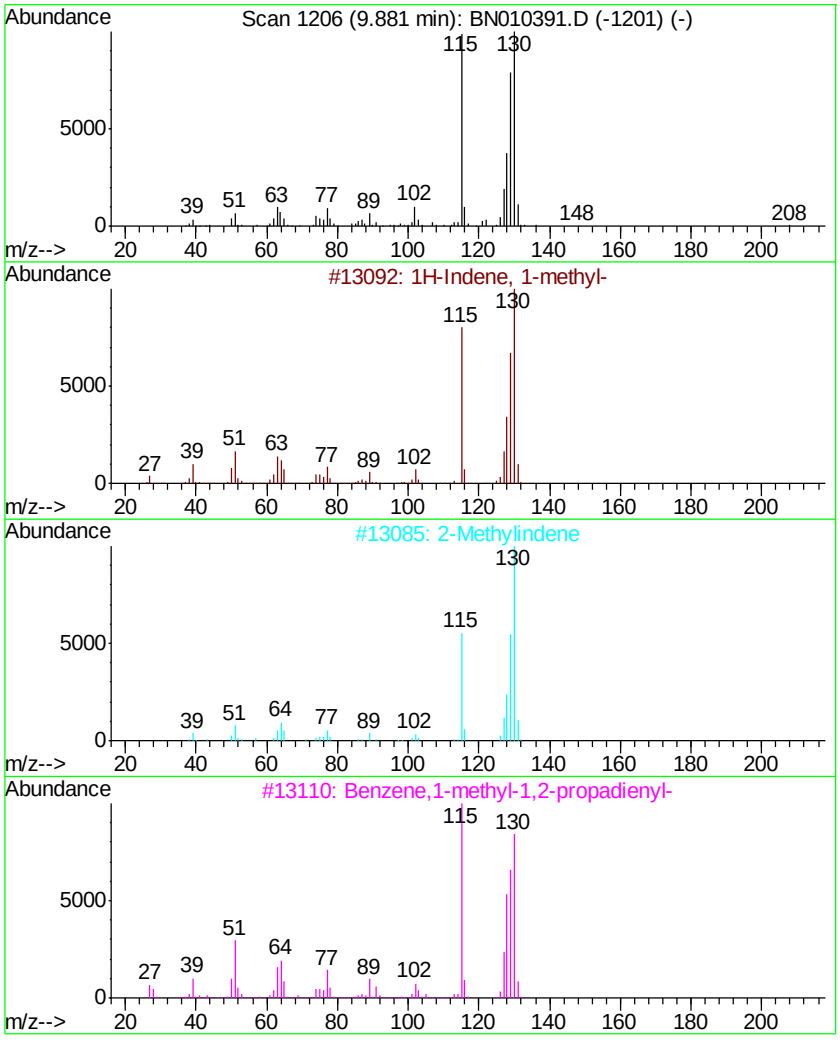
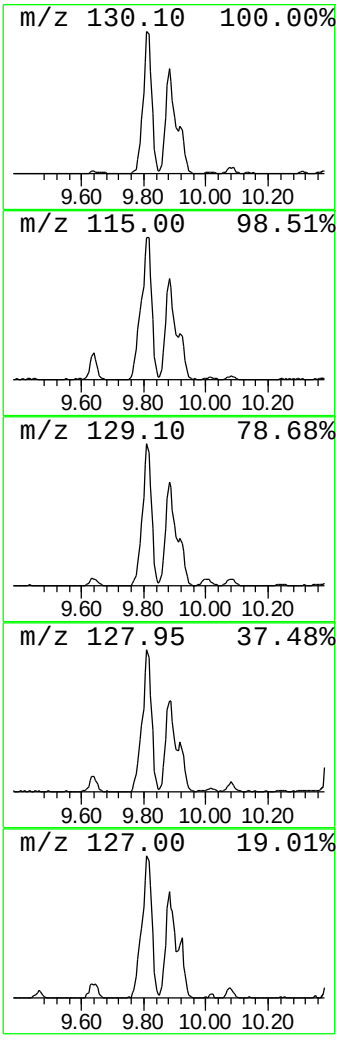
Peak Number 23 1H-Indene, 1-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.88	7.90 ng/ul	1125640	Naphthalene-d8	10.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	95
2	2-Methylindene	130	C10H10	002177-47-1	95
3	Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	94
4	Benzene, 1-butynyl-	130	C10H10	000622-76-4	93
5	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	93





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Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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Quant Title : SVOA CALIBRATION

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

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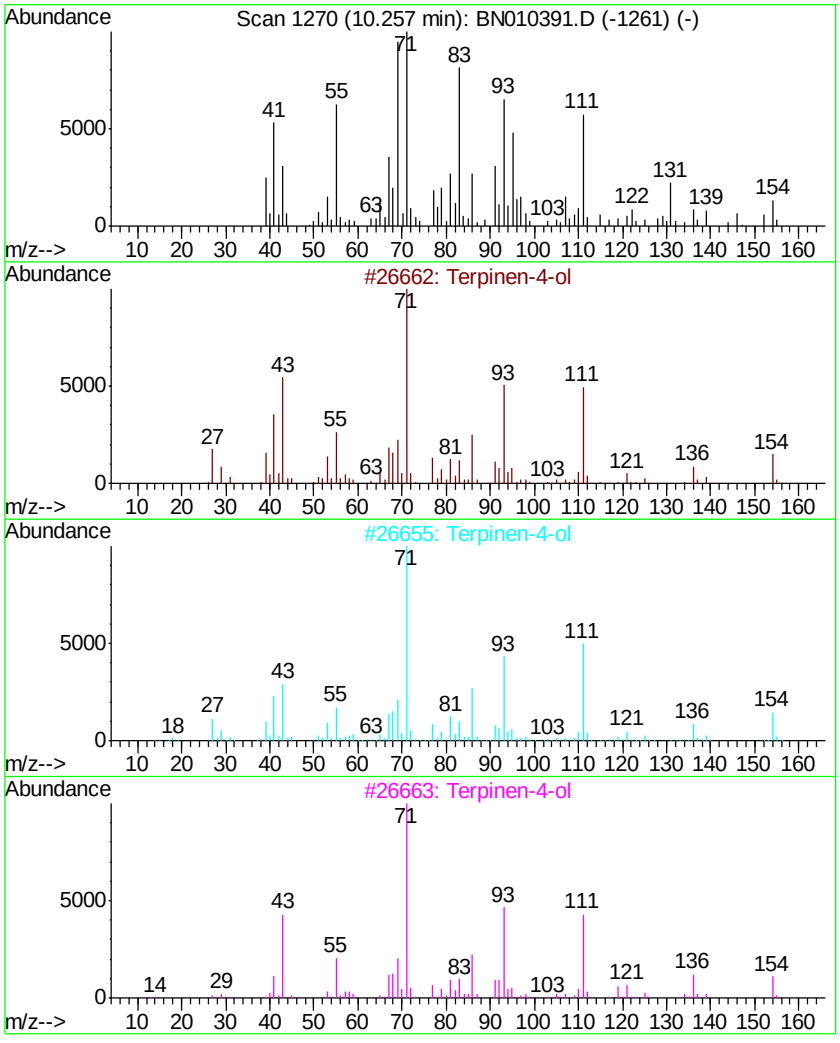
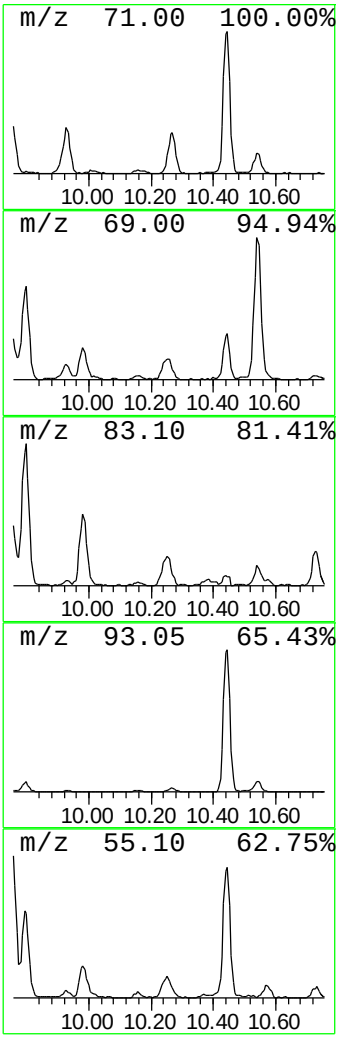
Peak Number 25 Terpinen-4-ol Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.26	3.30 ng/ul	470102	Naphthalene-d8	10.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Terpinen-4-ol	154	C10H18O	000562-74-3	86
2	Terpinen-4-ol	154	C10H18O	000562-74-3	50
3	Terpinen-4-ol	154	C10H18O	000562-74-3	45
4	Terpinen-4-ol	154	C10H18O	000562-74-3	38
5	3-Cyclohexen-1-ol, 4-methyl-1-(1...	154	C10H18O	020126-76-5	38





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Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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Quant Title : SVOA CALIBRATION

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

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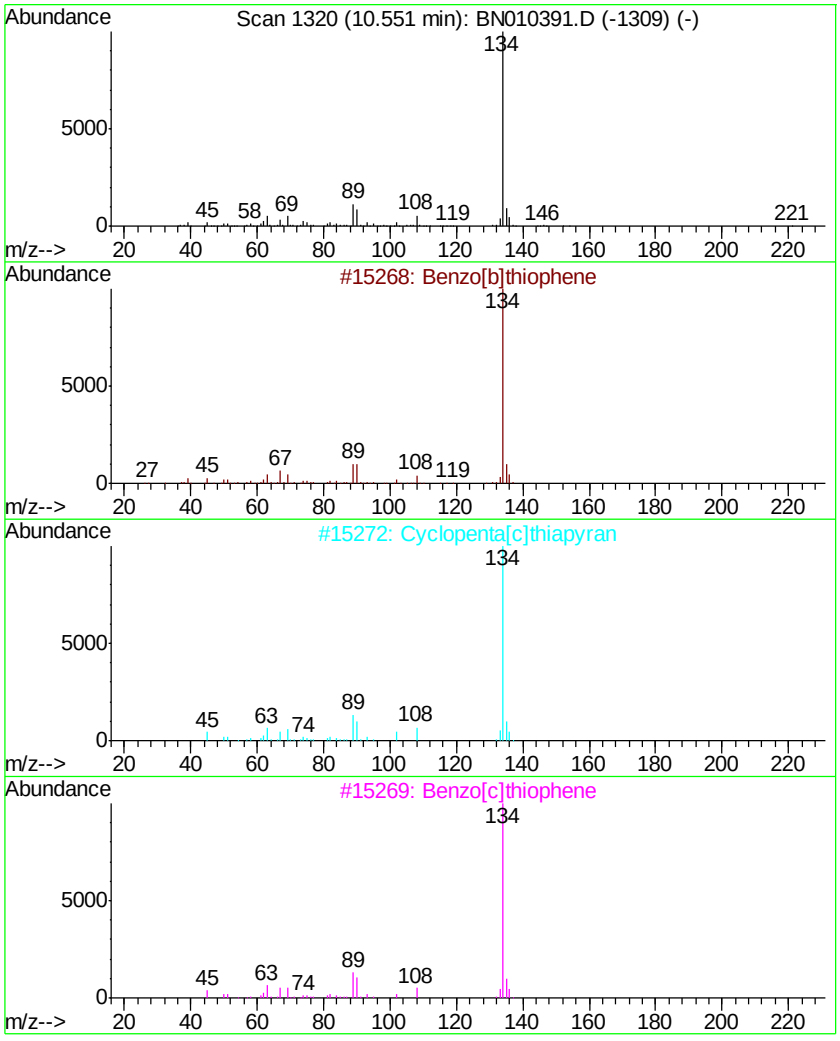
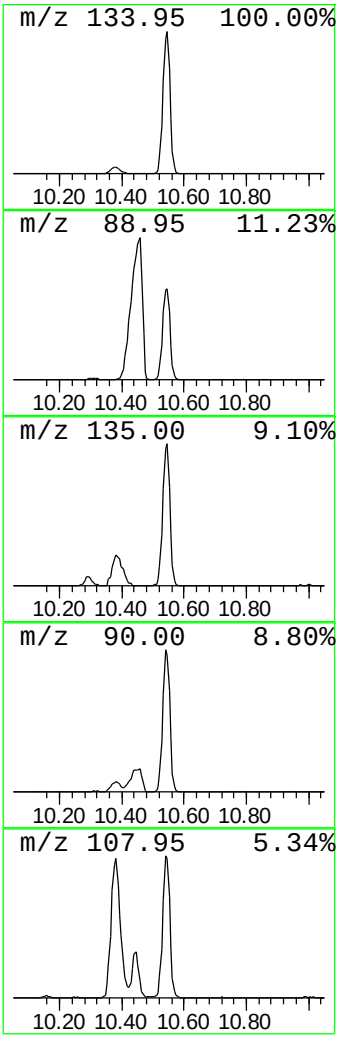
Peak Number 26 Benzo[b]thiophene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.55	44.56 ng/ul	6349280	Naphthalene-d8	10.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene	134	C8H6S	000095-15-8	94
2	Cyclopenta[c]thiapyran	134	C8H6S	000270-63-3	94
3	Benzo[c]thiophene	134	C8H6S	000270-82-6	94
4	Benzo[b]thiophene	134	C8H6S	000095-15-8	91
5	Benzo[b]thiophene	134	C8H6S	000095-15-8	86





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Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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Quant Title : SVOA CALIBRATION

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

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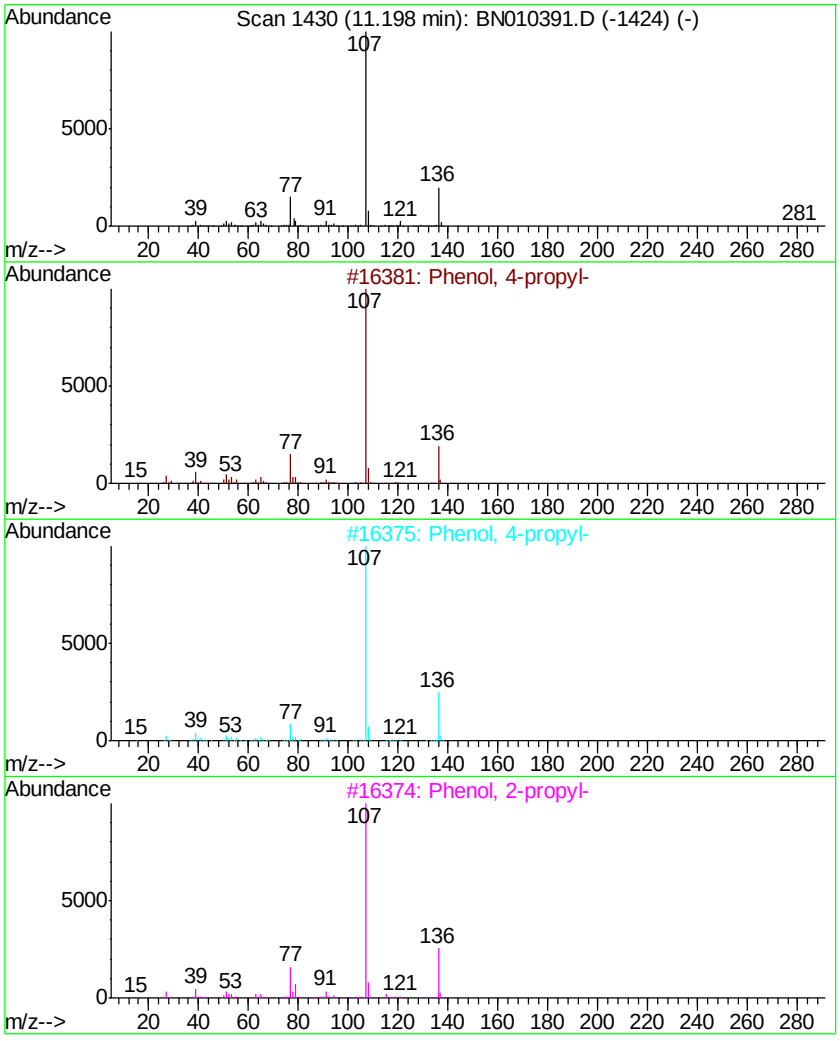
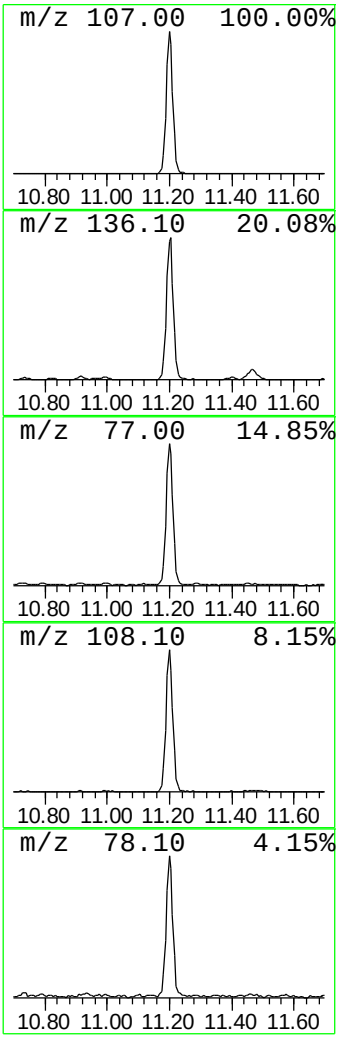
Peak Number 27 Phenol, 4-propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.20	22.48 ng/ul	3203410	Naphthalene-d8	10.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-propyl-	136	C9H12O	000645-56-7	91
2	Phenol, 4-propyl-	136	C9H12O	000645-56-7	80
3	Phenol, 2-propyl-	136	C9H12O	000644-35-9	64
4	Phenol, 2-propyl-	136	C9H12O	000644-35-9	52
5	Phenol, 2-ethyl-	122	C8H10O	000090-00-6	50





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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

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Quant Title : SVOA CALIBRATION

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

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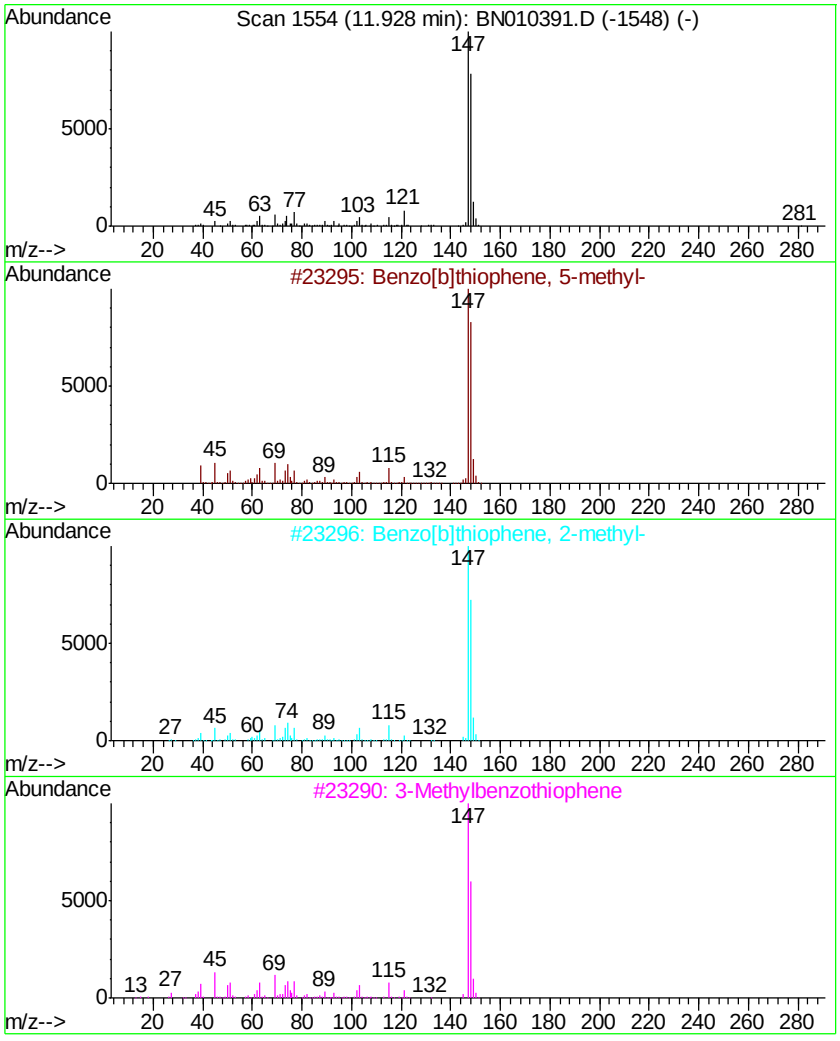
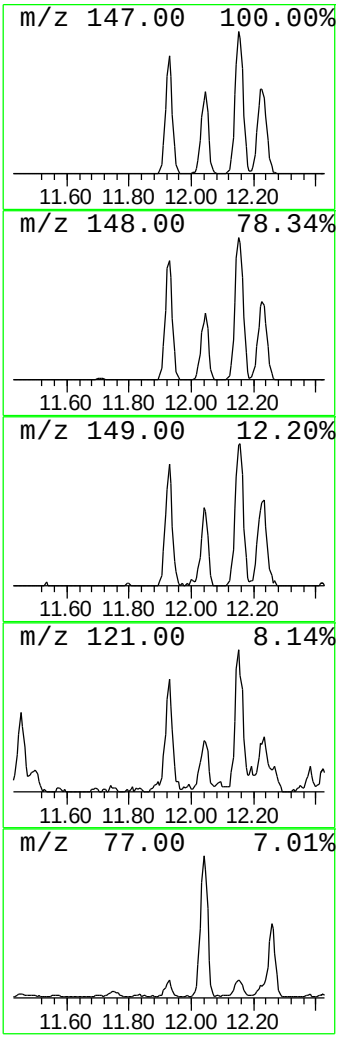
Peak Number 28 Benzo[b]thiophene, 5-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	5.80 ng/ul	825785	Naphthalene-d8	10.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	91
2	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
3	3-Methylbenzothiophene	148	C9H8S	001455-18-1	91
4	3-Methylbenzothiophene	148	C9H8S	001455-18-1	87
5	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	86





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Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

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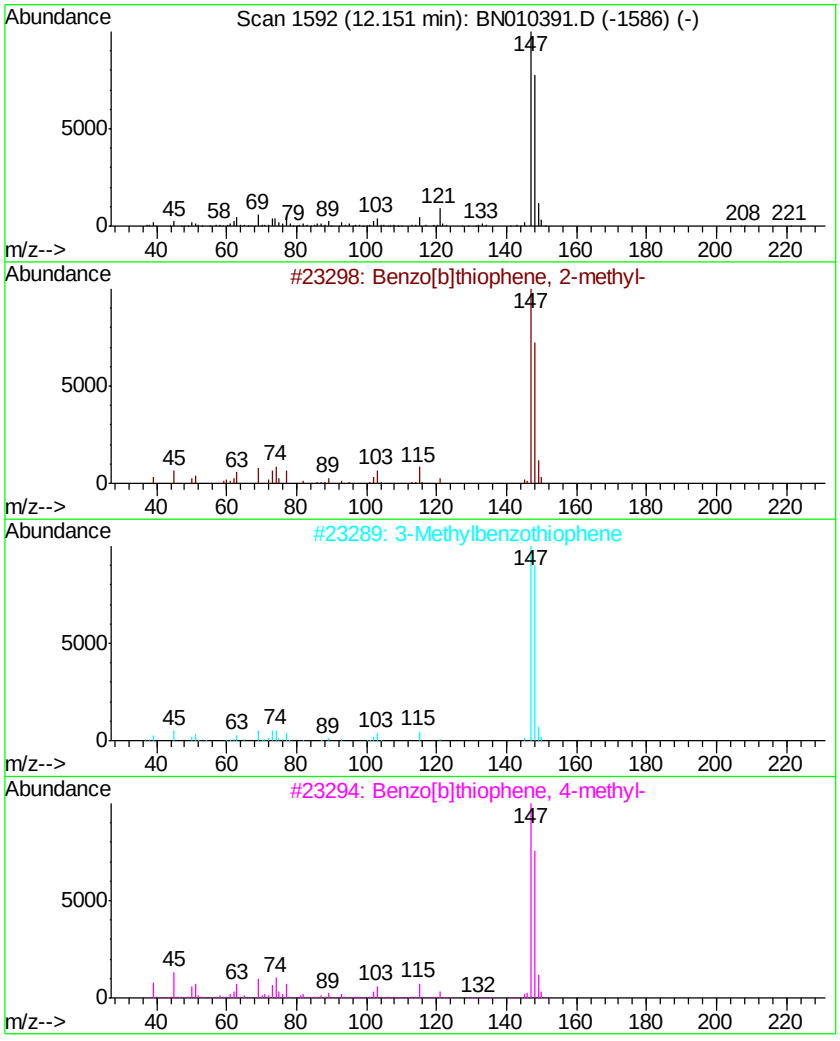
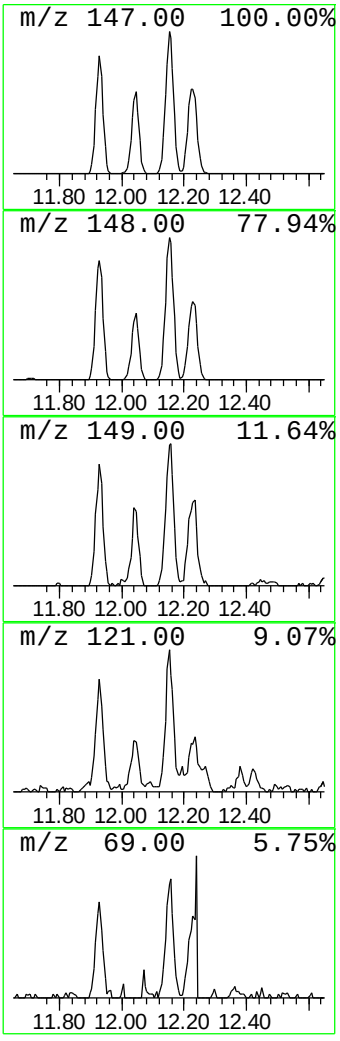
Peak Number 29 Benzo[b]thiophene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	8.19 ng/ul	1166510	Naphthalene-d8	10.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	90
2	3-Methylbenzothiophene	148	C9H8S	001455-18-1	90
3	Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	90
4	Benzo[b]thiophene, 5-methyl-	148	C9H8S	014315-14-1	90
5	Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	90





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Data File : BN010391.D
Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

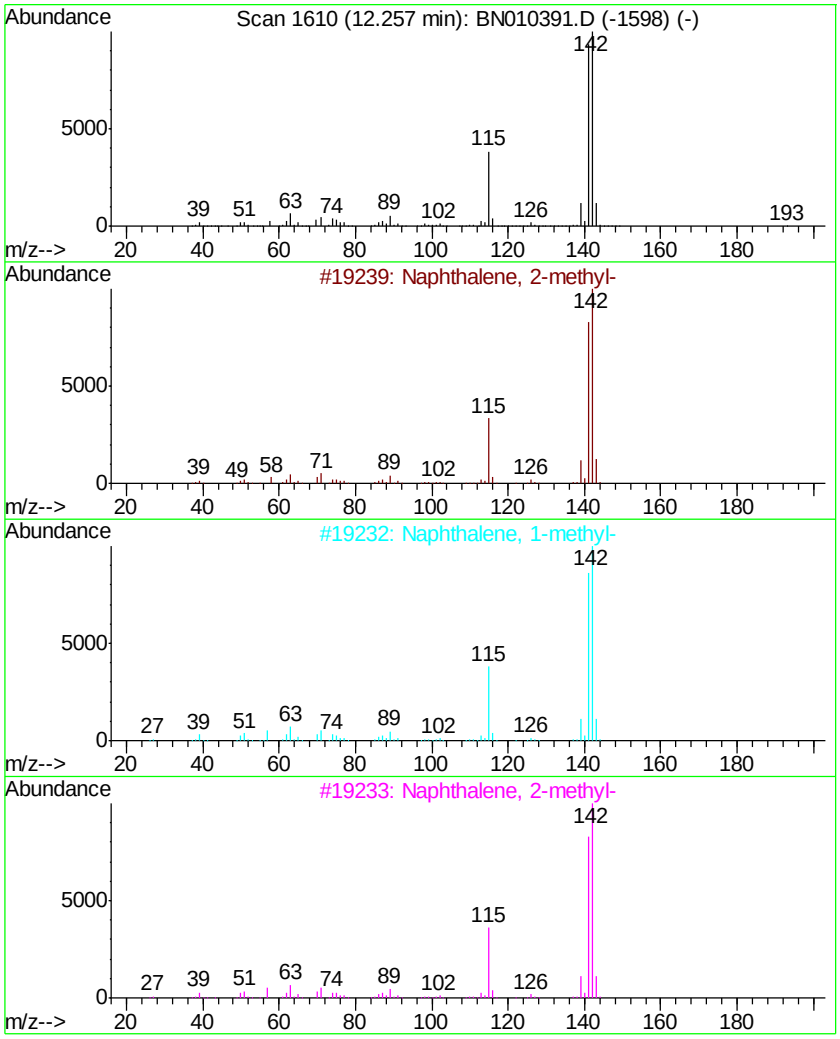
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

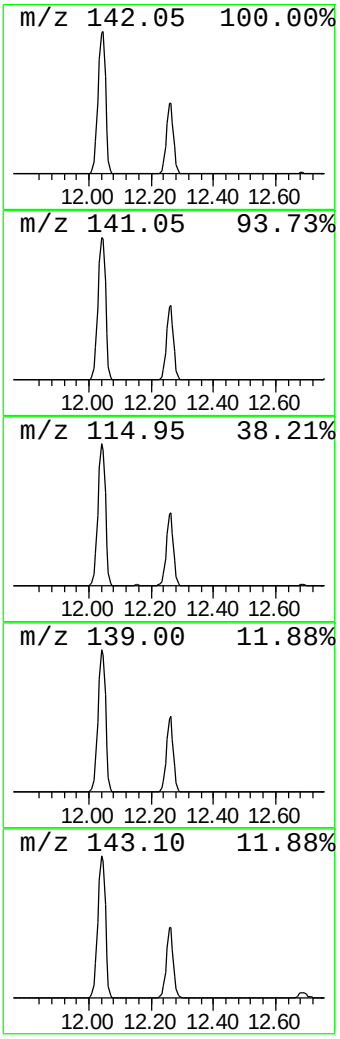
****

Peak Number 30 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	119.37 ng/ul	17007100	Naphthalene-d8	10.38

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	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
5	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN040120\
Data File : BN010391.D
Acq On : 02 Apr 2020 14:21
Operator : CG/JU
Sample : L2065-11DL 10X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : DBF07DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN040120MA.M
Quant Title : SVOA CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1,3-dime...	5.31	4.5	ng/ul	589512	1	7.61	2606660	20.0
Bicyclo[4.2.0]oct...	5.65	2.0	ng/ul	267606	1	7.61	2606660	20.0
(1R)-2,6,6-Trimet...	6.33	12.8	ng/ul	1663630	1	7.61	2606660	20.0
Camphene	6.63	2.9	ng/ul	383488	1	7.61	2606660	20.0
Benzene, 1-ethyl-...	6.73	4.6	ng/ul	598815	1	7.61	2606660	20.0
Benzene, 1-ethyl-...	6.78	2.7	ng/ul	354463	1	7.61	2606660	20.0
Mesitylene	6.86	5.0	ng/ul	656982	1	7.61	2606660	20.0
.beta.-Pinene	7.07	5.2	ng/ul	670834	1	7.61	2606660	20.0
Benzene, 1-ethyl-...	7.28	16.3	ng/ul	2120380	1	7.61	2606660	20.0
Benzofuran	7.34	9.7	ng/ul	1263050	1	7.61	2606660	20.0
p-Cymene	7.76	17.3	ng/ul	2258030	1	7.61	2606660	20.0
Cyclohexene, 1-me...	7.84	6.2	ng/ul	802872	1	7.61	2606660	20.0
Indane	7.98	38.2	ng/ul	4979150	1	7.61	2606660	20.0
Indene	8.14	99.9	ng/ul	13022400	1	7.61	2606660	20.0
Bicyclo[2.2.1]hep...	8.84	11.0	ng/ul	1436610	1	7.61	2606660	20.0
Benzofuran, 2-met...	8.95	5.8	ng/ul	756418	1	7.61	2606660	20.0
4-Aminobenzyl cya...	9.07	15.0	ng/ul	2142050	2	10.38	2849580	20.0
Benzofuran, 7-met...	9.13	3.8	ng/ul	534360	2	10.38	2849580	20.0
1H-Indene, 2,3-di...	9.63	4.7	ng/ul	670785	2	10.38	2849580	20.0
Cyclohexanemethan...	9.76	23.9	ng/ul	3412280	2	10.38	2849580	20.0
(+)-2-Bornanone	9.80	29.8	ng/ul	4248290	2	10.38	2849580	20.0
1H-Indene, 1-methyl-	9.88	7.9	ng/ul	1125640	2	10.38	2849580	20.0
Terpinen-4-ol	10.26	3.3	ng/ul	470102	2	10.38	2849580	20.0
Benzo[b]thiophene	10.55	44.6	ng/ul	6349280	2	10.38	2849580	20.0
Phenol, 4-propyl-	11.20	22.5	ng/ul	3203410	2	10.38	2849580	20.0
Benzo[b]thiophene...	11.93	5.8	ng/ul	825785	2	10.38	2849580	20.0
Benzo[b]thiophene...	12.15	8.2	ng/ul	1166510	2	10.38	2849580	20.0
Naphthalene, 1-me...	12.26	119.4	ng/ul	17007100	2	10.38	2849580	20.0