LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CROSSWICKS-DRUM-2

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.098	507	512	517	rVB	156198	177287	1.68%	0.204%
2	5.363	552	557	561	rBV	2991497	3050295	28.95%	3.505%
3	5.487	570	578	579	rBV	7091564	10536875	100.00%	12.108%
4	5.504	579	581	585	rVB	7026421	6522941	61.91%	7.495%
5	5.740	615	621	624	rBV	6570625	7522346	71.39%	8.644%
6	6.340	717	723	727	rBV	624757	515918	4.90%	0.593%
7	6.410	730	735	737	rBV	4010629	3622653	34.38%	4.163%
8	6.439	737	740	743	rVB	2336666	1849601	17.55%	2.125%
9	6.487	743	748	750	rBV	2920639	2720428	25.82%	3.126%
10	6.528	750	755	758	rVV	3090645	3592491	34.09%	4.128%
11	6.569	758	762	765	rVB	3198628	2591929	24.60%	2.978%
12	6.651	772	776	780	rBV	3191068	3407381	32.34%	3.915%
13	6.716	781	787	790	rVV	8235094	9349237	88.73%	10.743%
14	6.875	811	814	817	rBV	859879	728423	6.91%	0.837%
15	6.910	817	820	822	rVV	248471	180082	1.71%	0.207%
16	6.939	822	825	828	rVV	3027463	2416594	22.93%	2.777%
17	7.034	837	841	843	rBV	2435361	2012446	19.10%	2.312%
18	7.057	843	845	849	rVB	1021740	797635	7.57%	0.917%
19	7.128	853	857	859	rBV2	332393	320514	3.04%	0.368%
20	7.151	859	861	865	rVB	611983	467399	4.44%	0.537%
21	7.198	865	869	871	rBV	912025	768190	7.29%	0.883%
22	7.269	878	881	890	rBV	241192	234522	2.23%	0.269%
23	7.339	890	893	896	rBV	338098	336290	3.19%	0.386%
24	7.369	896	898	901	rVB	264251	211401	2.01%	0.243%
25	7.416	902	906	907	rBV	515780	522329	4.96%	0.600%
26	7.434	907	909	915	rVB	1875620	1797921	17.06%	2.066%
27	7.575	928	933	937	rBV2	135736	166629	1.58%	0.191%
28	8.157	1027	1032	1034	rBV	805963	726586	6.90%	0.835%
29	8.181	1034	1036	1039	rVB	353538	270427	2.57%	0.311%
30	8.545	1094	1098	1102	rBV3	148101	194865	1.85%	0.224%
31	8.622	1109	1111	1115	rVB2	152423	156552	1.49%	0.180%
32	8.722	1121	1128	1130	rBV5	152346	282300	2.68%	0.324%
33	8.786	1136	1139	1142	rBV3	198290	278980	2.65%	0.321%
34	8.875	1151	1154	1156	rVB	478088	394947	3.75%	0.454%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CROSSWICKS-DRUM-2

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	8.910	1157	1160	1162	rBV2	257323	230581	2.19%	0.265%
36	8.969	1168	1170	1172	rBV	324095	259148	2.46%	0.298%
37	9.039	1179	1182	1186	rBV4	315946	519024	4.93%	0.596%
38	9.122	1194	1196	1198	rBV2	215005	224099	2.13%	0.258%
39	9.157	1200	1202	1205	rVV3	482065	443559	4.21%	0.510%
40	9.186	1205	1207	1210	rVB2	410545	348328	3.31%	0.400%
41	9.233	1211	1215	1218	rBV	2319329	1956493	18.57%	2.248%
42	9.286	1222	1224	1225	rBV	250884	163355	1.55%	0.188%
43	9.322	1226	1230	1233	rVV3	953193	1168821	11.09%	1.343%
44	9.492	1257	1259	1261	rBV2	455054	499647	4.74%	0.574%
45	9.633	1280	1283	1285	rBV2	1621883	1796405	17.05%	2.064%
46	9.798	1308	1311	1314	rBV2	1288892	2094366	19.88%	2.407%
47	9.839	1316	1318	1321	rVB	2434179	2423656	23.00%	2.785%
48	9.928	1327	1333	1337	rBV7	1687521	3607095	34.23%	4.145%
49	14.904	2176	2179	2181	rBV2	465930	524894	4.98%	0.603%
50	14.968	2188	2190	2193	rBV2	234376	266664	2.53%	0.306%
51	15.539	2283	2287	2291	rVB	516903	716724	6.80%	0.824%
52	15.592	2294	2296	2303	rBV8	82643	174055	1.65%	0.200%
53	15.798	2328	2331	2342	rVB9	100912	298714	2.83%	0.343%
54	15.880	2342	2345	2354	rBV9	104050	325849	3.09%	0.374%
55	17.309	2584	2588	2595	rVB	136097	260258	2.47%	0.299%

Sum of corrected areas: 87026149

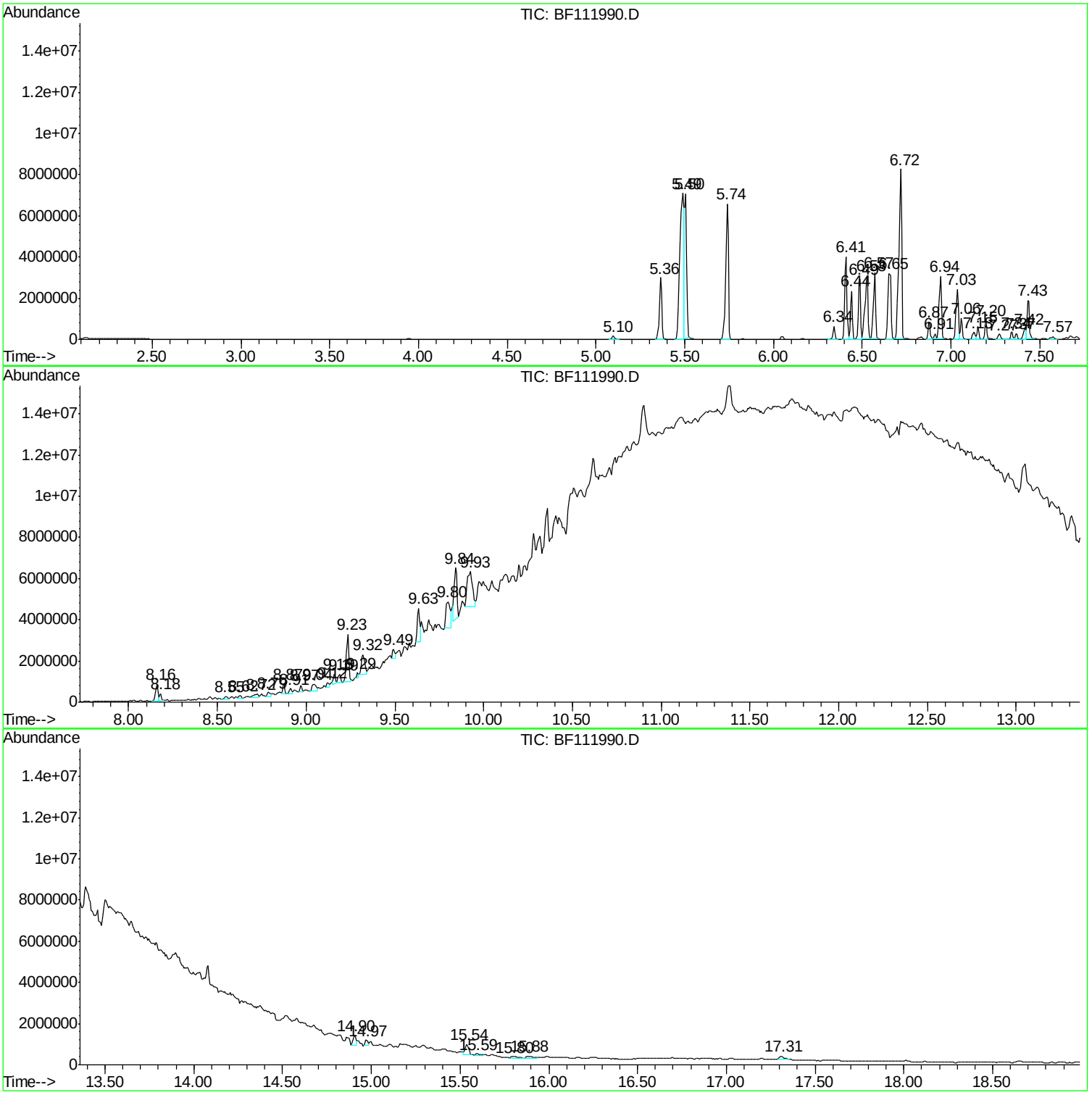
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CROSSWICKS-DRUM-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CROSSWICKS-DRUM-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

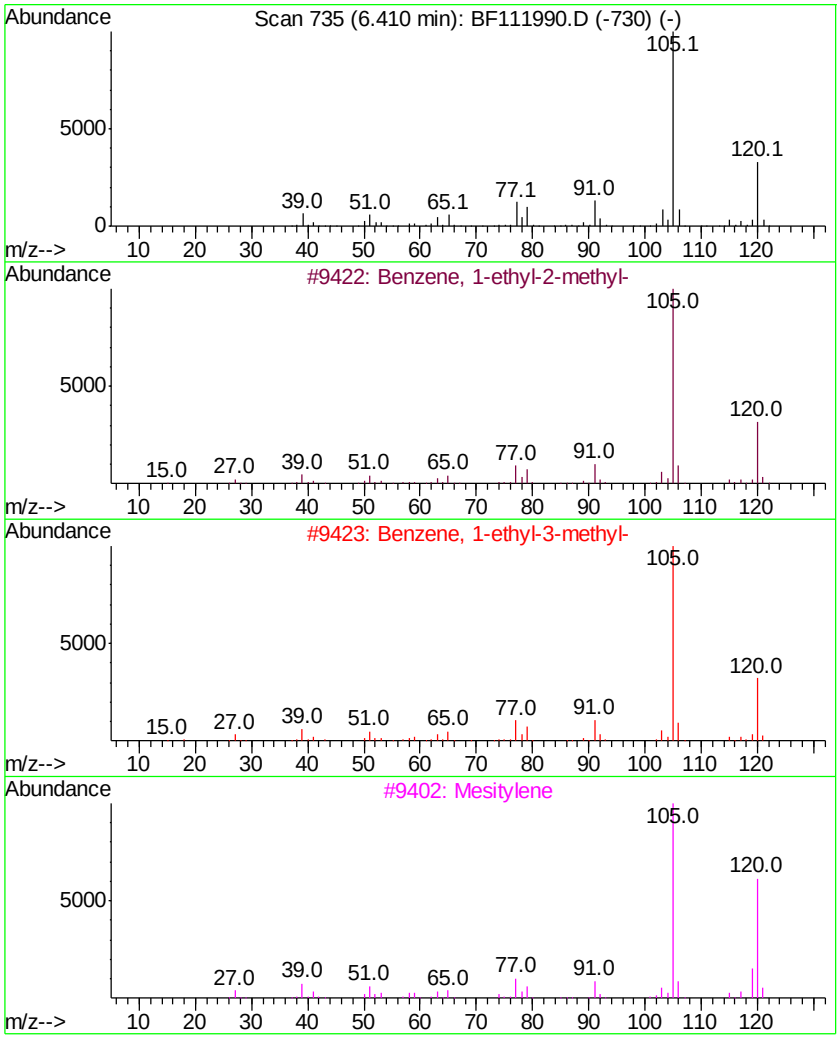
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

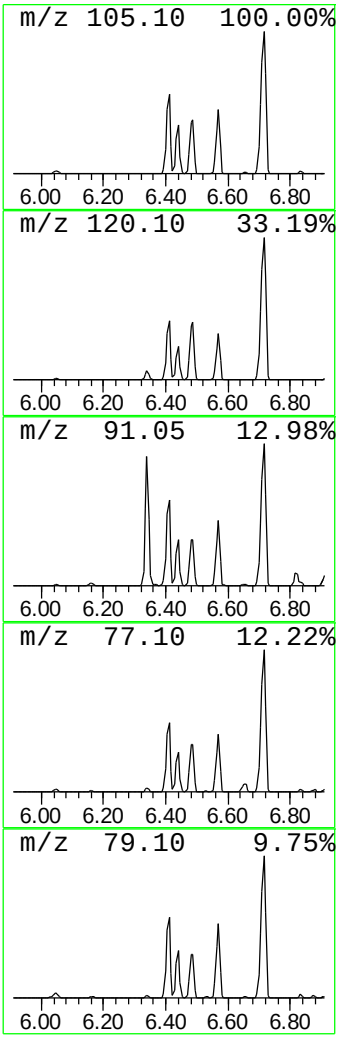
****

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.41	99.47 ng	3622650	1,4-Dichlorobenzene-d4	6.87

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-3-methyl-	120	C9H12	000620-14-4	95
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CROSSWICKS-DRUM-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

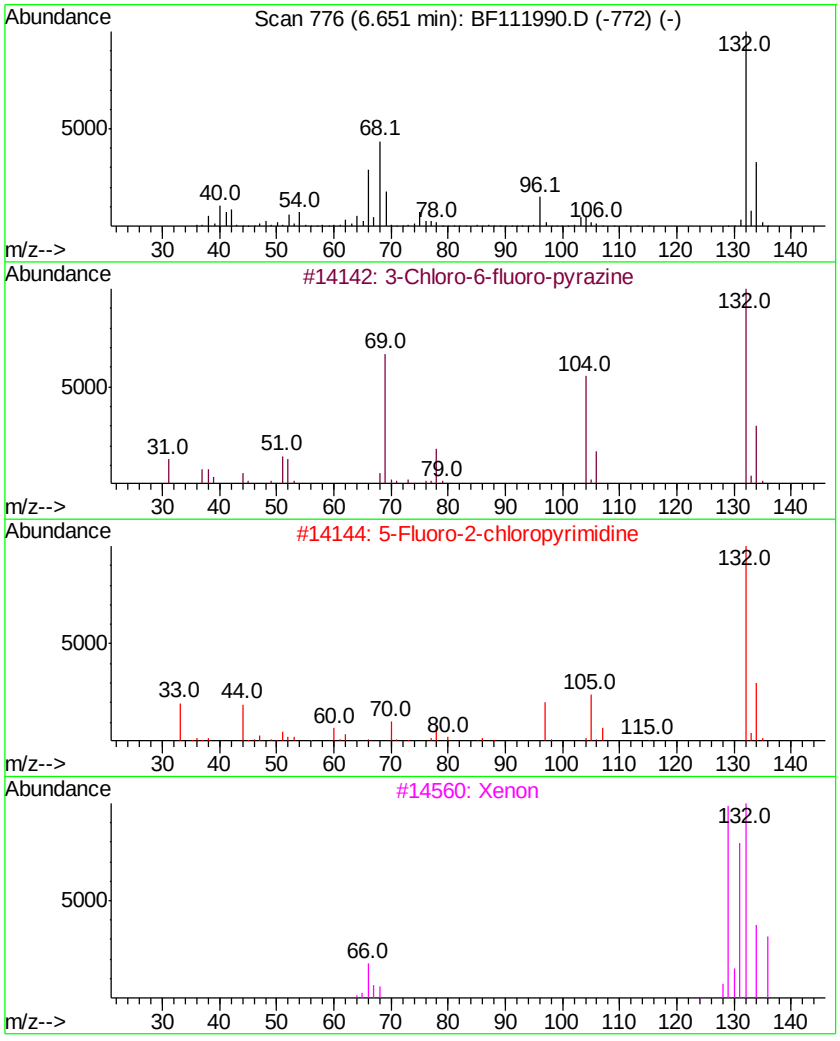
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

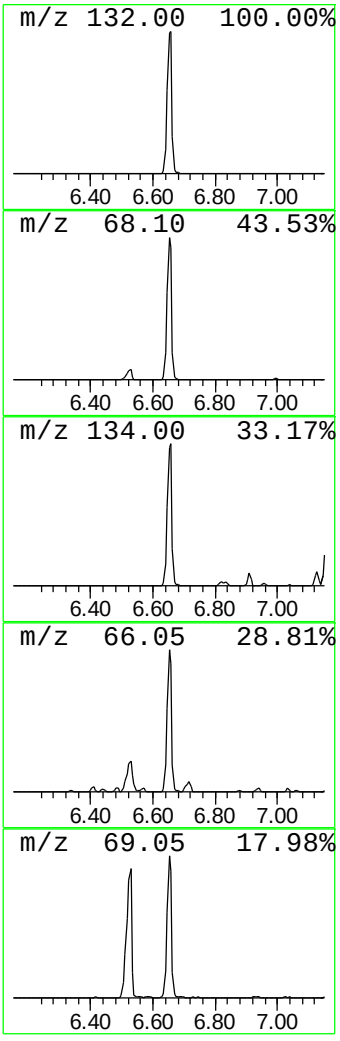
****

Peak Number 6 unknown6.65 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	93.56 ng	3407380	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3	Xenon	132	Xe	007440-63-3	9
4	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5	Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CROSSWICKS-DRUM-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

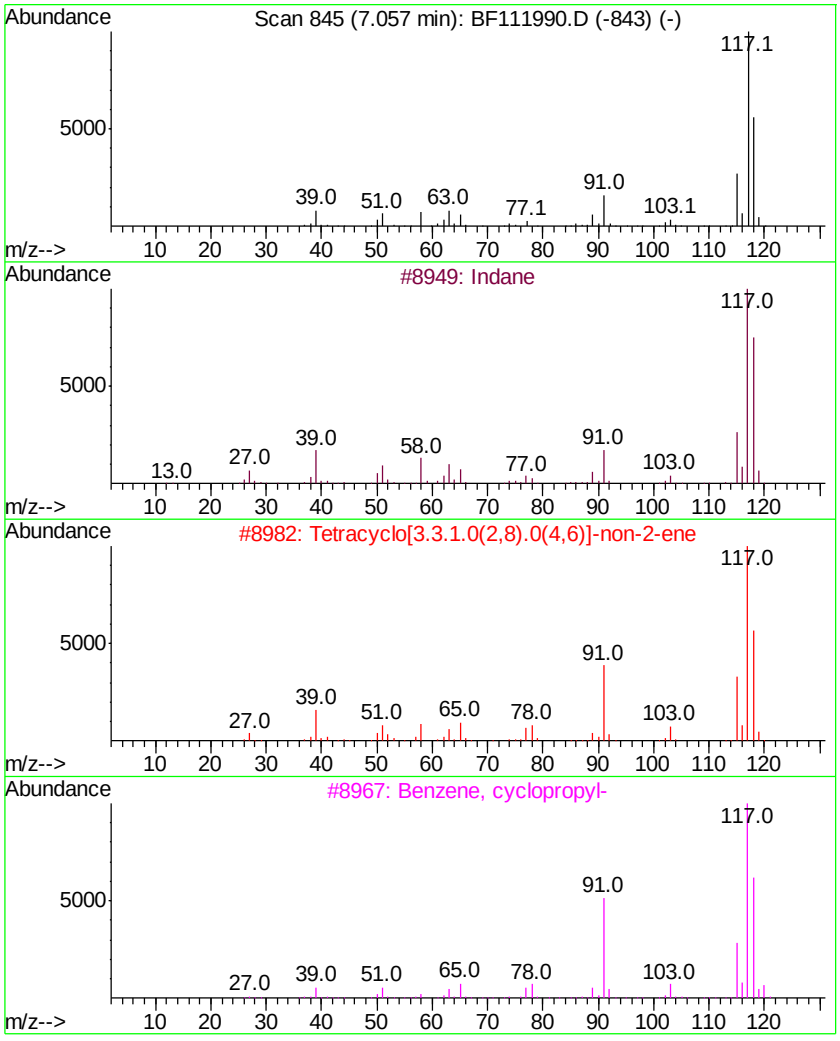
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

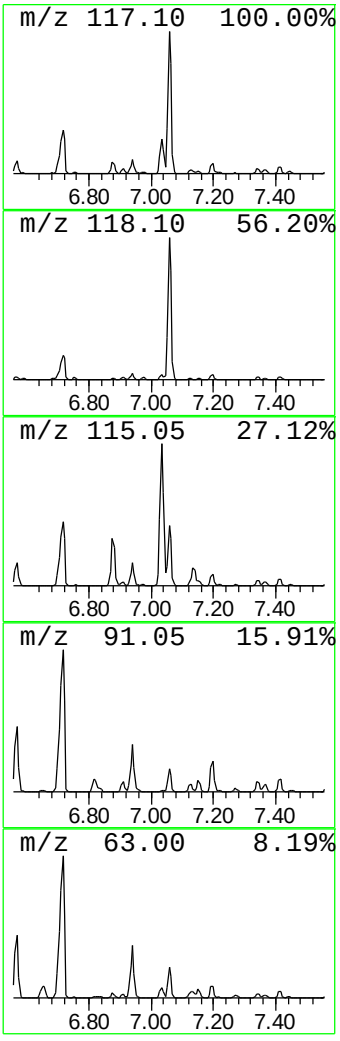
****

Peak Number 10 Indane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	21.90 ng	797635	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	91
2	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	83
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	83
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
5	trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	53





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CROSSWICKS-DRUM-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

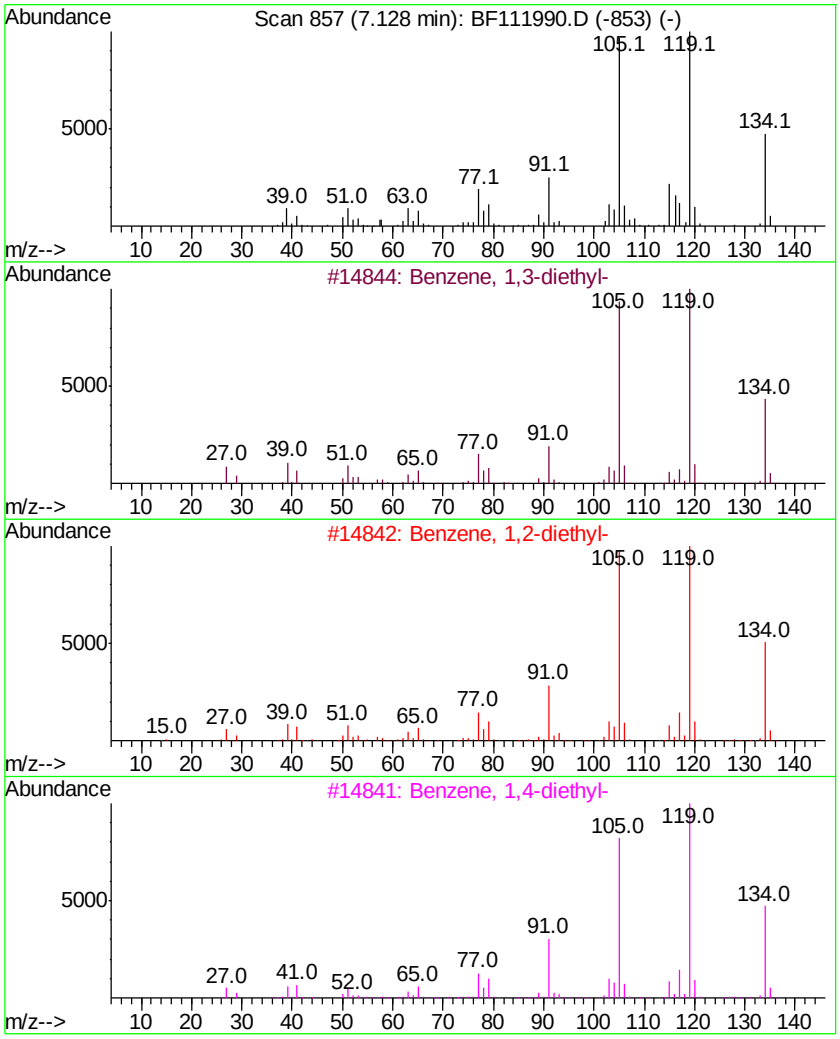
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

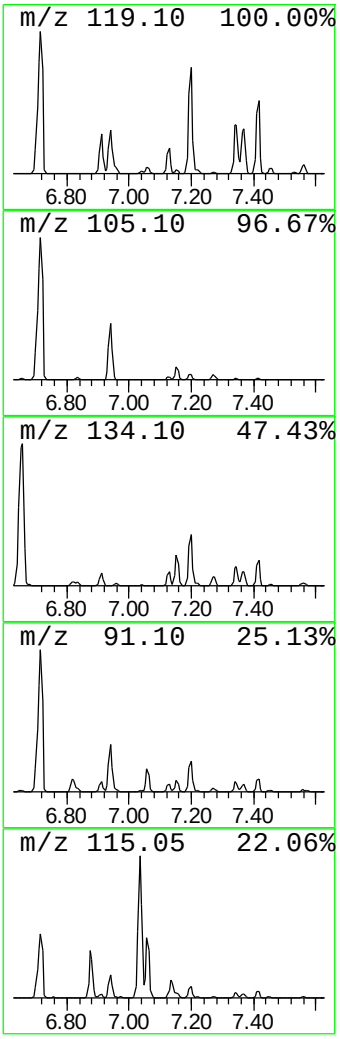
****

Peak Number 11 Benzene, 1,3-diethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.13	8.80 ng	320514	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	94
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	83
5	1,3,8-p-Menthatriene	134	C10H14	018368-95-1	74





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CROSSWICKS-DRUM-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

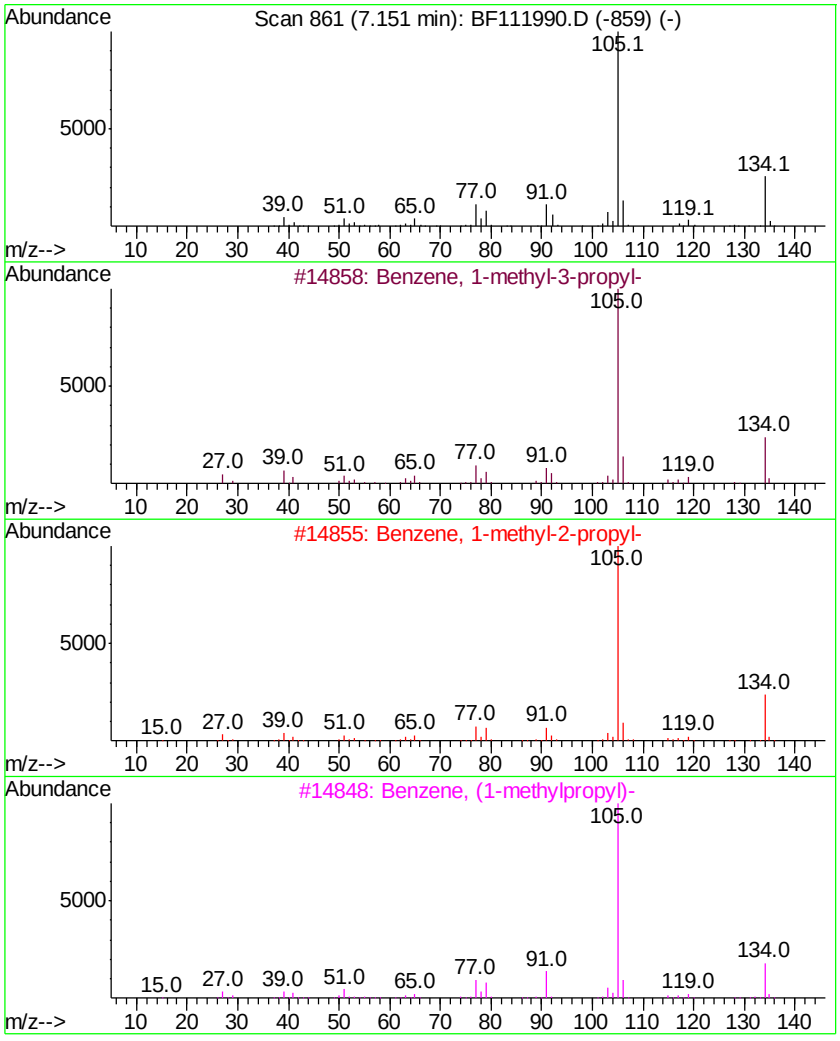
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

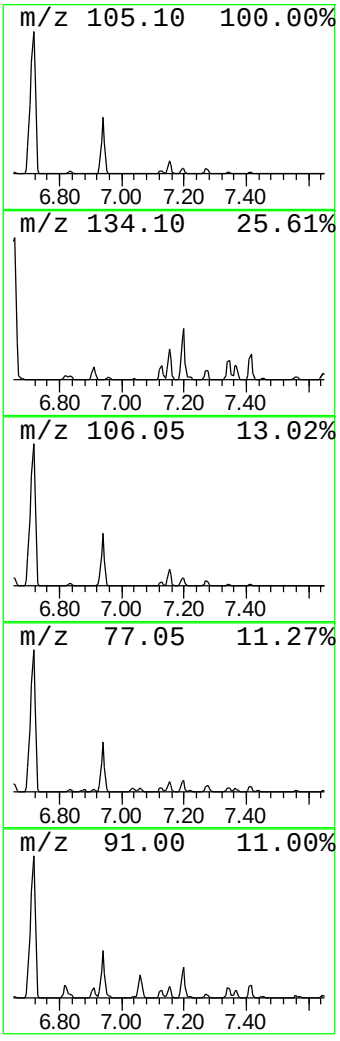
****

Peak Number 12 Benzene, 1-methyl-3-propyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.15	12.83 ng	467399	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	74





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CROSSWICKS-DRUM-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

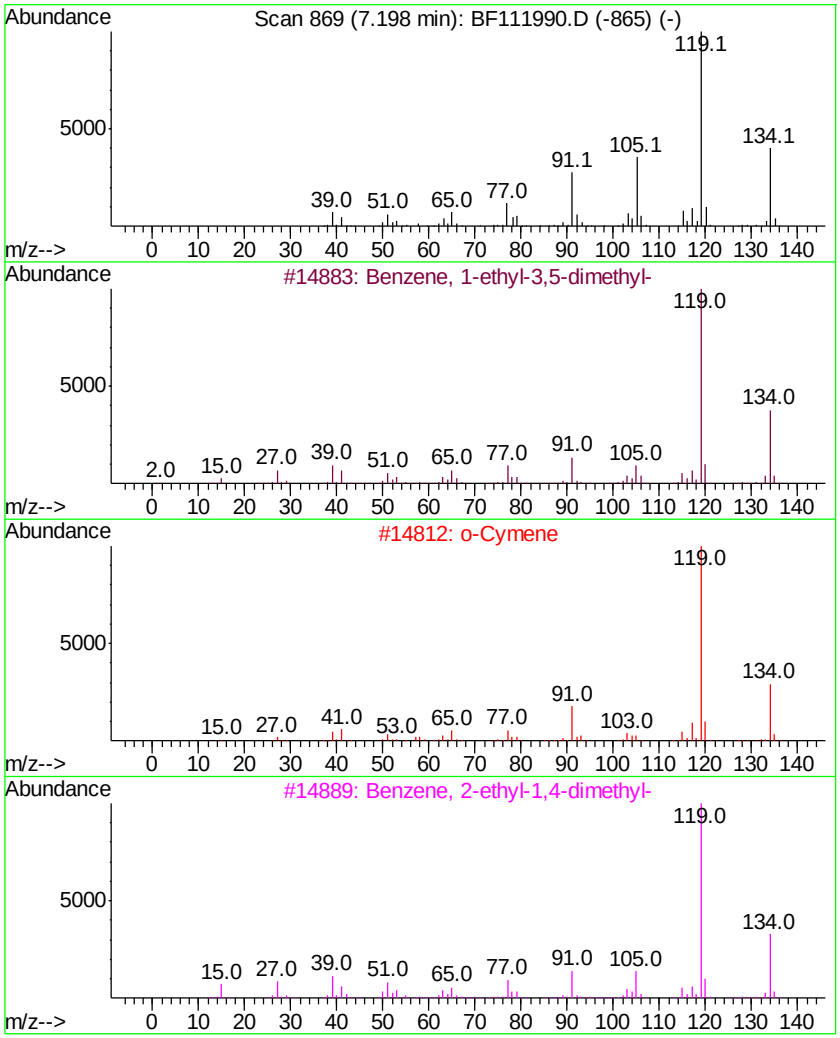
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

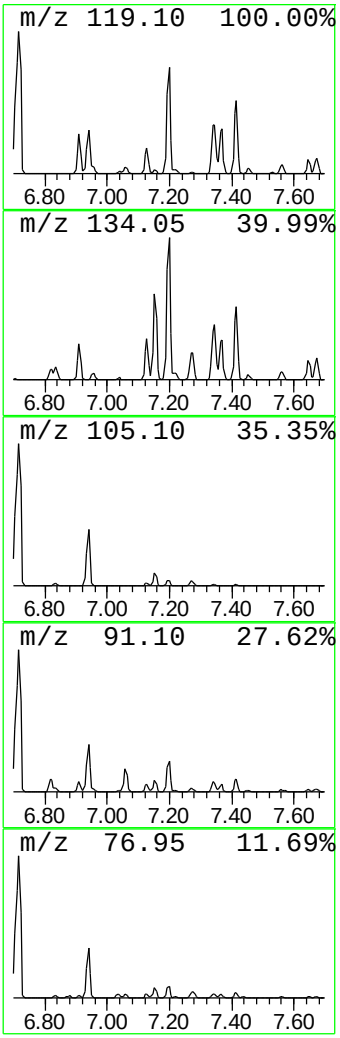
****

Peak Number 13 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	21.09 ng	768190	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
2	o-Cymene	134	C10H14	000527-84-4	93
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CROSSWICKS-DRUM-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

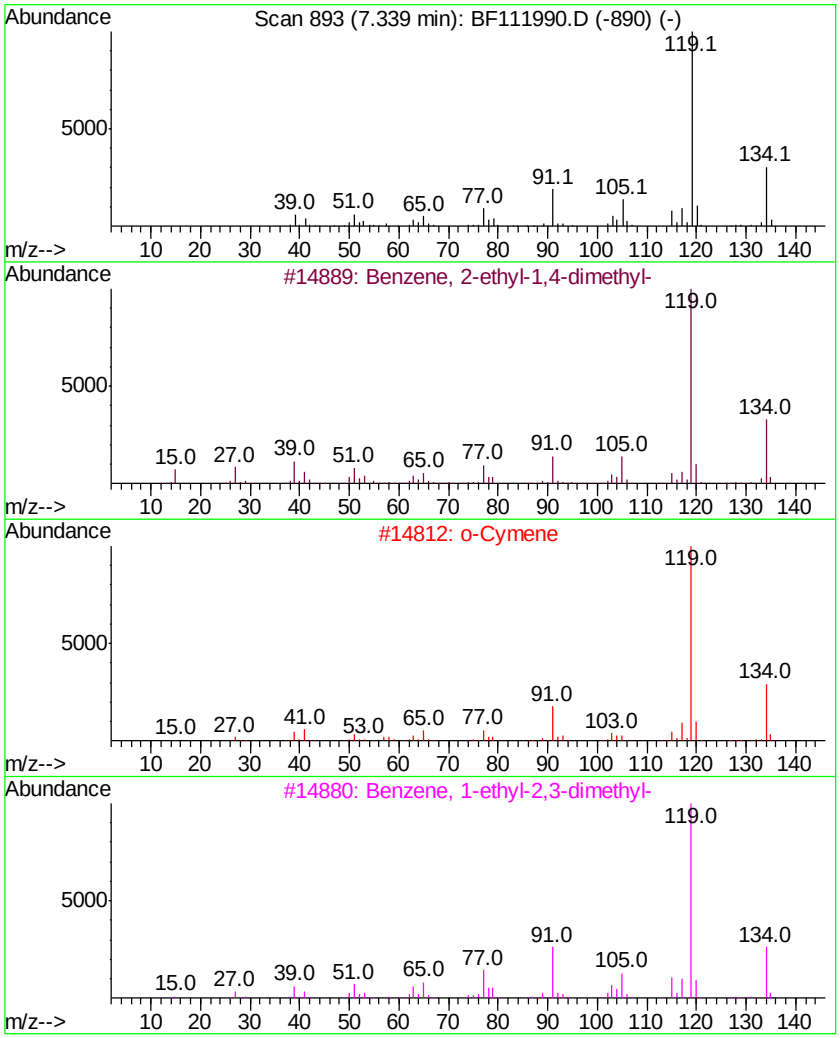
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

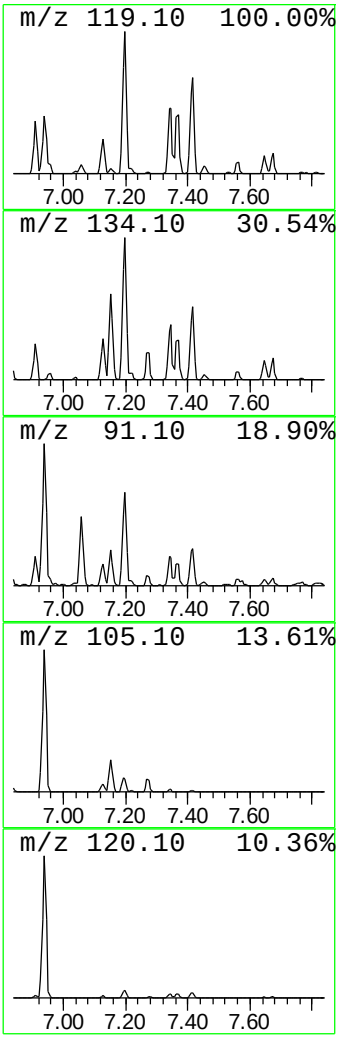
****

Peak Number 14 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.34	9.23 ng	336290	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CROSSWICKS-DRUM-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

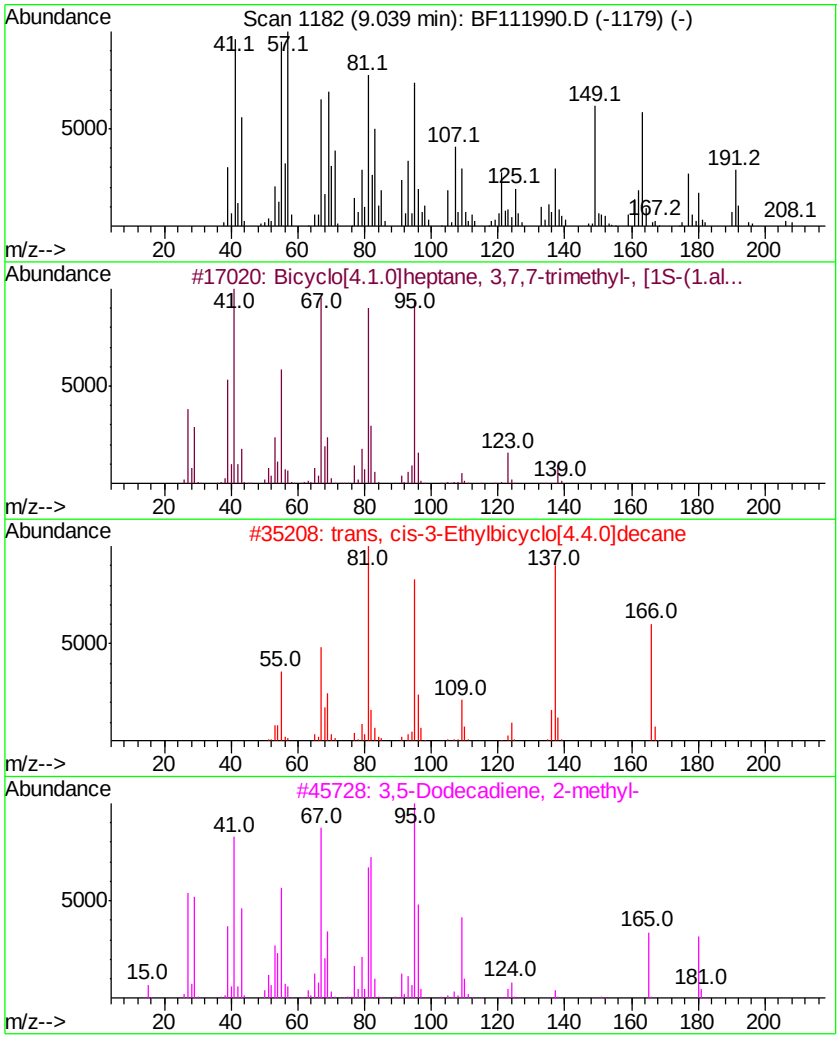
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

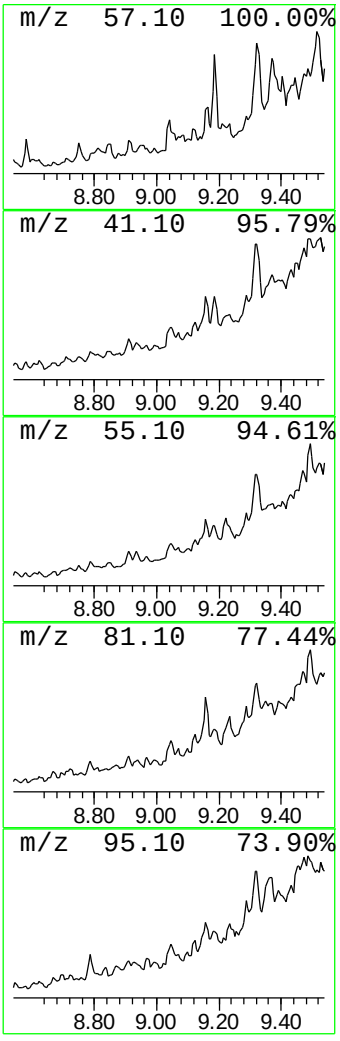
****

Peak Number 15 unknown9.04 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.04	14.29 ng	519024	Naphthalene-d8	8.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.1.0]heptane, 3,7,7-tri...	138	C10H18	006069-97-2	49
2	trans, cis-3-Ethylbicyclo[4.4.0]...	166	C12H22	066660-43-3	46
3	3,5-Dodecadiene, 2-methyl-	180	C13H24	055638-51-2	41
4	trans,cis-2,8-Dimethylspiro[5.5]...	180	C13H24	095472-52-9	38
5	cis,trans-2,9-Dimethylspiro[5.5]...	180	C13H24	095530-74-8	38





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Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CROSSWICKS-DRUM-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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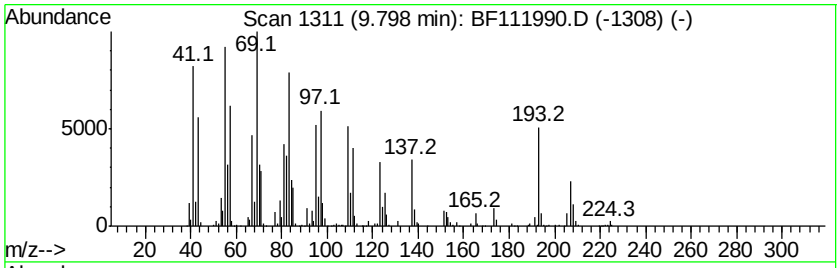
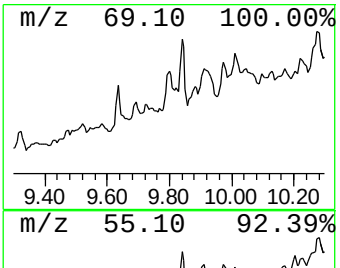
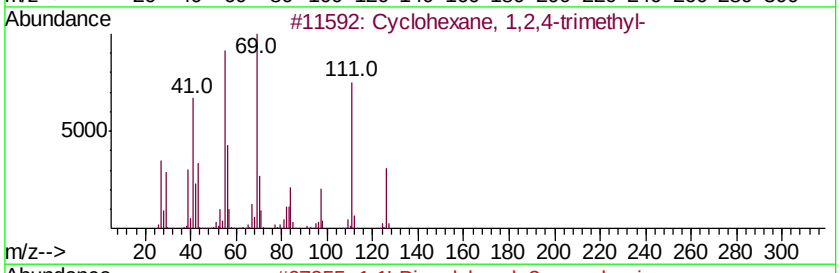
Peak Number 16 unknown9.80 Concentration Rank 16

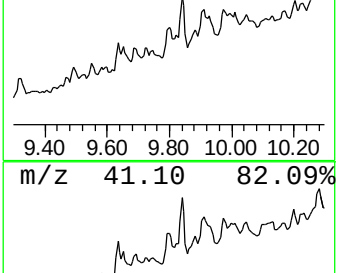
R.T.	EstConc	Area	Relative to ISTD	R.T.
9.80	11.61 ng	2094370	Acenaphthene-d10	9.93

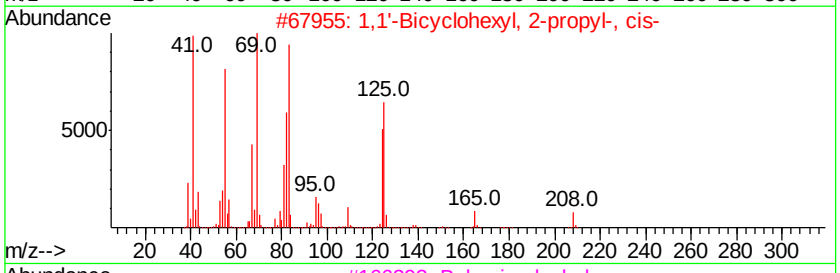
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	38
2	1,1'-Bicyclohexyl, 2-propyl-, cis-	208	C15H28	054934-88-2	38
3	Behenic alcohol	326	C22H46O	000661-19-8	38
4	1-Heneicosanol	312	C21H44O	015594-90-8	30
5	3-Hexene, 2,2,5,5-tetramethyl-, ...	140	C10H20	000692-47-7	30

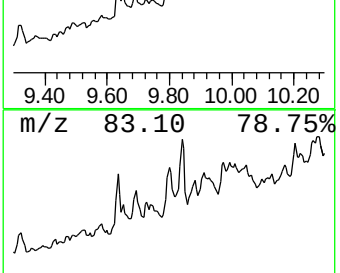


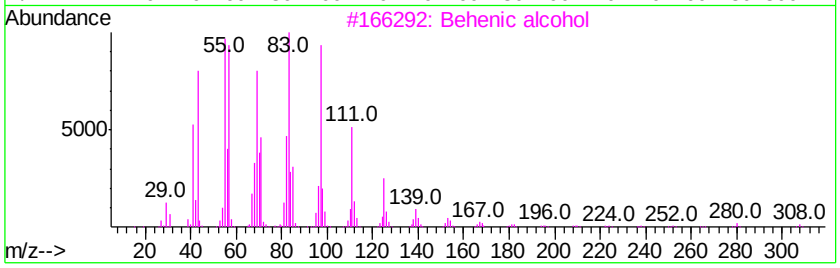


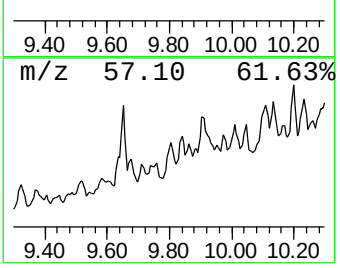












Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011019\
Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CROSSWICKS-DRUM-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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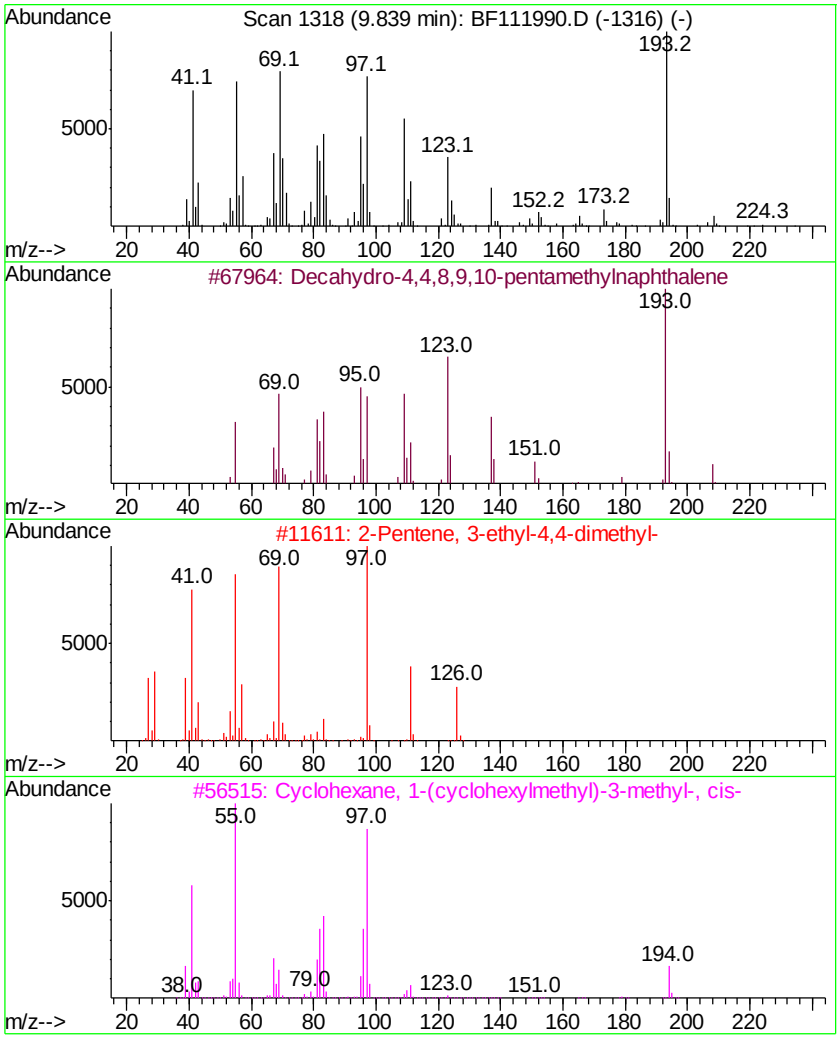
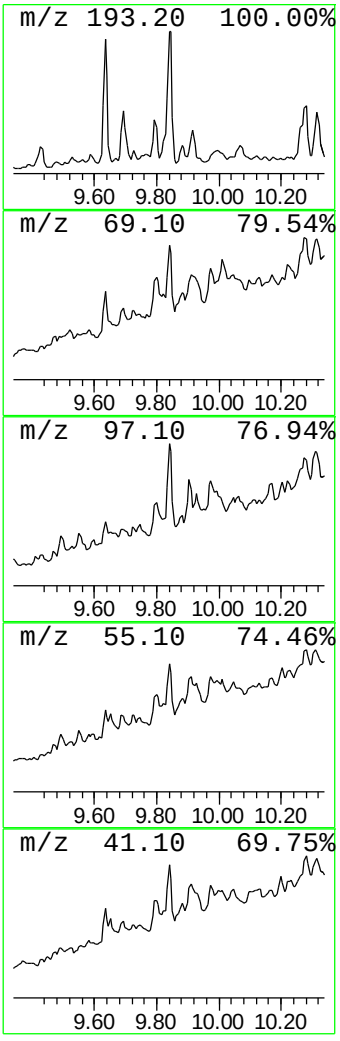
Peak Number 17 unknown9.84 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.84	13.44 ng	2423660	Acenaphthene-d10	9.93

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	30
2	2-Pentene, 3-ethyl-4,4-dimethyl-	126	C9H18	053907-59-8	30
3	Cyclohexane, 1-(cyclohexylmethyl...	194	C14H26	054823-96-0	18
4	Cyclopentanecarboxylic acid, 3-p...	324	C21H40O2	1000280-59-1	14
5	(5-Ethylcyclopent-1-enyl)methanol	126	C8H14O	036431-59-1	14





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Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : CROSSWICKS-DRUM-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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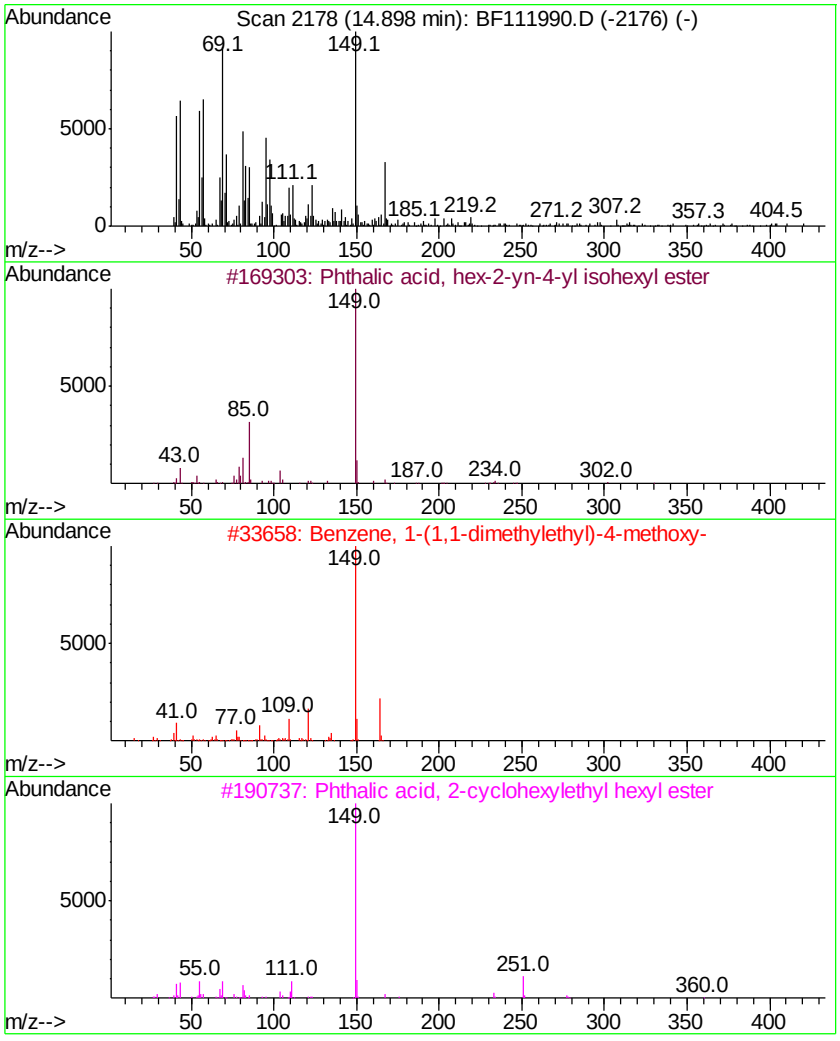
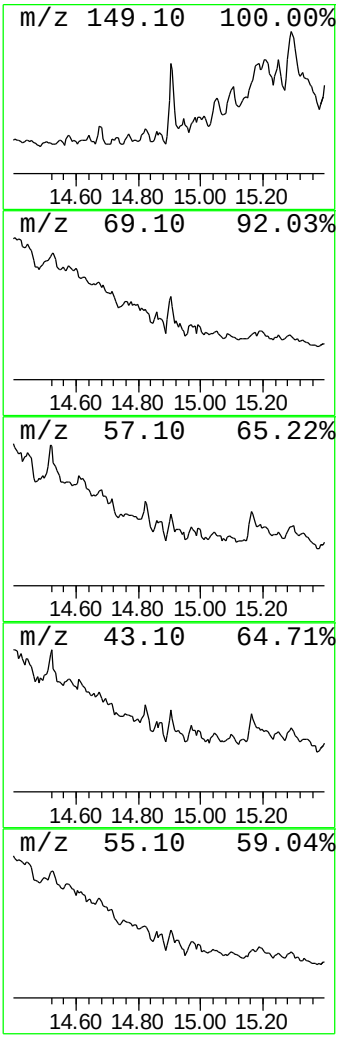
Peak Number 18 unknown14.90 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.90	14.65 ng	524894	Perylene-d12	15.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phthalic acid, hex-2-yn-4-yl iso...	330	C20H26O4	1000315-20-0	45
2	Benzene, 1-(1,1-dimethylethyl)-4...	164	C11H16O	005396-38-3	38
3	Phthalic acid, 2-cyclohexylethyl...	360	C22H32O4	1000309-87-6	38
4	Phthalic acid, 2-cyclohexylethyl...	318	C19H26O4	1000309-08-5	38
5	Benzothiazole, 2-methyl-	149	C8H7NS	000120-75-2	38





Library Search Compound Report

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Data File : BF111990.D
Acq On : 10 Jan 2019 16:33
Operator : JU/SJ
Sample : K1094-03
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CROSSWICKS-DRUM-2

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF010719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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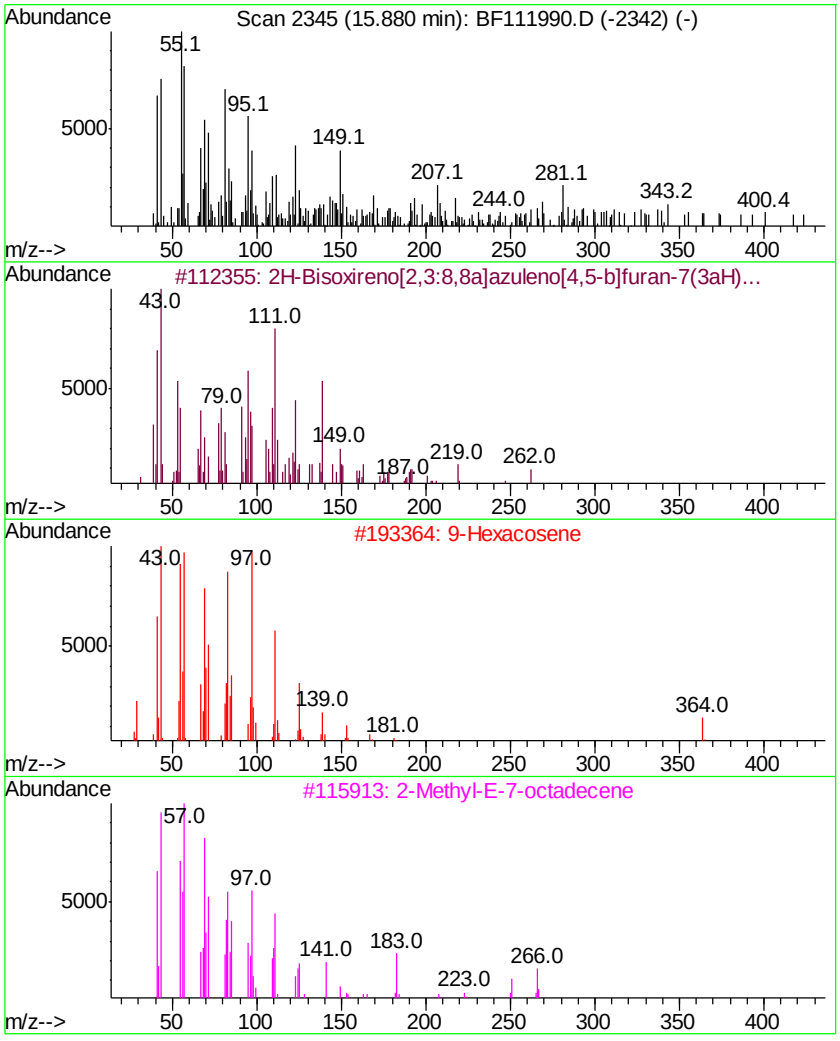
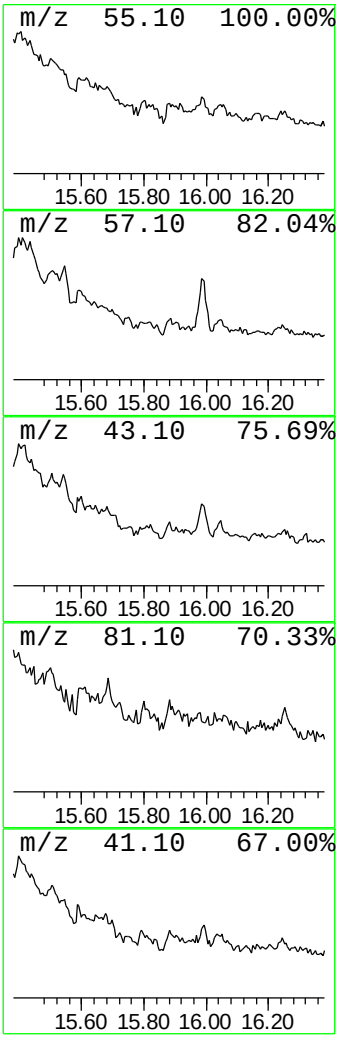
Peak Number 20 unknown15.88 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.88	9.09 ng	325849	Perylene-d12	15.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Bisoxireno[2,3:8,8a]azuleno[4...	262	C15H18O4	139343-89-8	38
2	9-Hexacosene	364	C26H52	071502-22-2	25
3	2-Methyl-E-7-octadecene	266	C19H38	1000130-92-3	25
4	2-Cyclohexen-1-one, 5-bromo-4,4-...	202	C8H11BrO	1000327-26-6	25
5	2,5,5,6,8a-Pentamethyl-trans-4a,...	208	C14H24O	1000215-77-8	25





Tentatively Identified Compound (LSC) summary

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Quant Title : ASP BNA STANDARDS FOR 5 POINT 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-ethyl-...	6.41	99.5	ng	3622650	1	6.87	728423	20.0
unknown6.65	6.65	93.6	ng	3407380	1	6.87	728423	20.0
Indane	7.06	21.9	ng	797635	1	6.87	728423	20.0
Benzene, 1,3-diet...	7.13	8.8	ng	320514	1	6.87	728423	20.0
Benzene, 1-methyl...	7.15	12.8	ng	467399	1	6.87	728423	20.0
Benzene, 1-ethyl-...	7.20	21.1	ng	768190	1	6.87	728423	20.0
Benzene, 2-ethyl-...	7.34	9.2	ng	336290	1	6.87	728423	20.0
unknown9.04	9.04	14.3	ng	519024	2	8.16	726586	20.0
unknown9.80	9.80	11.6	ng	2094370	3	9.93	3607100	20.0
unknown9.84	9.84	13.4	ng	2423660	3	9.93	3607100	20.0
unknown14.90	14.90	14.7	ng	524894	6	15.54	716724	20.0
unknown15.88	15.88	9.1	ng	325849	6	15.54	716724	20.0