LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
Data File : BF107623.D
Acq On : 24 Jul 2018 17:11
Operator : JU/SJ
Sample : J4096-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-10

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-BF072018.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	5.190	482	485	498	rBV	267778	345243	7.45%	0.680%
2	5.601	551	555	579	rBV	1268937	2142151	46.20%	4.219%
3	6.242	660	664	670	rVB	315214	276358	5.96%	0.544%
4	6.601	721	725	733	rBV	1444875	1944007	41.93%	3.829%
5	6.660	733	735	739	rVB	163430	179657	3.87%	0.354%
6	6.742	745	749	771	rVB	1938347	2386872	51.48%	4.701%
7	6.966	784	787	796	rBV	1345192	1453004	31.34%	2.862%
8	7.037	796	799	802	rVV	95950	112200	2.42%	0.221%
9	7.072	802	805	808	rVV	223414	175931	3.79%	0.346%
10	7.119	810	813	828	rVB	1544537	1626431	35.08%	3.203%
11	7.531	879	883	900	rBV	1024467	1360659	29.35%	2.680%
12	8.248	1001	1005	1025	rBV	1826524	1977981	42.66%	3.895%
13	8.978	1124	1129	1133	rBV2	78921	81657	1.76%	0.161%
14	9.319	1184	1187	1197	rBV	2227785	2181314	47.05%	4.296%
15	9.713	1248	1254	1262	rVB	201408	212990	4.59%	0.419%
16	10.007	1297	1304	1308	rBV	1877663	1916637	41.34%	3.775%
17	10.042	1308	1310	1318	rVB	195191	201284	4.34%	0.396%
18	10.154	1326	1329	1332	rBV2	51406	54676	1.18%	0.108%
19	10.213	1335	1339	1342	rBV2	101650	119542	2.58%	0.235%
20	10.560	1394	1398	1404	rBV	421238	427580	9.22%	0.842%
21	10.642	1407	1412	1415	rVV3	125838	161304	3.48%	0.318%
22	10.713	1422	1424	1429	rVB	69017	71201	1.54%	0.140%
23	10.801	1435	1439	1453	rBV	1212062	1402662	30.25%	2.762%
24	10.907	1453	1457	1464	rVB2	168744	219515	4.73%	0.432%
25	10.972	1465	1468	1469	rBV	88773	72276	1.56%	0.142%
26	11.001	1469	1473	1477	rVB5	67133	108004	2.33%	0.213%
27	11.036	1477	1479	1482	rBV	68844	71260	1.54%	0.140%
28	11.107	1486	1491	1495	rBV3	110344	195097	4.21%	0.384%
29	11.189	1502	1505	1507	rBV	103871	82920	1.79%	0.163%
30	11.372	1533	1536	1538	rBV	186990	160559	3.46%	0.316%
31	11.401	1538	1541	1548	rVB	311818	289289	6.24%	0.570%
32	11.507	1555	1559	1561	rBV	1588937	1827570	39.42%	3.599%
33	11.530	1561	1563	1569	rVV	4432570	4360368	94.04%	8.587%
34	11.583	1569	1572	1575	rVV	732519	791805	17.08%	1.559%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
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Acq On : 24 Jul 2018 17:11
Operator : JU/SJ
Sample : J4096-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-10

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-BF072018.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.613	1575	1577	1582	rVB2	132865	158771	3.42%	0.313%
36	11.724	1593	1596	1597	rBV2	144735	128731	2.78%	0.254%
37	12.007	1641	1644	1646	rBV	406103	398563	8.60%	0.785%
38	12.036	1646	1649	1653	rVB2	320386	360001	7.76%	0.709%
39	12.119	1660	1663	1672	rVB2	895929	1123014	24.22%	2.212%
40	12.301	1691	1694	1696	rBV	258713	247241	5.33%	0.487%
41	12.330	1696	1699	1702	rVV	282316	263614	5.69%	0.519%
42	12.677	1755	1758	1761	rBV3	182496	222395	4.80%	0.438%
43	12.730	1762	1767	1769	rVV	4494803	4636623	100.00%	9.131%
44	12.754	1769	1771	1774	rVB	1724249	1479164	31.90%	2.913%
45	12.960	1799	1806	1811	rBV	3502907	4269293	92.08%	8.408%
46	13.060	1820	1823	1824	rBV	392653	420599	9.07%	0.828%
47	13.089	1825	1828	1831	rVB	1501982	1413401	30.48%	2.784%
48	13.348	1869	1872	1875	rBV2	582348	541669	11.68%	1.067%
49	14.113	1999	2002	2005	rBV	1065607	1042510	22.48%	2.053%
50	14.160	2006	2010	2013	rVV2	1693734	2780054	59.96%	5.475%
51	14.189	2013	2015	2021	rVB	1397315	1385081	29.87%	2.728%
52	15.636	2258	2261	2266	rVB	713103	916175	19.76%	1.804%

Sum of corrected areas: 50776903

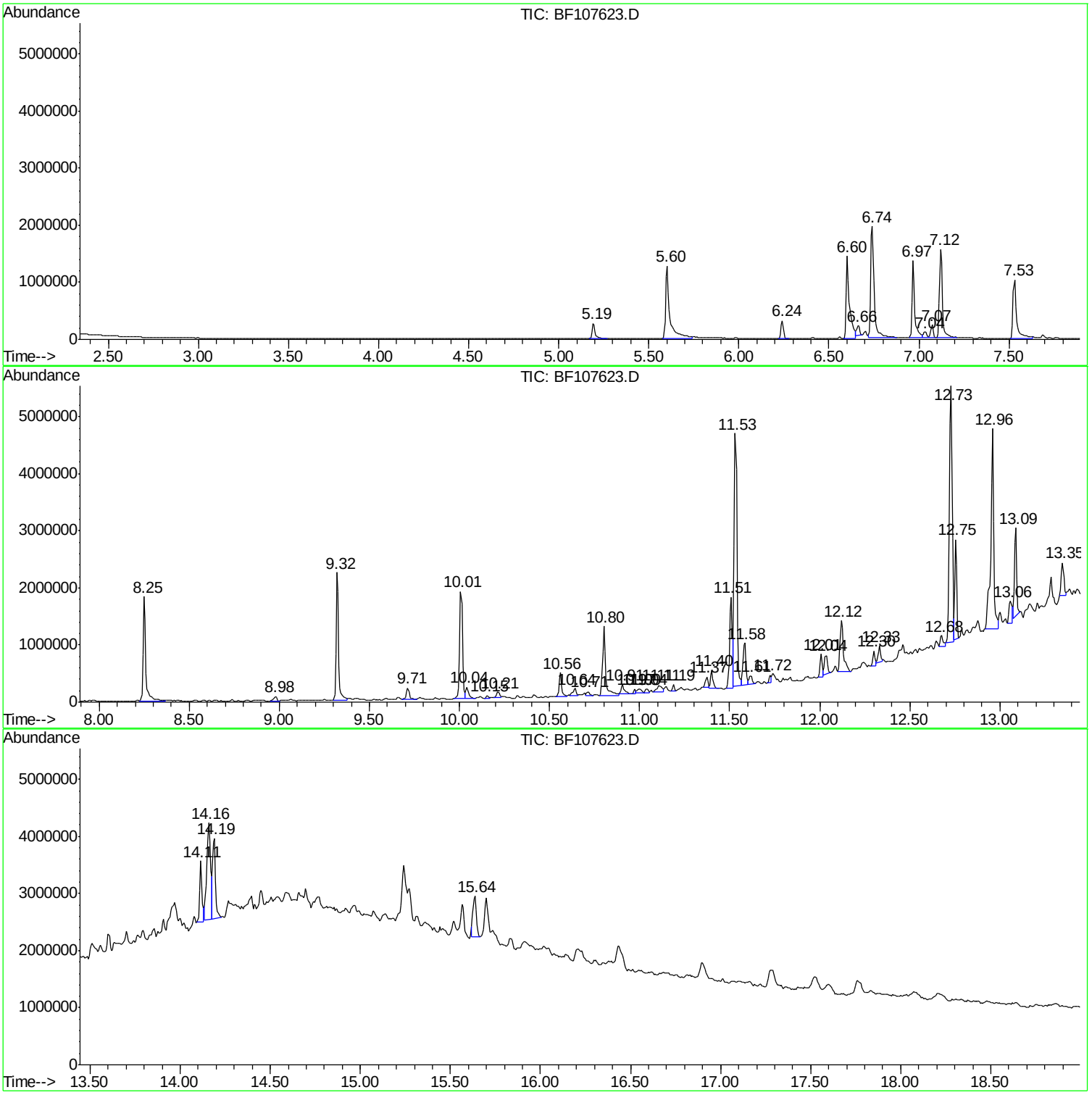
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
Data File : BF107623.D
Acq On : 24 Jul 2018 17:11
Operator : JU/SJ
Sample : J4096-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
Data File : BF107623.D
Acq On : 24 Jul 2018 17:11
Operator : JU/SJ
Sample : J4096-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

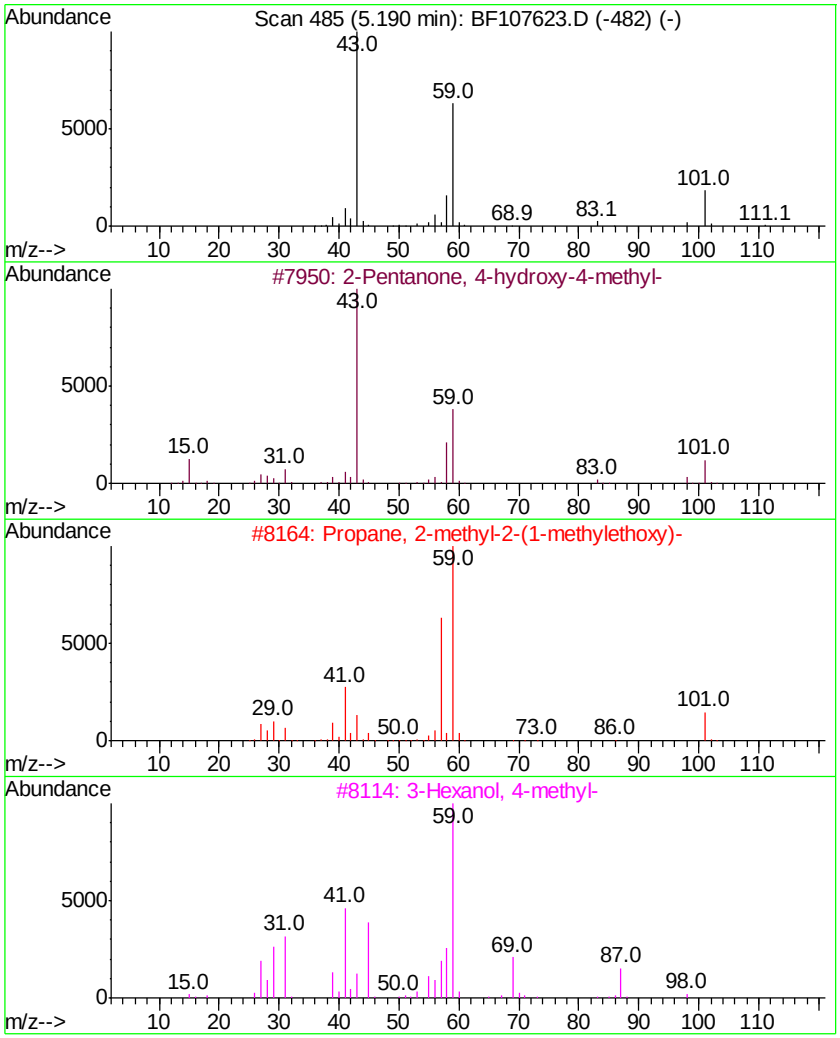
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

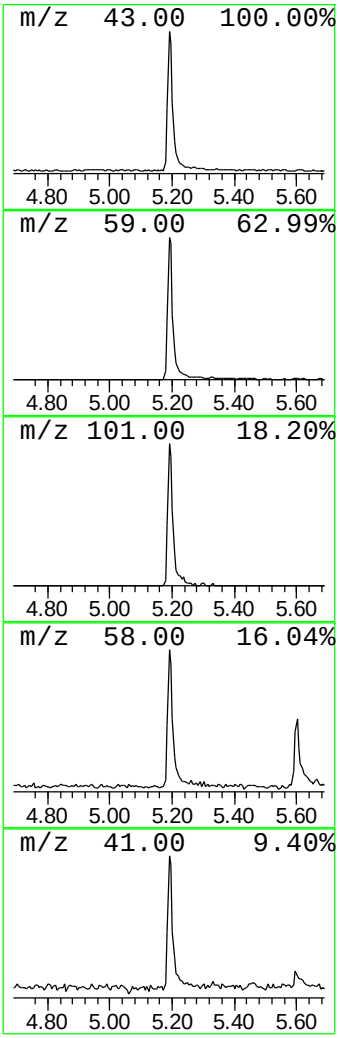
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	4.75 ng	345243	1,4-Dichlorobenzene-d4	6.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	53
3	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
Data File : BF107623.D
Acq On : 24 Jul 2018 17:11
Operator : JU/SJ
Sample : J4096-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

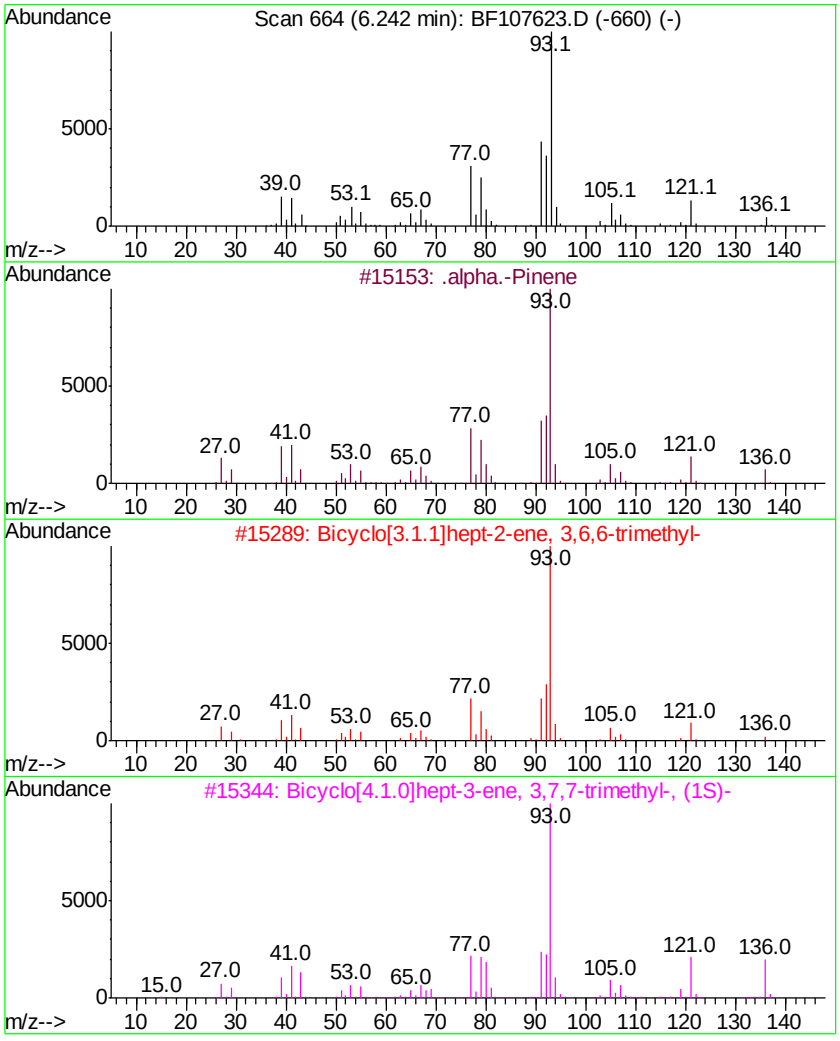
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

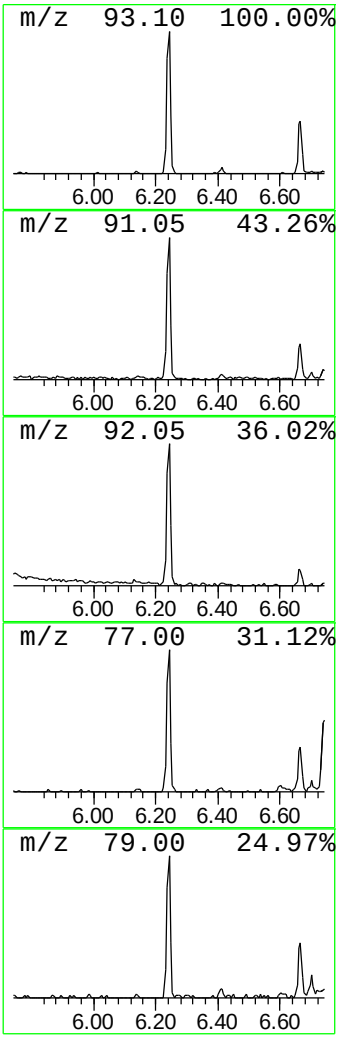
****

Peak Number 2 .alpha.-Pinene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.24	3.80 ng	276358	1,4-Dichlorobenzene-d4	6.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Pinene	136	C10H16	000080-56-8	95
2	Bicyclo[3.1.1]hept-2-ene, 3,6,6-...	136	C10H16	004889-83-2	94
3	Bicyclo[4.1.0]hept-3-ene, 3,7,7-...	136	C10H16	000498-15-7	91
4	1,3,6-Octatriene, 3,7-dimethyl-,...	136	C10H16	003779-61-1	91
5	1R-.alpha.-Pinene	136	C10H16	007785-70-8	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
Data File : BF107623.D
Acq On : 24 Jul 2018 17:11
Operator : JU/SJ
Sample : J4096-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

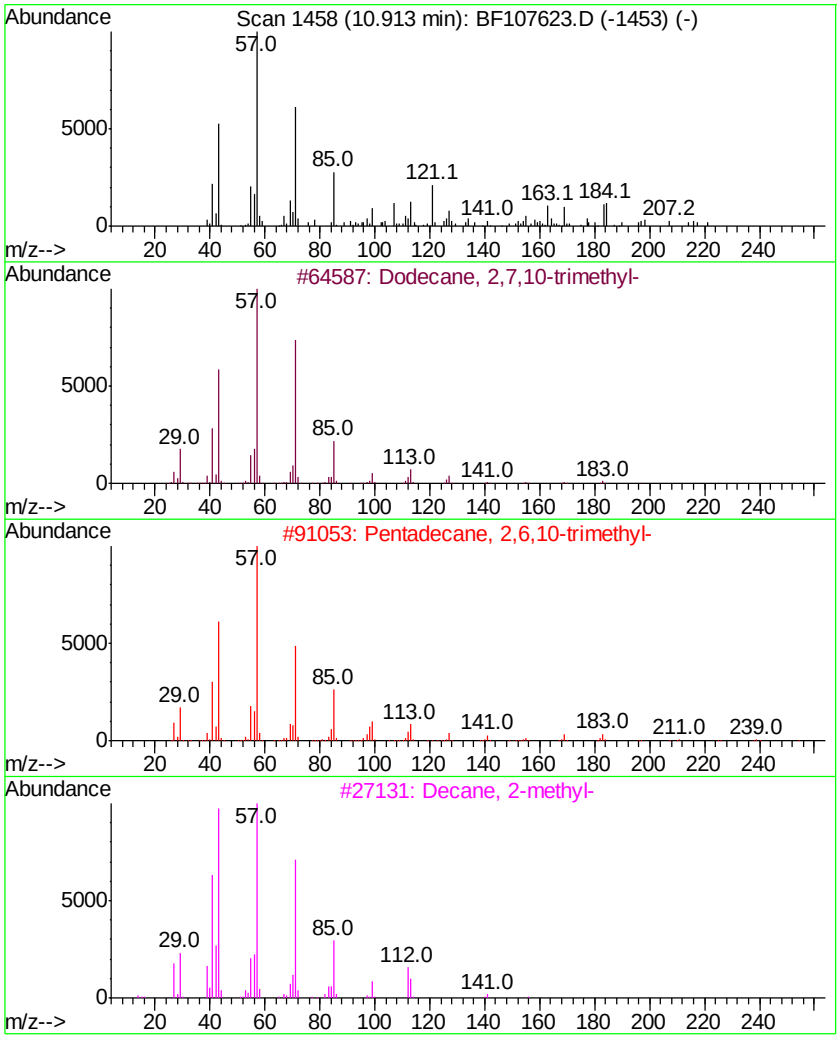
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

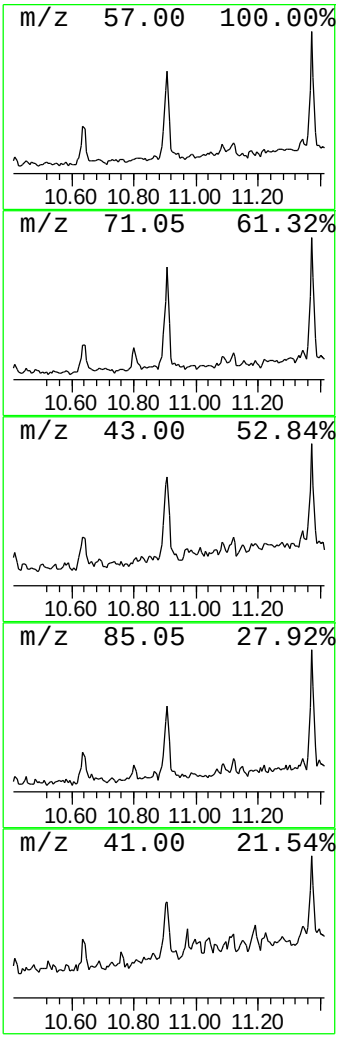
****

Peak Number 3 Dodecane, 2,7,10-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.91	2.40 ng	219515	Phenanthrene-d10	11.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	86
2	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	86
3	Decane, 2-methyl-	156	C11H24	006975-98-0	80
4	Hexadecane, 3-methyl-	240	C17H36	006418-43-5	78
5	Tridecane, 5-propyl-	226	C16H34	055045-11-9	78





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
Data File : BF107623.D
Acq On : 24 Jul 2018 17:11
Operator : JU/SJ
Sample : J4096-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

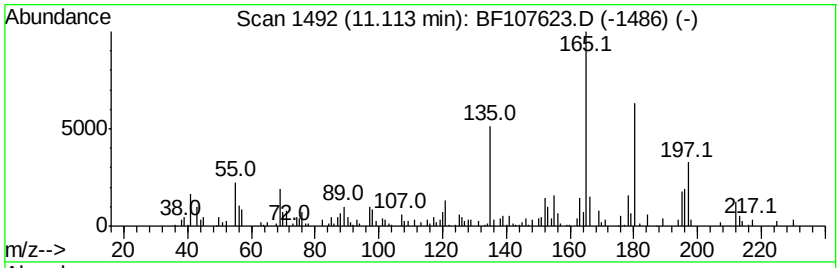
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

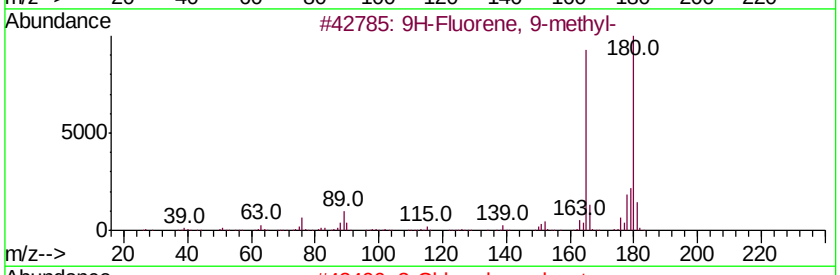
****

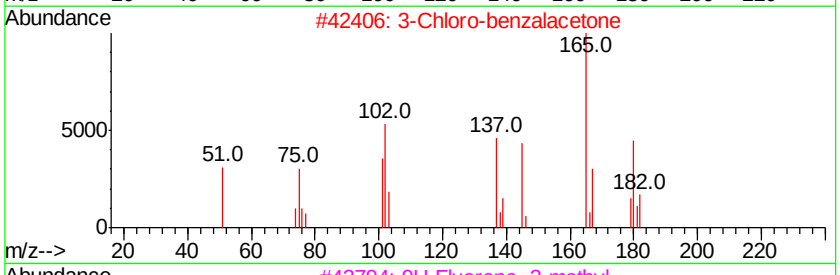
Peak Number 4 9H-Fluorene, 9-methyl- Concentration Rank 12

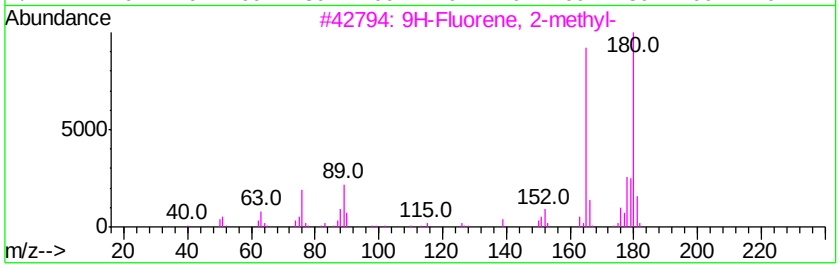
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.11	2.14 ng	195097	Phenanthrene-d10	11.51

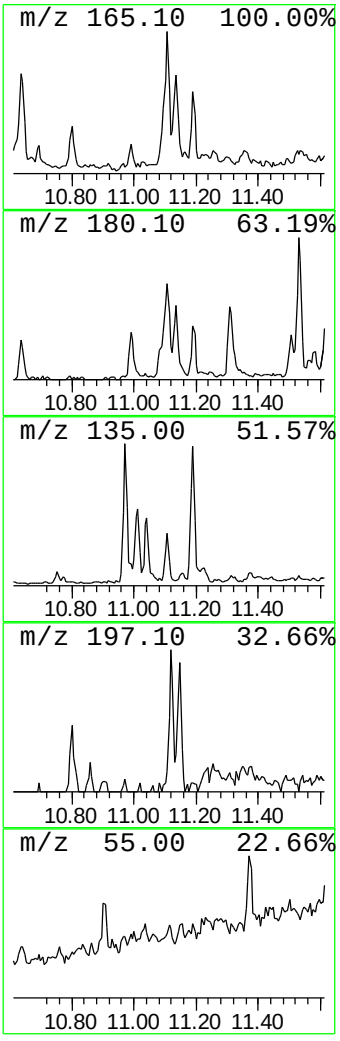
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	64
2	3-Chloro-benzalacetone	180	C10H9ClO	1000146-90-5	55
3	9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	53
4	Benzenamine, N-ethyl-N-methyl-4-...	180	C9H12N2O2	056269-48-8	49
5	Ethanone, 1-(3,4-dimethoxyphenyl)-	180	C10H12O3	001131-62-0	46











Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
Data File : BF107623.D
Acq On : 24 Jul 2018 17:11
Operator : JU/SJ
Sample : J4096-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

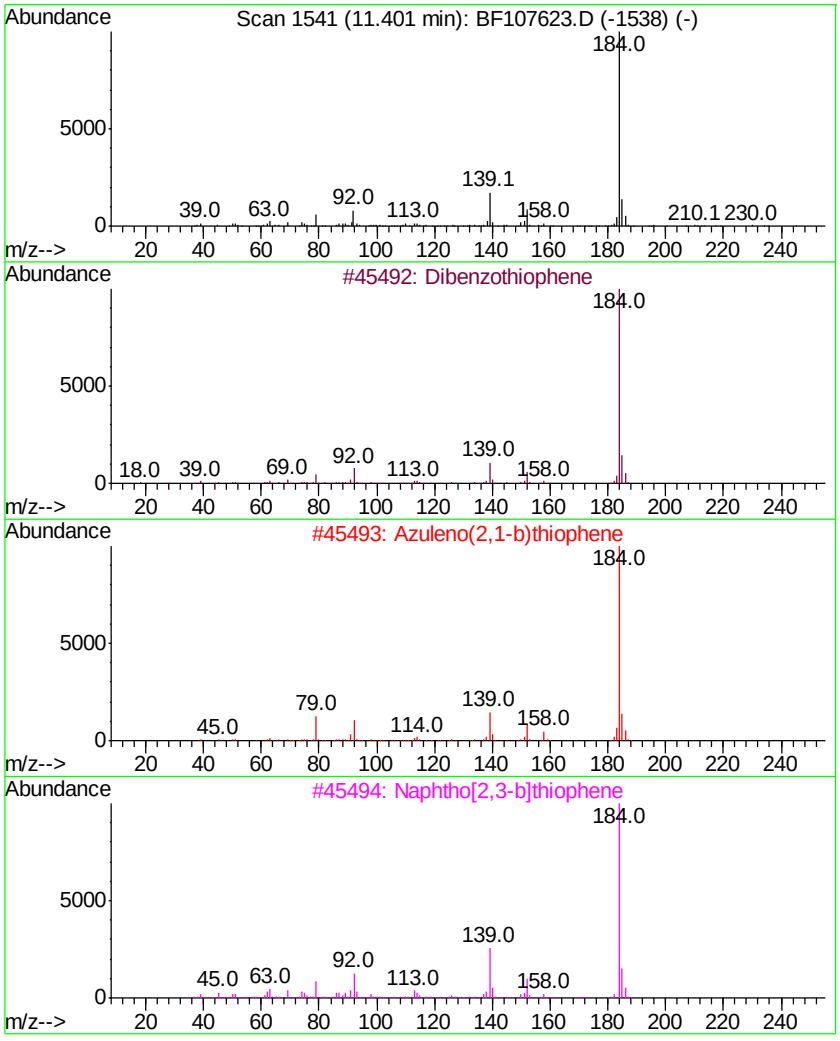
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

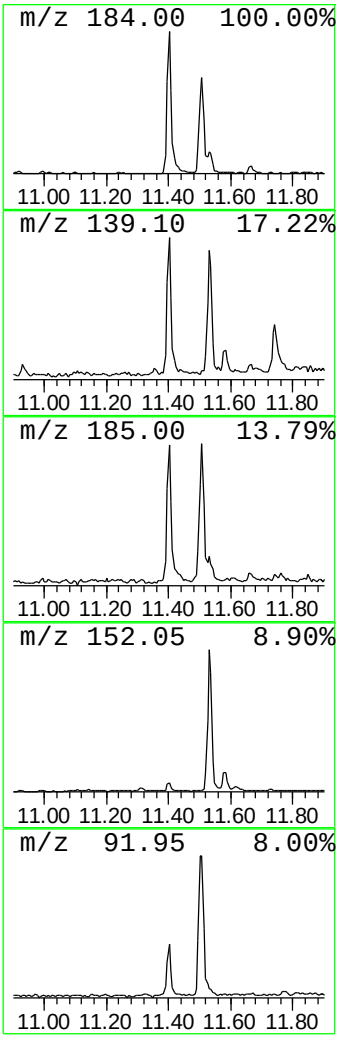
****

Peak Number 5 Dibenzothiophene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.40	3.17 ng	289289	Phenanthrene-d10	11.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dibenzothiophene	184	C12H8S	000132-65-0	97
2	Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	87
3	Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	86
4	Dibenzothiophene, 5-oxide	200	C12H8OS	001013-23-6	72
5	1,1'-Biphenyl, 2-methoxy-	184	C13H12O	000086-26-0	72





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
Data File : BF107623.D
Acq On : 24 Jul 2018 17:11
Operator : JU/SJ
Sample : J4096-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

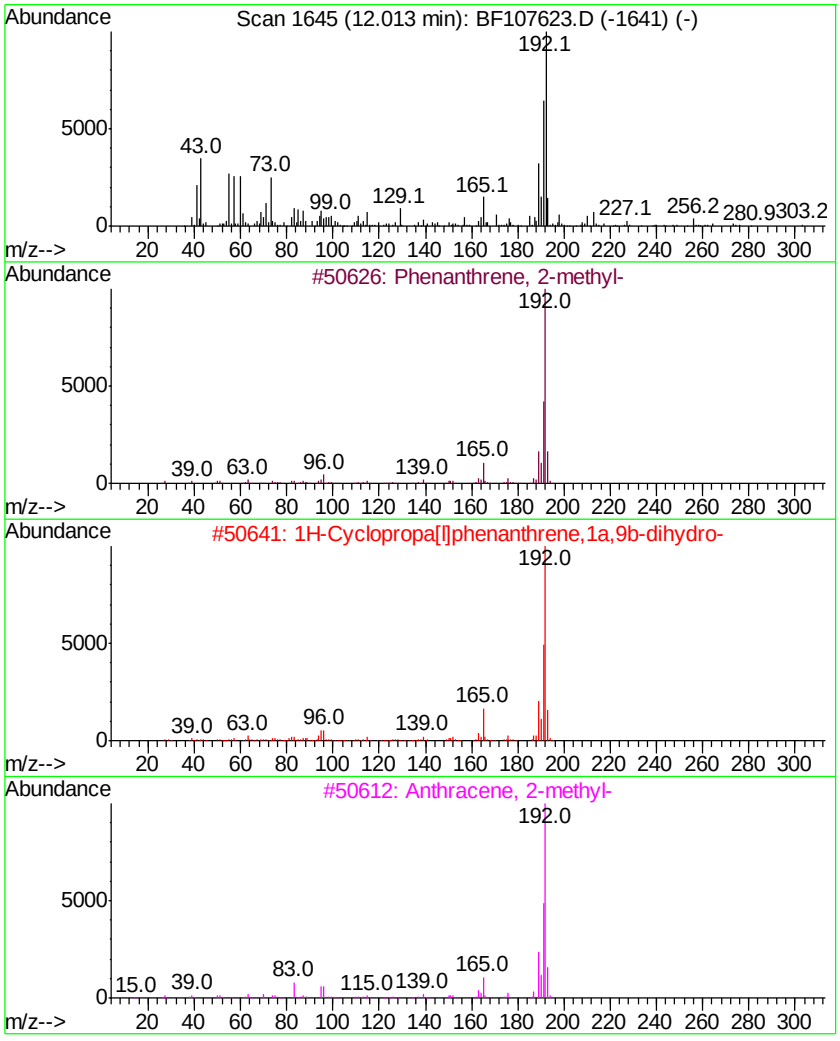
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

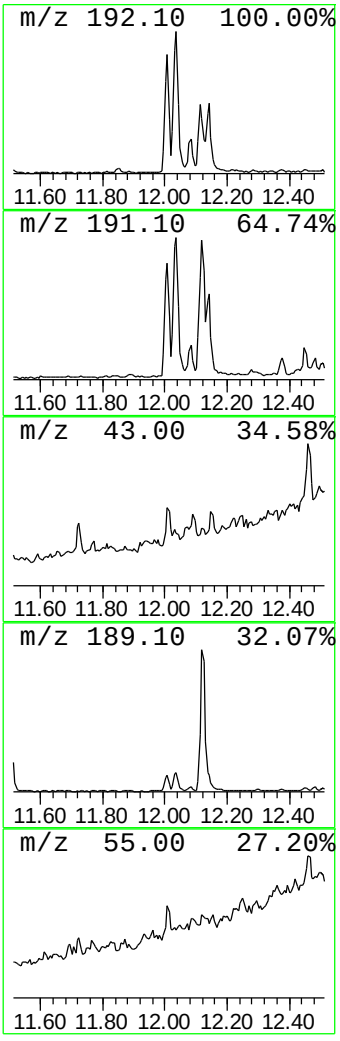
****

Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	4.36 ng	398563	Phenanthrene-d10	11.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	94
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	92





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
Data File : BF107623.D
Acq On : 24 Jul 2018 17:11
Operator : JU/SJ
Sample : J4096-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

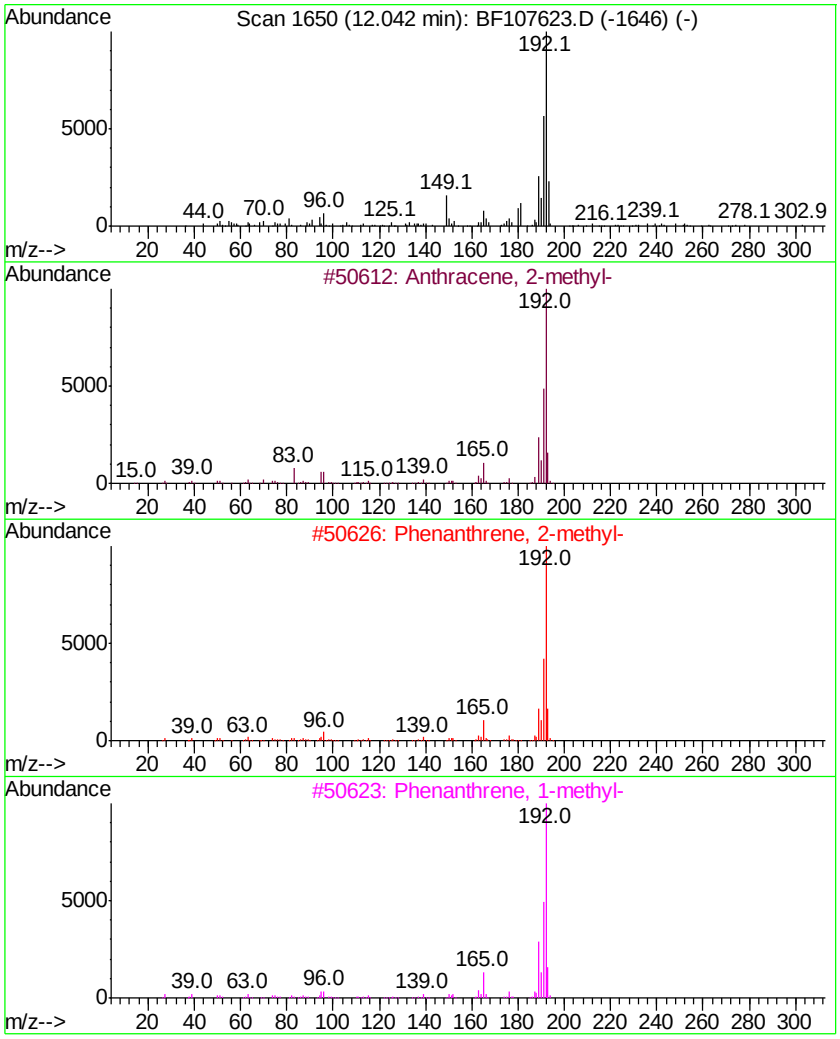
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

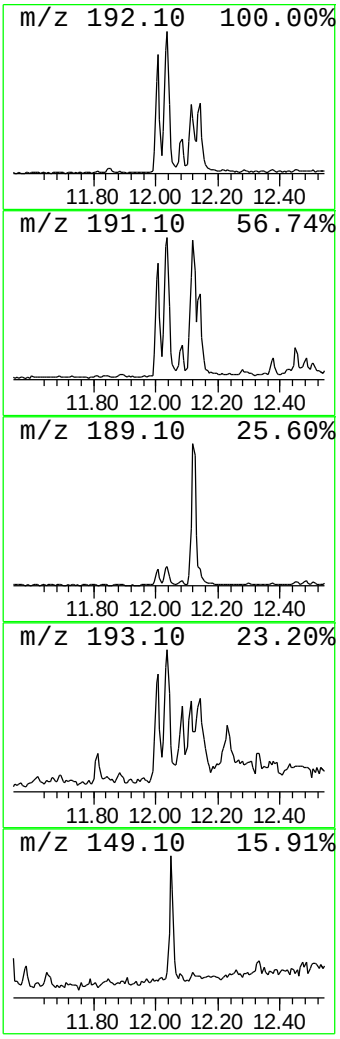
****

Peak Number 7 Anthracene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.04	3.94 ng	360001	Phenanthrene-d10	11.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 2-methyl-	192	C15H12	000613-12-7	97
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	92
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	90





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072418\
Data File : BF107623.D
Acq On : 24 Jul 2018 17:11
Operator : JU/SJ
Sample : J4096-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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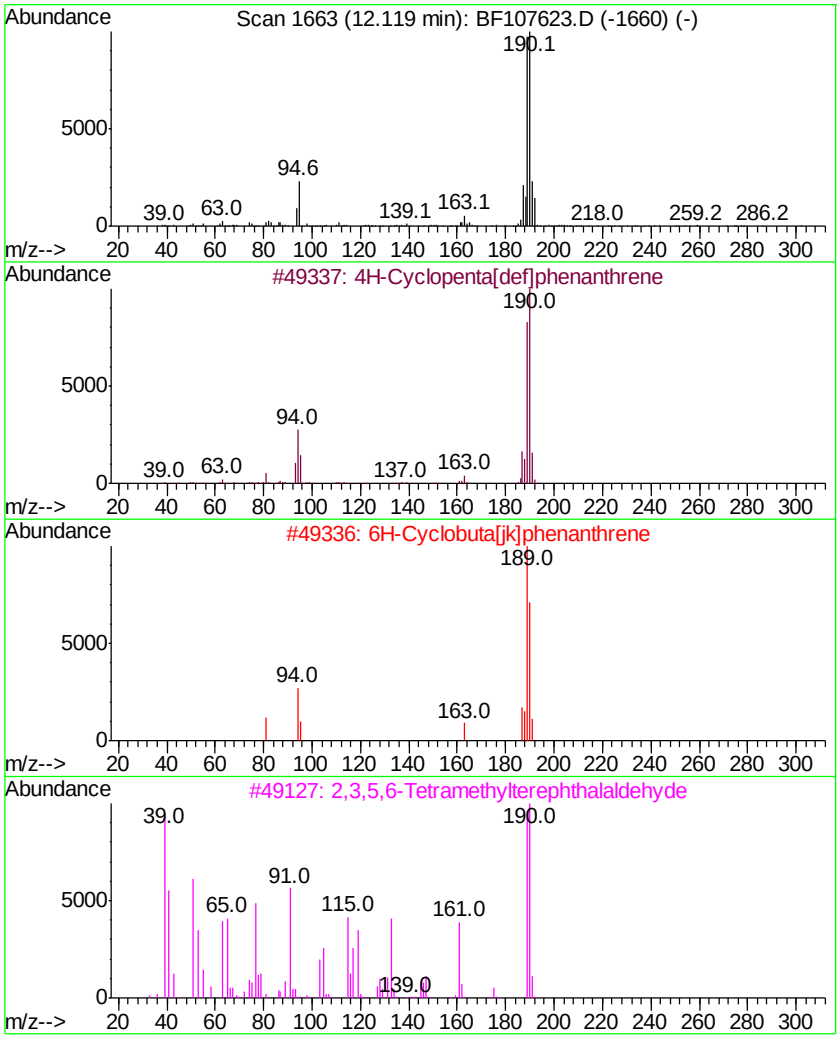
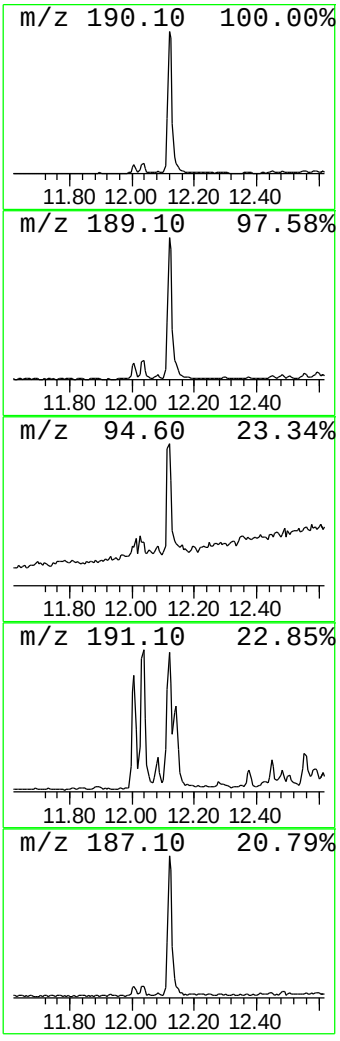
Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.12	12.29 ng	1123010	Phenanthrene-d10	11.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	94
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	50
3	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
5	1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	25





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Operator : JU/SJ
Sample : J4096-08 2X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-10

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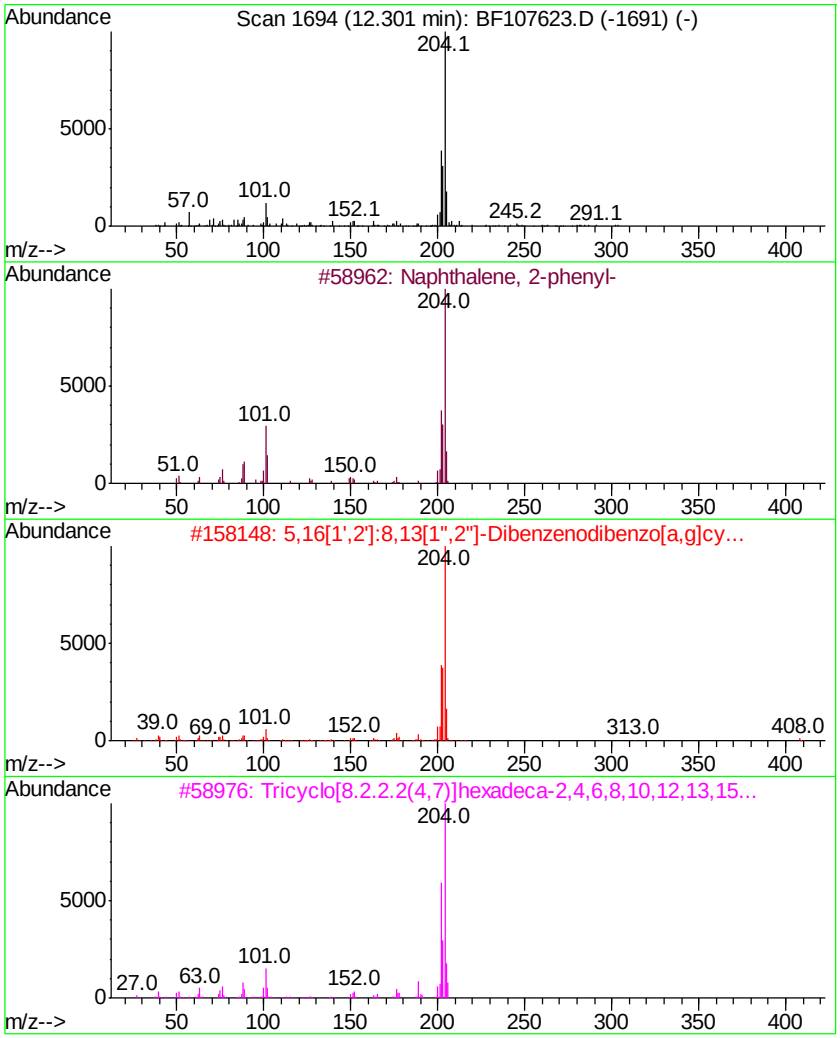
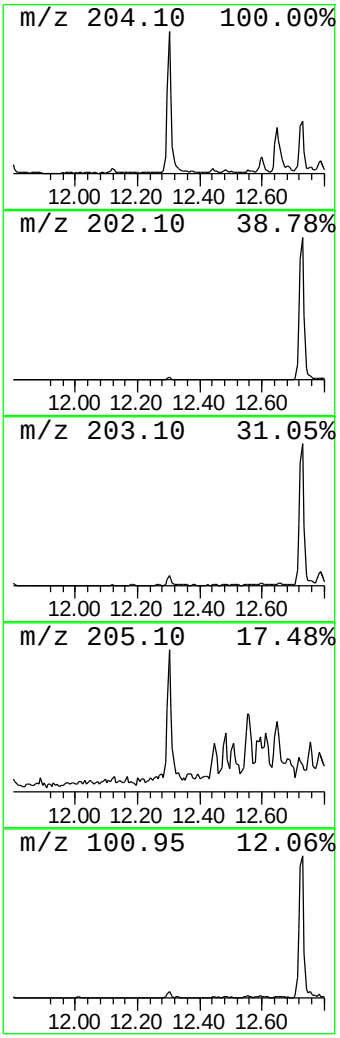
Peak Number 9 Naphthalene, 2-phenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.30	2.71 ng	247241	Phenanthrene-d10	11.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
2	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	80
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	53
4	2-Phenylnaphthalene	204	C16H12	035465-71-5	52
5	Pyrazol-5-ol, 3-(3,4-methylenedi...	204	C10H8N2O3	1000271-76-1	50





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Instrument : BNA_F
ClientSampleId : DW-10

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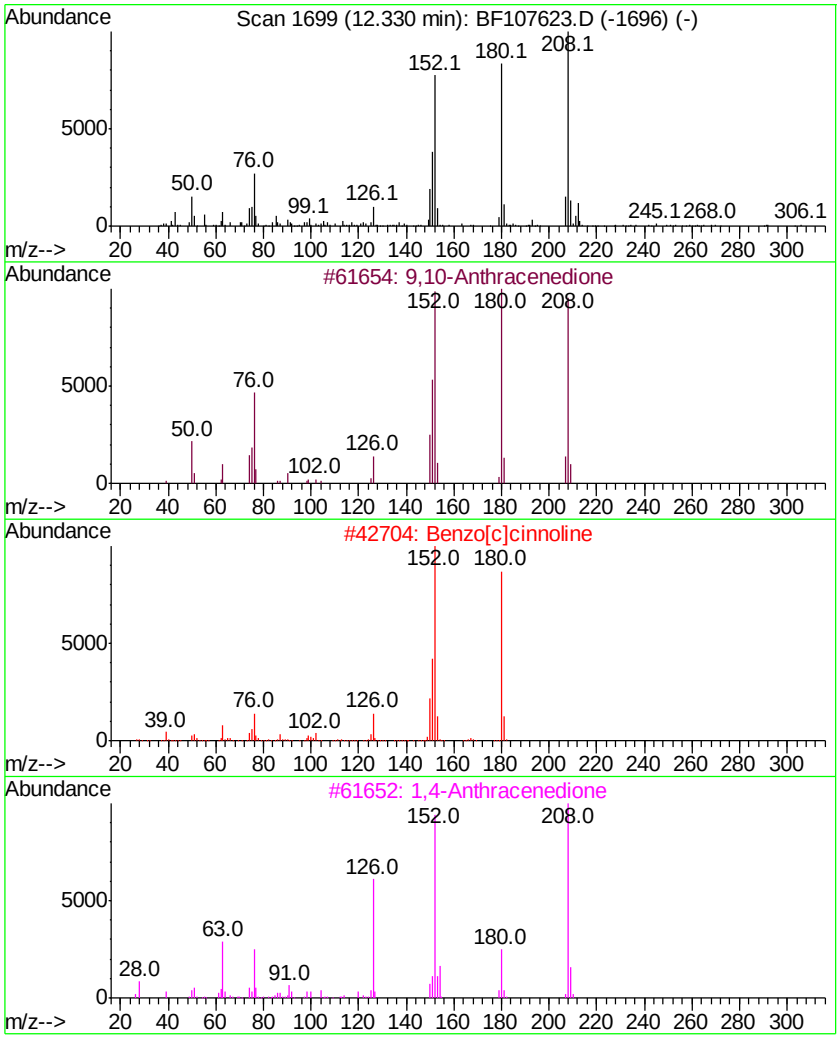
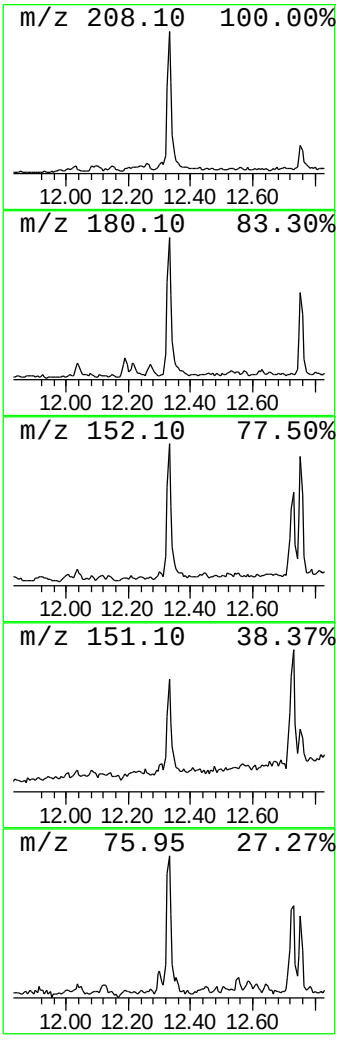
Peak Number 10 9,10-Anthracenedione Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.33	2.88 ng	263614	Phenanthrene-d10	11.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	64
3	1,4-Anthracenedione	208	C14H8O2	000635-12-1	47
4	1H-Phenalen-1-one	180	C13H8O	000548-39-0	46
5	2,5-Cyclohexadiene-1,4-dione, 2,...	208	C6H2Cl2O4	000087-88-7	43





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ClientSampleId : DW-10

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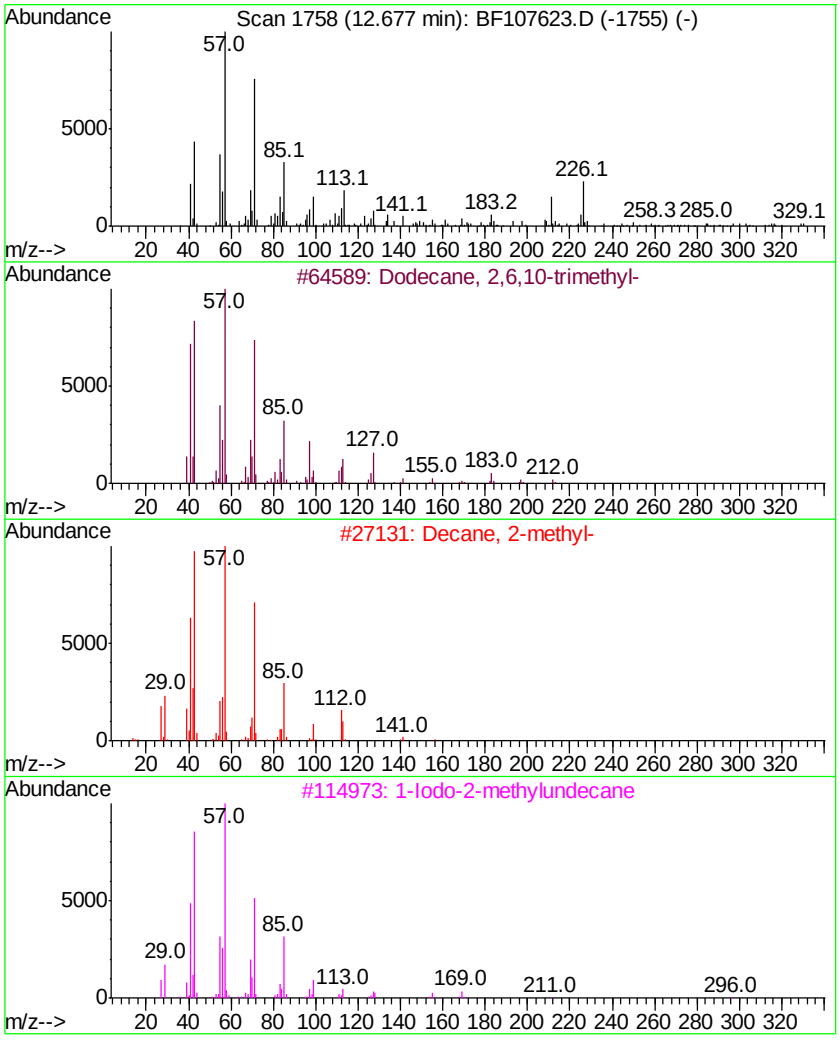
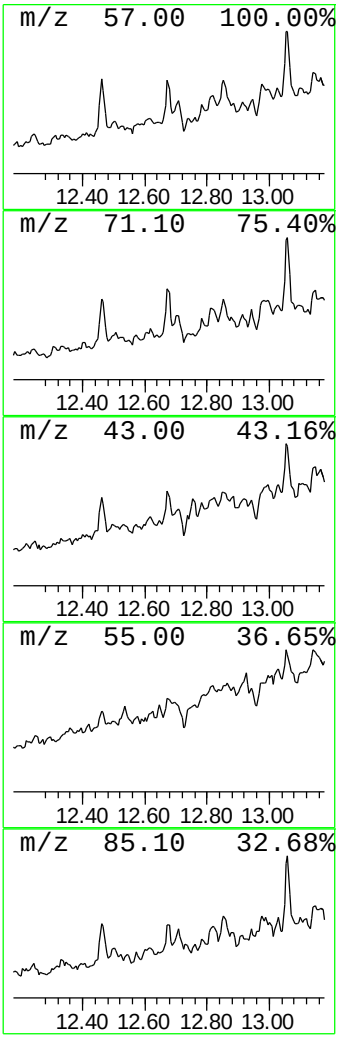
Peak Number 11 Dodecane, 2,6,10-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.68	2.43 ng	222395	Phenanthrene-d10	11.51

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	76
2	Decane, 2-methyl-	156	C11H24	006975-98-0	64
3	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	64
4	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	64
5	Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	64





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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DW-10

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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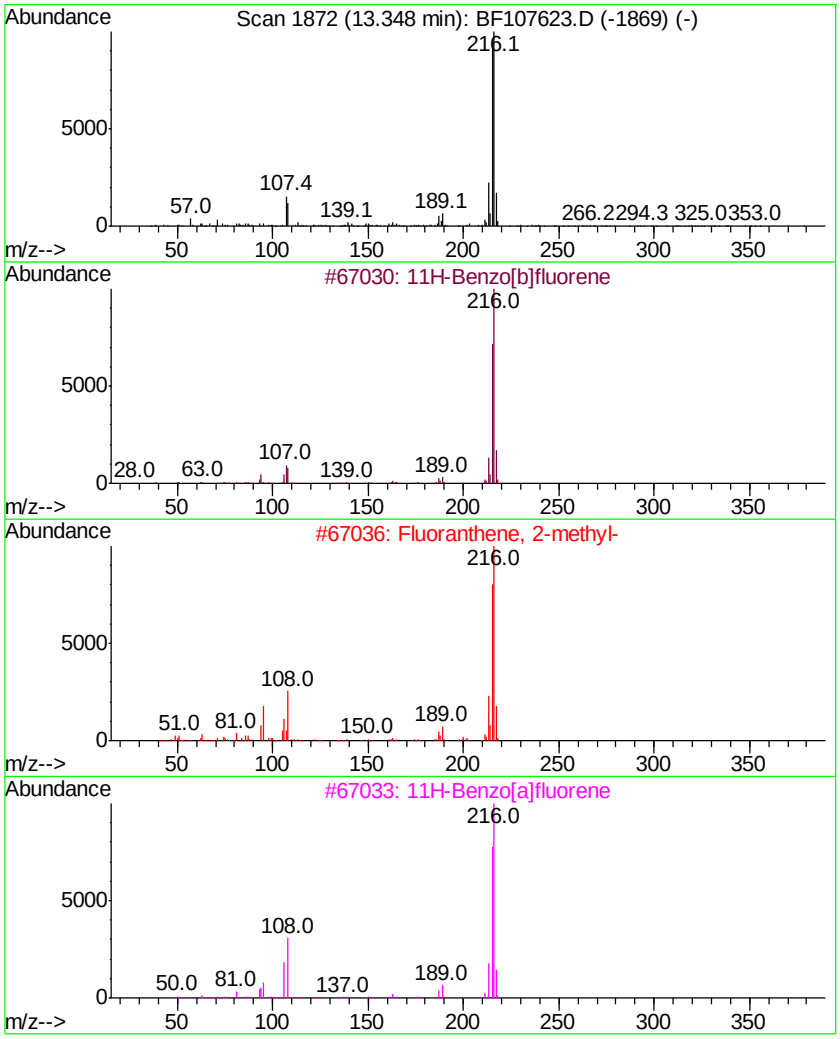
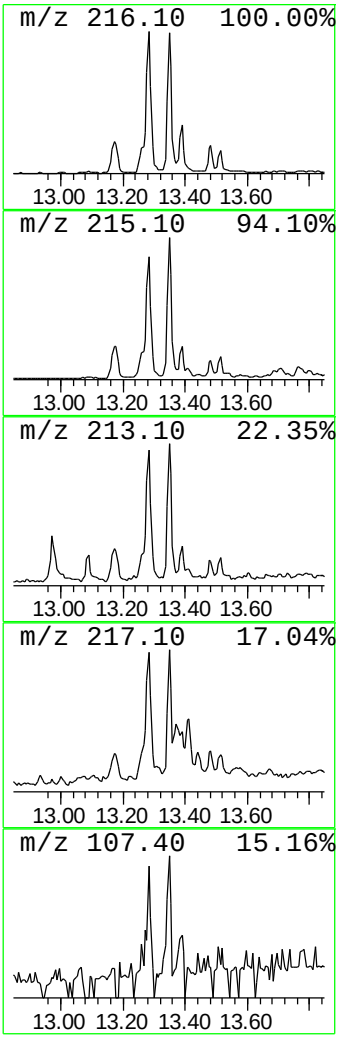
Peak Number 12 11H-Benzo[b]fluorene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.35	3.90 ng	541669	Chrysene-d12	14.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
4	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72
5	Benzoic acid, 3-(3,5-dimethyl-1-...	216	C12H12N2O2	312531-88-7	64





Tentatively Identified Compound (LSC) summary

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Quant Title : ASP BNA STANDARDS FOR 5 POINT 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
2-Pentanone, 4-hy...	5.19	4.8	ng	345243	1	6.97	1453000	20.0
.alpha.-Pinene	6.24	3.8	ng	276358	1	6.97	1453000	20.0
Dodecane, 2,7,10-...	10.91	2.4	ng	219515	4	11.51	1827570	20.0
9H-Fluorene, 9-me...	11.11	2.1	ng	195097	4	11.51	1827570	20.0
Dibenzothiophene	11.40	3.2	ng	289289	4	11.51	1827570	20.0
Phenanthrene, 2-m...	12.01	4.4	ng	398563	4	11.51	1827570	20.0
Anthracene, 2-met...	12.04	3.9	ng	360001	4	11.51	1827570	20.0
4H-Cyclopenta[def...	12.12	12.3	ng	1123010	4	11.51	1827570	20.0
Naphthalene, 2-ph...	12.30	2.7	ng	247241	4	11.51	1827570	20.0
9,10-Anthracenedione	12.33	2.9	ng	263614	4	11.51	1827570	20.0
Dodecane, 2,6,10-...	12.68	2.4	ng	222395	4	11.51	1827570	20.0
11H-Benzo[b]fluorene	13.35	3.9	ng	541669	5	14.16	2780050	20.0