

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078507.D
 Acq On : 14 Apr 2015 19:02
 Operator : TP/IZ
 Sample : G1828-03 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB06A

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

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.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	1.496	26	27	34	rVB2	9924	22392	1.27%	0.263%
2	1.598	34	36	69	rVB	708387	1760850	100.00%	20.653%
3	5.062	337	339	348	rBV	43860	70870	4.02%	0.831%
4	6.410	456	457	465	rVB	59045	80934	4.60%	0.949%
5	6.525	465	467	470	rBV	74678	82924	4.71%	0.973%
6	6.799	489	491	494	rVB	104975	111130	6.31%	1.303%
7	6.982	506	507	510	rBV	42113	57386	3.26%	0.673%
8	7.508	550	553	557	rVB	50037	46761	2.66%	0.548%
9	8.376	627	629	631	rBV	121695	163294	9.27%	1.915%
10	9.725	745	747	750	rBV	84624	104198	5.92%	1.222%
11	10.365	801	803	806	rBV	22820	20849	1.18%	0.245%
12	10.536	816	818	821	rBV	200128	211069	11.99%	2.476%
13	11.508	902	903	906	rVV2	50413	68097	3.87%	0.799%
14	12.365	976	978	979	rBV	224807	236653	13.44%	2.776%
15	12.388	979	980	983	rVB	127254	117365	6.67%	1.377%
16	12.457	983	986	988	rBV	81524	84687	4.81%	0.993%
17	12.982	1030	1032	1034	rBV	13495	21135	1.20%	0.248%
18	13.017	1034	1035	1038	rVB	20703	29058	1.65%	0.341%
19	13.085	1038	1041	1042	rBV	14591	20096	1.14%	0.236%
20	13.120	1042	1044	1045	rBV	37691	45387	2.58%	0.532%
21	13.371	1064	1066	1069	rVV2	12910	30133	1.71%	0.353%
22	13.668	1090	1092	1094	rBV	21633	33080	1.88%	0.388%
23	13.760	1099	1100	1104	rVB2	17303	22700	1.29%	0.266%
24	13.851	1106	1108	1111	rBV	401310	435817	24.75%	5.112%
25	14.034	1122	1124	1127	rBV2	21020	30170	1.71%	0.354%
26	14.125	1129	1132	1135	rBV	449345	461810	26.23%	5.417%
27	14.194	1135	1138	1140	rBV2	14666	18975	1.08%	0.223%
28	14.240	1140	1142	1144	rBV	80344	98282	5.58%	1.153%
29	14.297	1144	1147	1148	rBV	25519	33587	1.91%	0.394%
30	14.354	1148	1152	1154	rBV	145636	182238	10.35%	2.138%
31	14.411	1154	1157	1159	rBV2	28640	52040	2.96%	0.610%
32	14.514	1164	1166	1171	rVB2	51863	136284	7.74%	1.599%
33	14.628	1174	1176	1178	rBV	23172	23881	1.36%	0.280%
34	14.674	1178	1180	1185	rBV2	55250	96211	5.46%	1.128%

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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

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35	14.777	1188	1189	1191	rVB	35777	37758	2.14%	0.443%
36	14.823	1191	1193	1194	rBV	18094	28382	1.61%	0.333%
37	14.994	1205	1208	1209	rBV	16676	21973	1.25%	0.258%
38	15.177	1221	1224	1227	rVB	56944	77597	4.41%	0.910%
39	15.303	1232	1235	1237	rBV	67930	109491	6.22%	1.284%
40	15.360	1237	1240	1242	rVV2	49090	109485	6.22%	1.284%
41	15.440	1245	1247	1249	rVB	39564	57315	3.25%	0.672%
42	15.508	1251	1253	1255	rBV2	23022	33956	1.93%	0.398%
43	15.611	1259	1262	1264	rVV2	359420	678024	38.51%	7.953%
44	15.657	1264	1266	1267	rVV	235359	314625	17.87%	3.690%
45	15.691	1267	1269	1271	rVB3	24652	40796	2.32%	0.479%
46	15.737	1271	1273	1275	rBV2	32224	46271	2.63%	0.543%
47	15.874	1283	1285	1287	rVB2	30408	40616	2.31%	0.476%
48	15.920	1287	1289	1291	rBV2	18955	37718	2.14%	0.442%
49	16.080	1301	1303	1304	rBV	21687	40733	2.31%	0.478%
50	16.114	1304	1306	1308	rVV	50969	84451	4.80%	0.991%
51	16.160	1308	1310	1313	rVV2	26725	44890	2.55%	0.527%
52	16.903	1372	1375	1381	rBV	283899	578590	32.86%	6.786%
53	17.017	1383	1385	1387	rVB	48418	65010	3.69%	0.763%
54	17.211	1400	1402	1406	rBV	136649	222602	12.64%	2.611%
55	17.280	1406	1408	1411	rVB	192705	215212	12.22%	2.524%
56	17.337	1411	1413	1415	rBV	190561	310878	17.65%	3.646%
57	18.583	1519	1522	1526	rVB2	102214	198571	11.28%	2.329%
58	18.926	1549	1552	1554	rBV	83977	120387	6.84%	1.412%

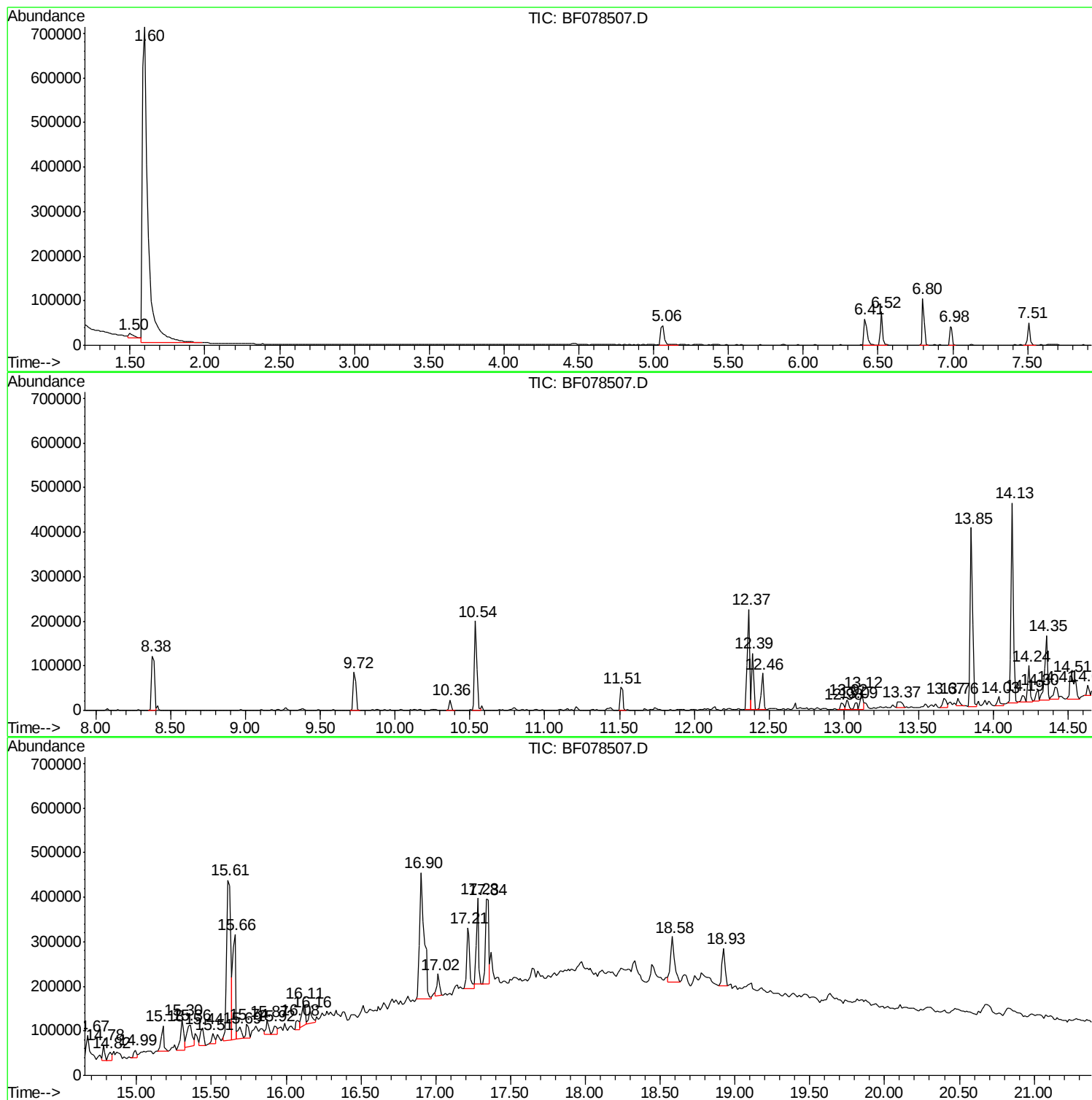
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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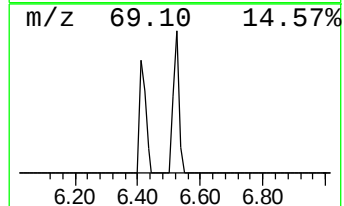
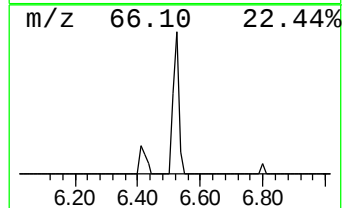
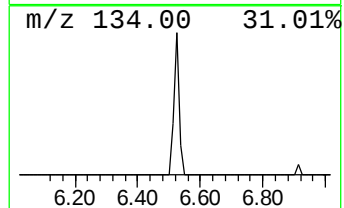
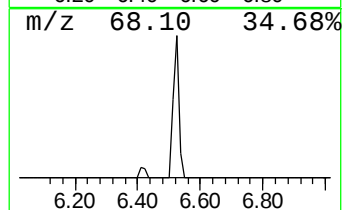
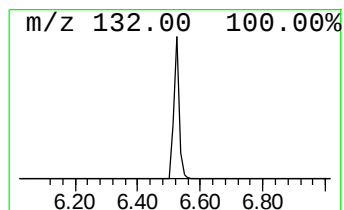
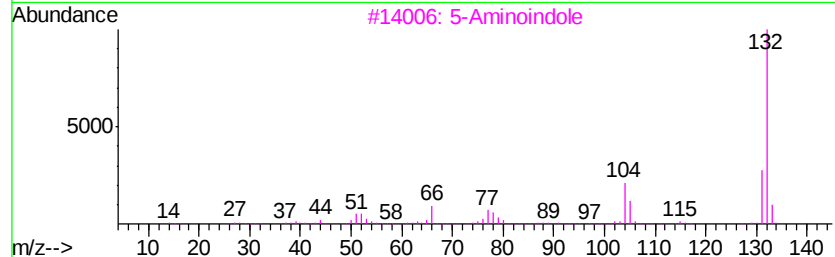
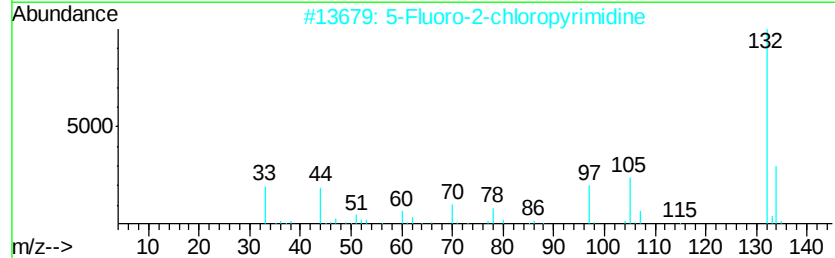
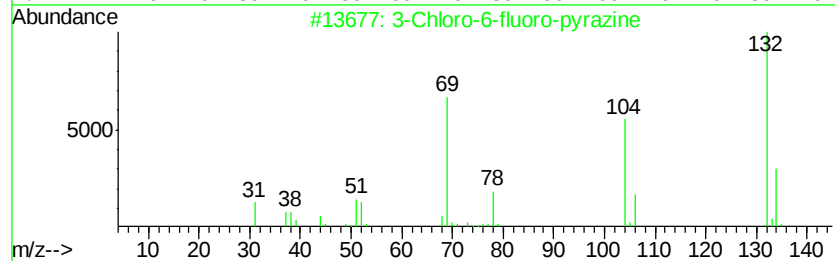
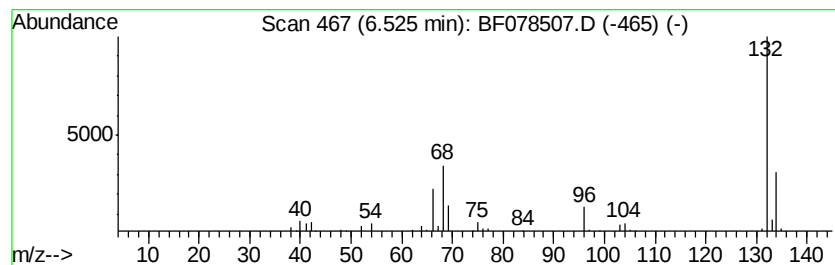
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown6.52 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	14.92 ng	82924	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	25
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	10
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9



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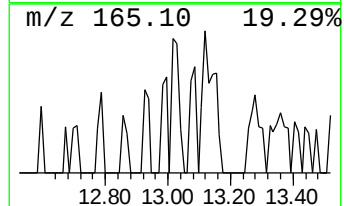
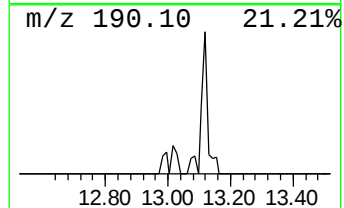
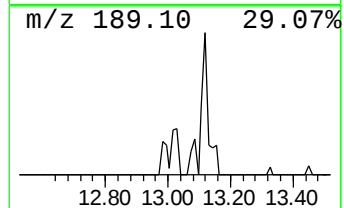
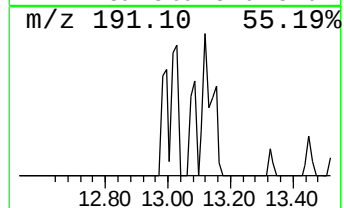
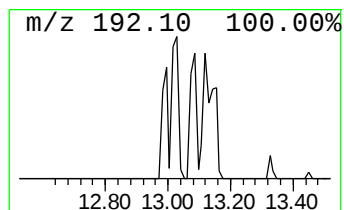
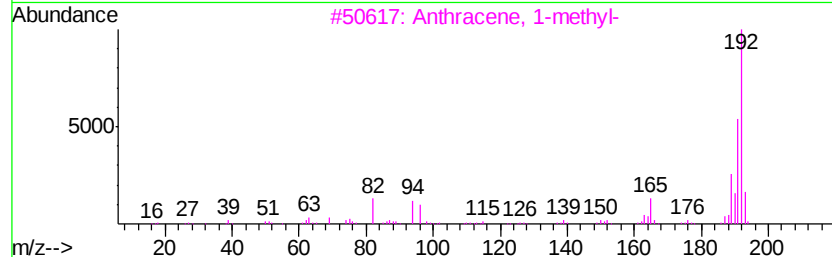
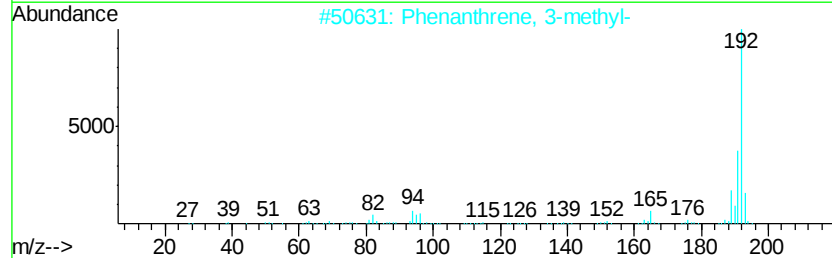
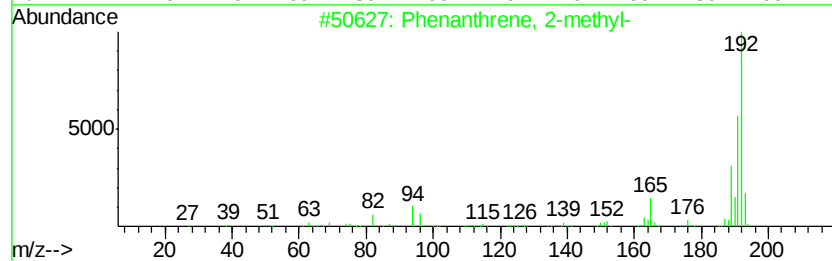
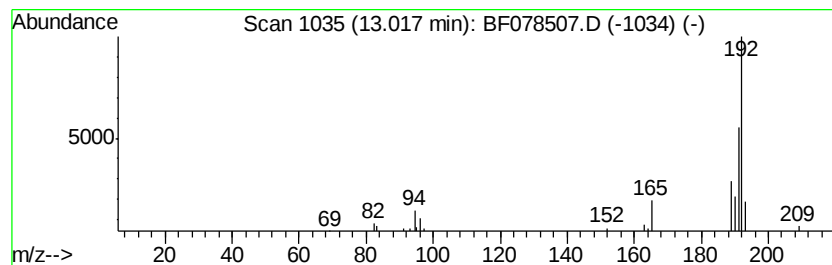
Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.02	2.46 ng	29058	Phenanthrene-d10		12.37
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	87
2	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	87
3	Anthracene, 1-methyl-	192	C15H12	000610-48-0	81
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	81
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	72



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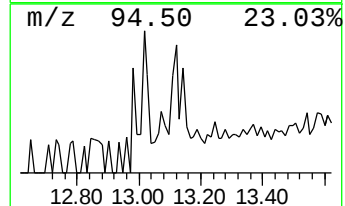
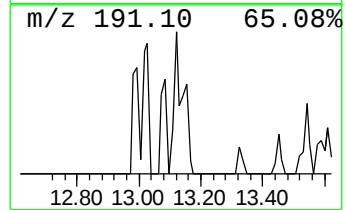
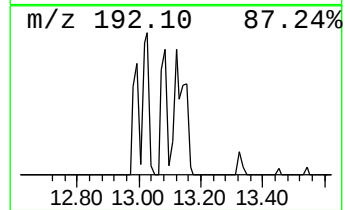
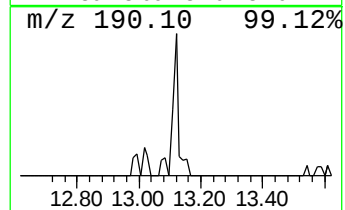
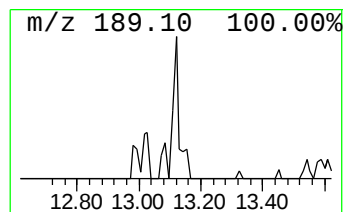
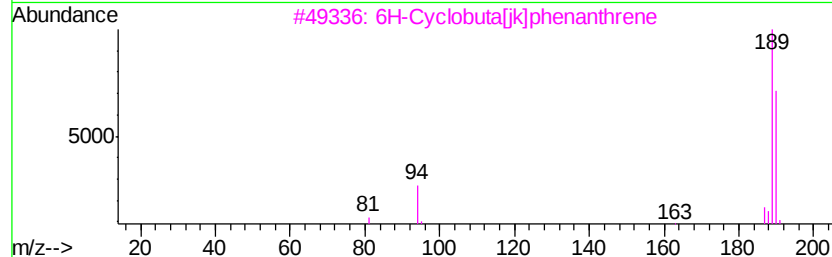
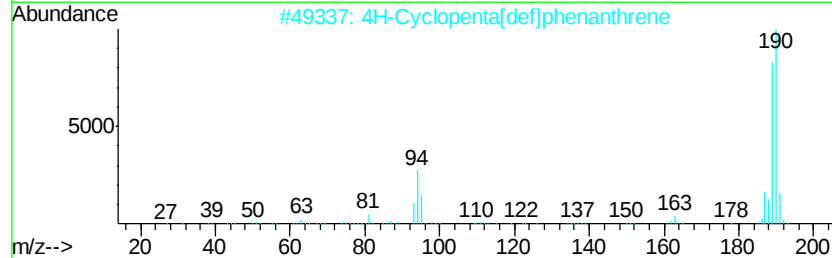
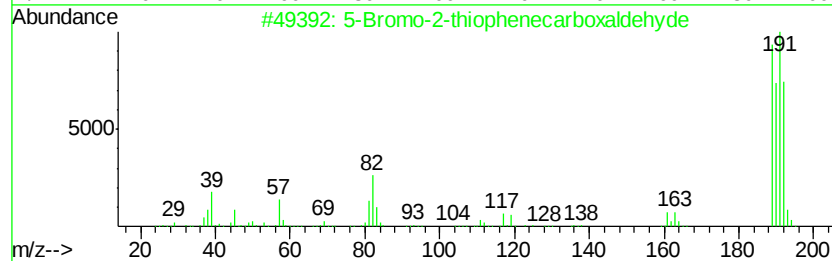
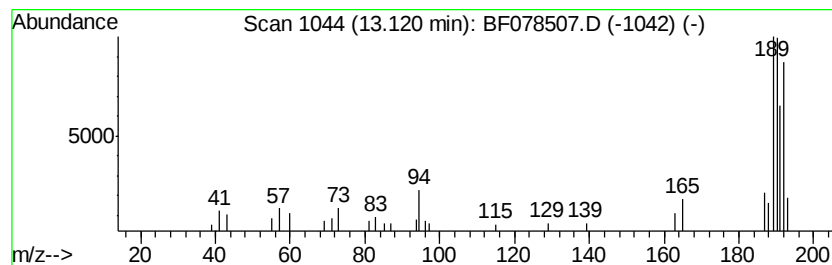
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 5-Bromo-2-thiophenecarboxal... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	3.84 ng	45387	Phenanthrene-d10	12.37

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Bromo-2-thiophenecarboxaldehyde	190	C5H3BrOS	004701-17-1	53
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	49
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	41
5	Methyl diselenide	190	C2H6Se2	007101-31-7	38



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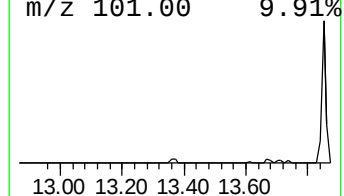
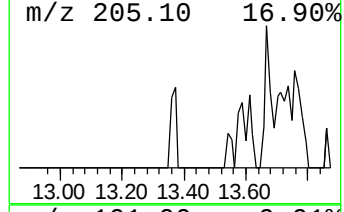
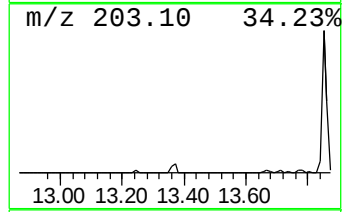
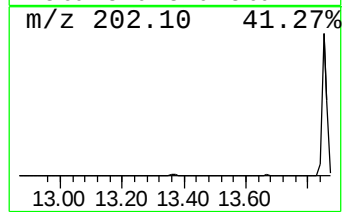
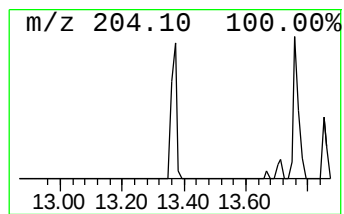
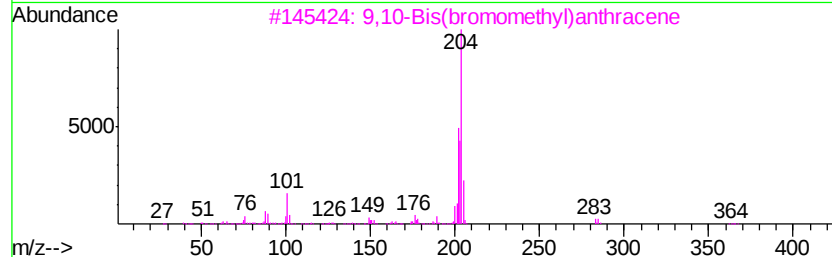
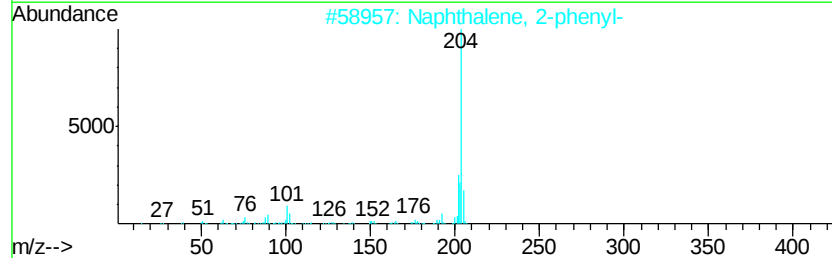
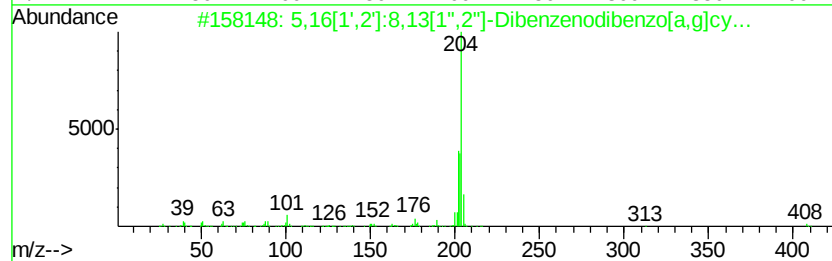
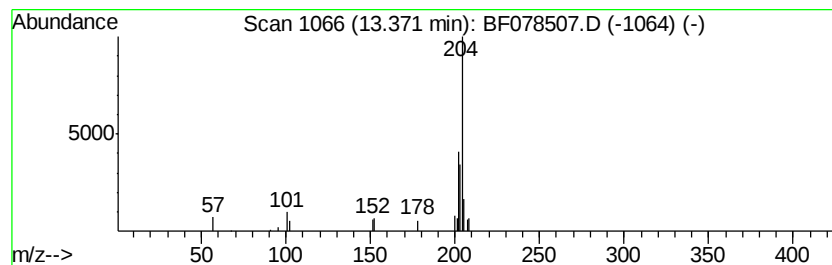
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.37	2.55 ng	30133	Phenanthrene-d10	12.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,16	[1',2']	:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	78
2	Naphthalene,	2-phenyl-		204	C16H12	000612-94-2	53
3	9,10-Bis(bromomethyl)anthracene			362	C16H12Br2	034373-96-1	53
4	Tricyclo[8.2.2.2(4,7)]hexadeca-2...			204	C16H12	006572-60-7	52
5	3,4-Biphenyldicarbonitrile			204	C14H8N2	004128-63-6	50



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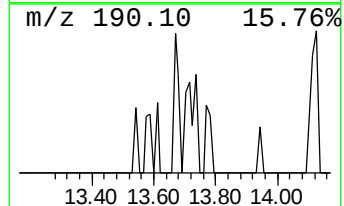
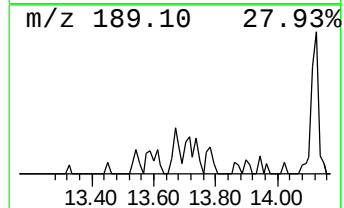
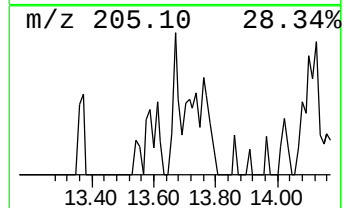
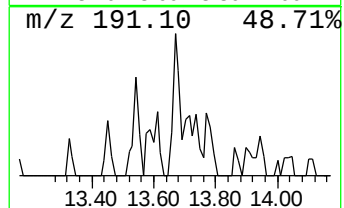
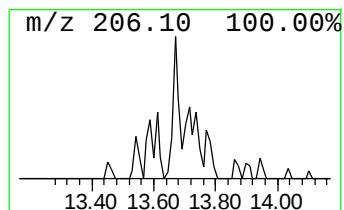
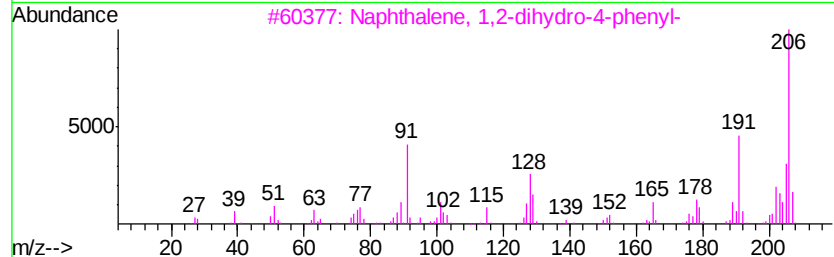
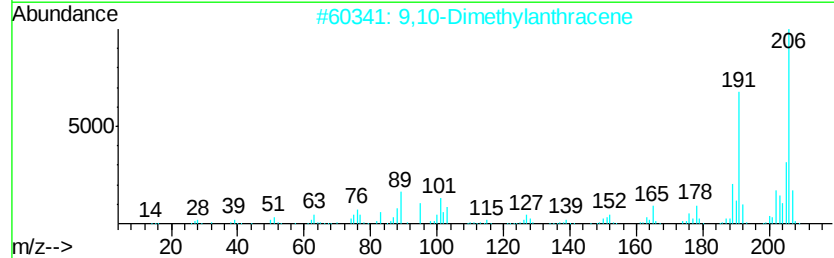
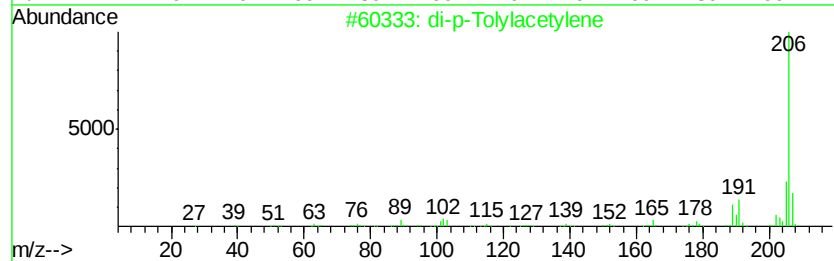
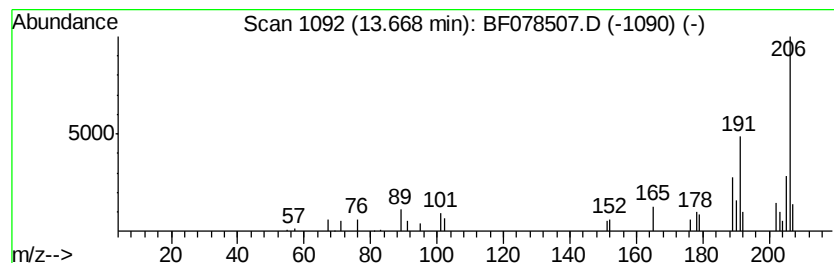
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 di-p-Tolylacetylene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	2.80 ng	33080	Phenanthrene-d10	12.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	di-p-Tolylacetylene	206	C16H14	002789-88-0	93
2		9,10-Dimethylantracene	206	C16H14	000781-43-1	91
3		Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	90
4		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	87
5		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078507.D
Acq On : 14 Apr 2015 19:02
Operator : TP/IZ
Sample : G1828-03 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HS-SB06A

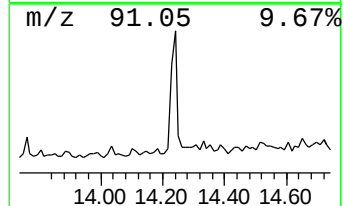
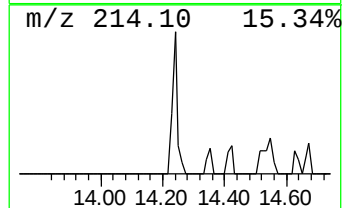
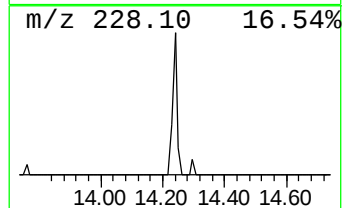
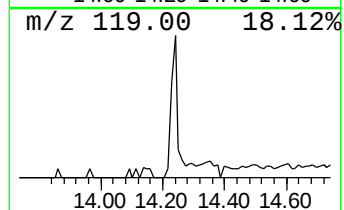
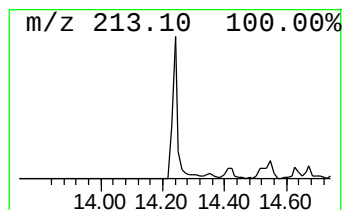
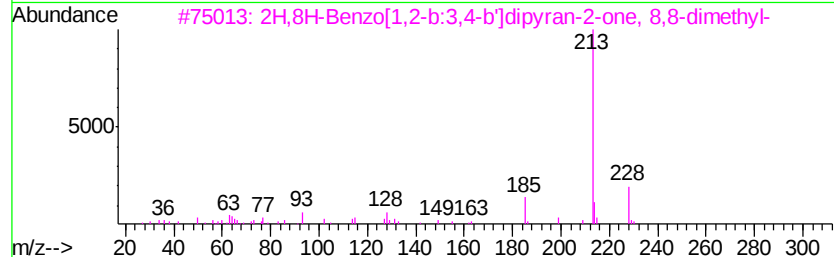
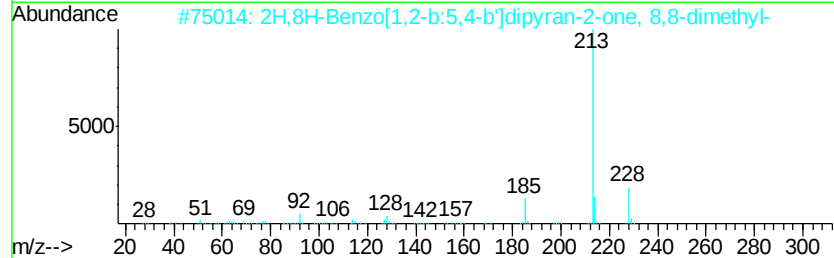
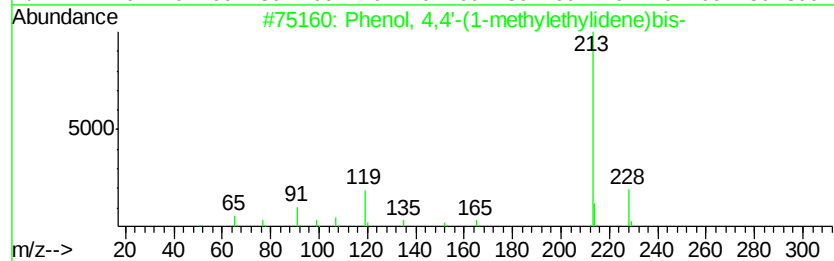
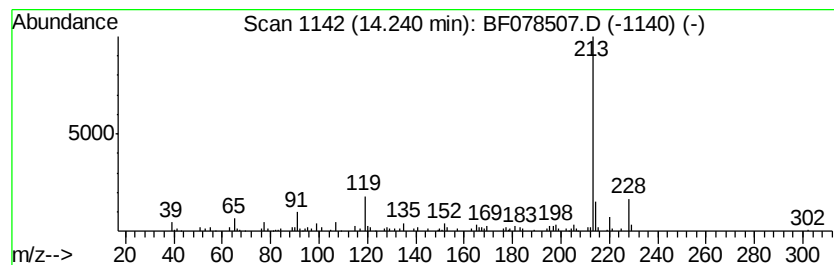
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenol, 4,4'-(1-methylethyl... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.24	2.90 ng	98282	Chrysene-d12	15.62

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	98
2		2H,8H-Benzo[1,2-b:5,4-b']dipyran...	228	C14H12O3	000553-19-5	64
3		2H,8H-Benzo[1,2-b:3,4-b']dipyran...	228	C14H12O3	000523-59-1	64
4		2,2-Bis-(4-hydroxyphenyl)-butane	242	C16H18O2	000077-40-7	59
5		Diphenolic acid	286	C17H18O4	000126-00-1	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078507.D
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Operator : TP/IZ
Sample : G1828-03 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

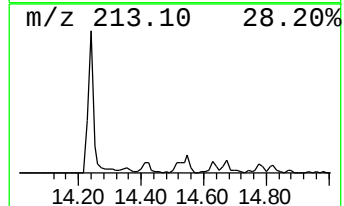
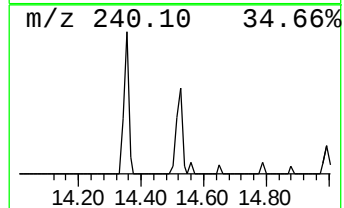
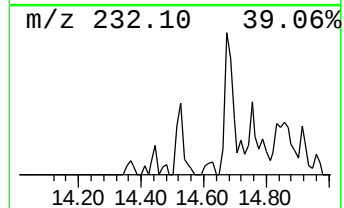
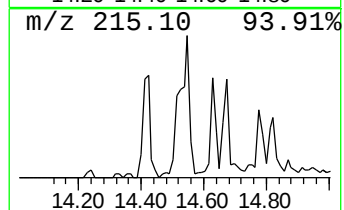
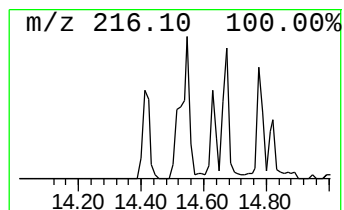
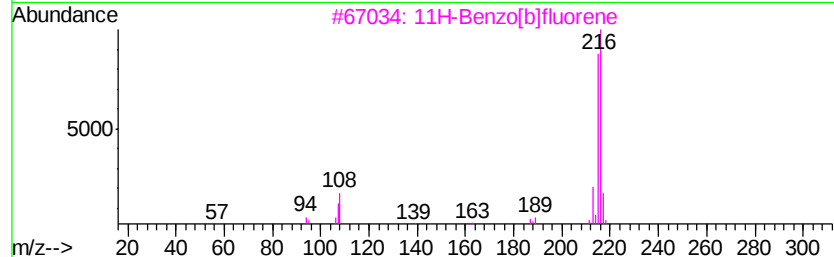
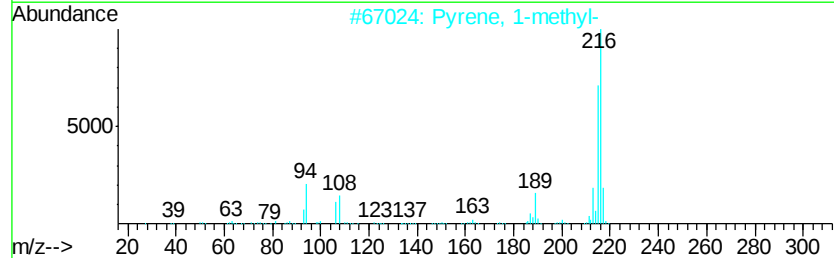
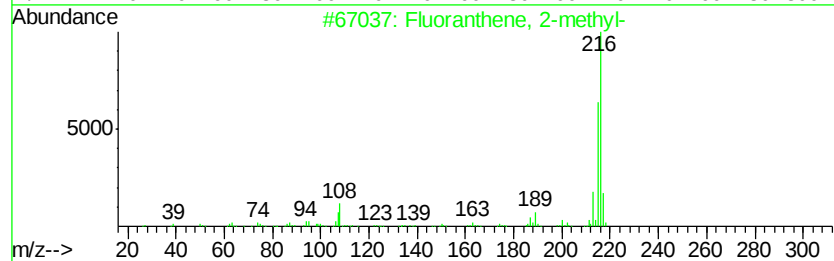
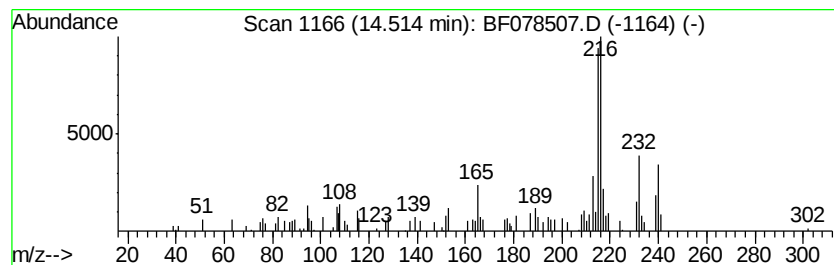
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Fluoranthene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.51	4.02 ng	136284	Chrysene-d12	15.62	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	60
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	53
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	49
4	3-Pyridinamine, 6-chloro-N-(phen...	216	C12H9ClN2	005325-71-3	47
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
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Acq On : 14 Apr 2015 19:02
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Sample : G1828-03 5X
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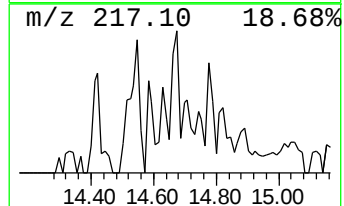
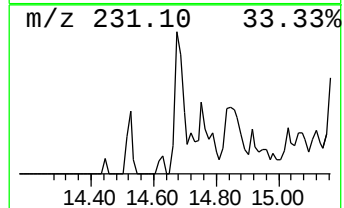
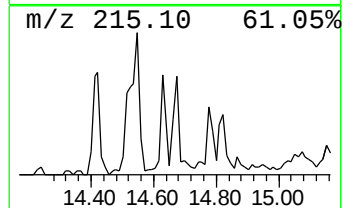
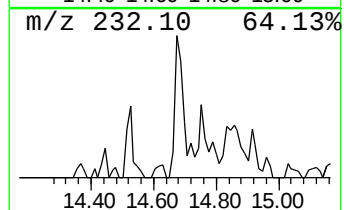
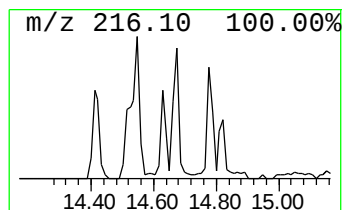
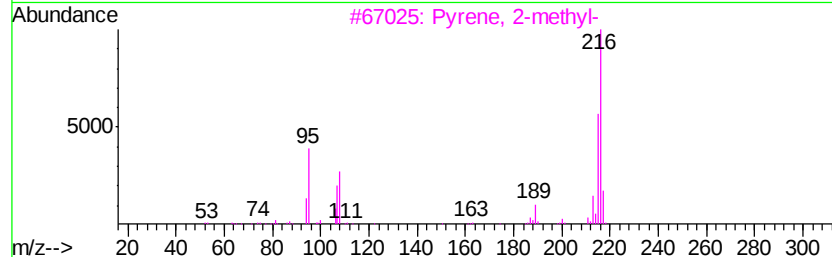
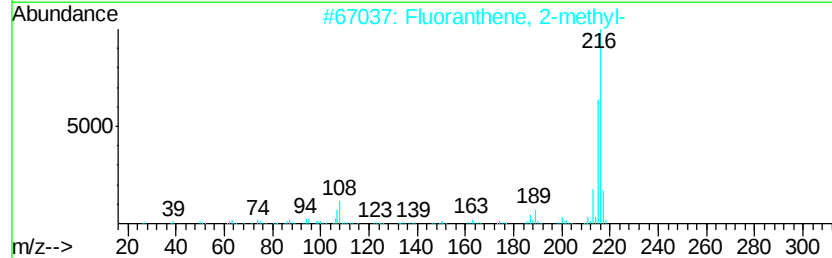
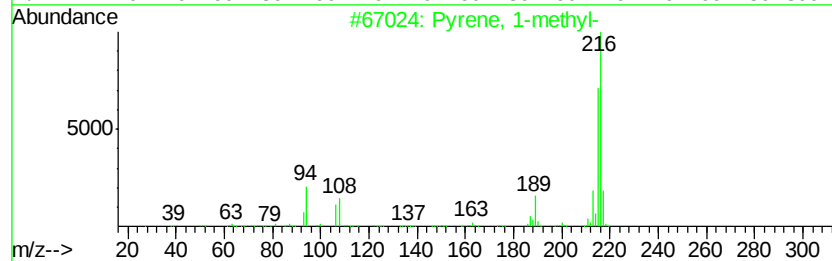
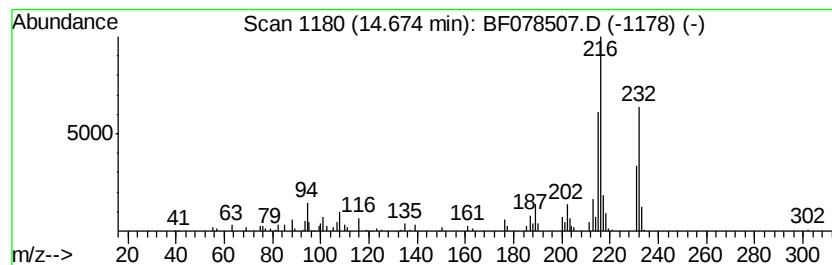
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Pyrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.67	2.84 ng	96211	Chrysene-d12	15.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Pyrene, 1-methyl-	216 C17H12	002381-21-7	91
2	Fluoranthene, 2-methyl-	216 C17H12	033543-31-6	70
3	Pyrene, 2-methyl-	216 C17H12	003442-78-2	60
4	Pyrene, 4-methyl-	216 C17H12	003353-12-6	55
5	11H-Benzo[a]fluorene	216 C17H12	000238-84-6	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078507.D
Acq On : 14 Apr 2015 19:02
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Sample : G1828-03 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

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ClientSampleId :
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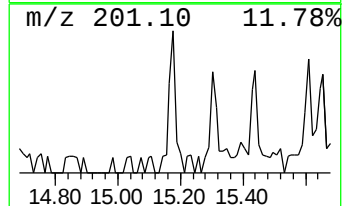
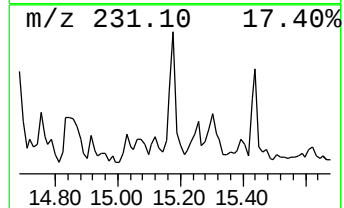
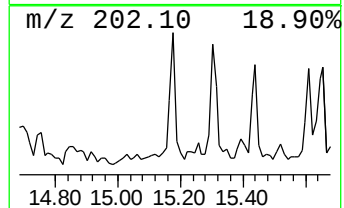
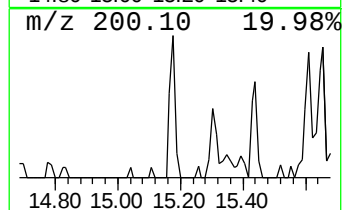
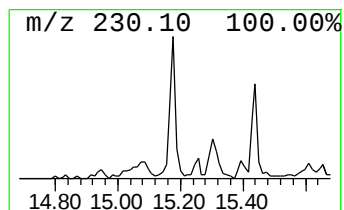
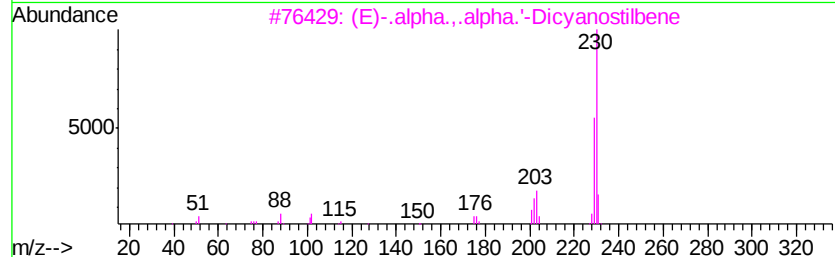
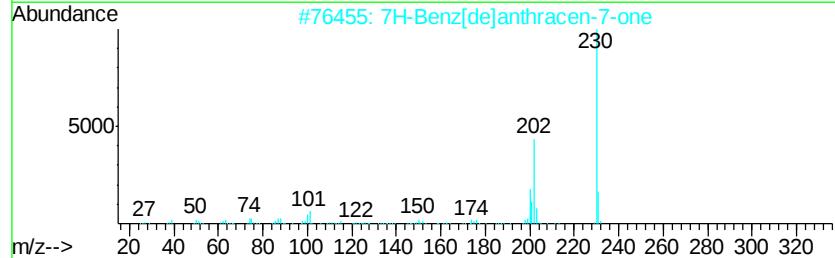
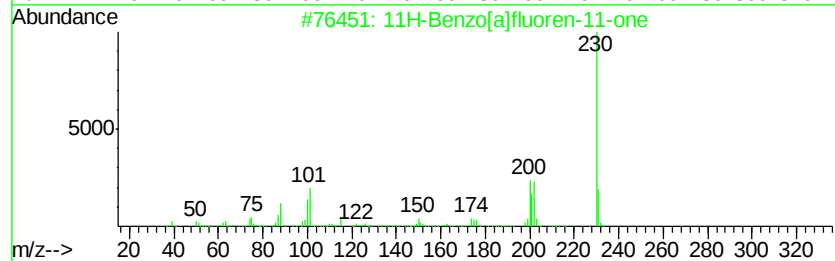
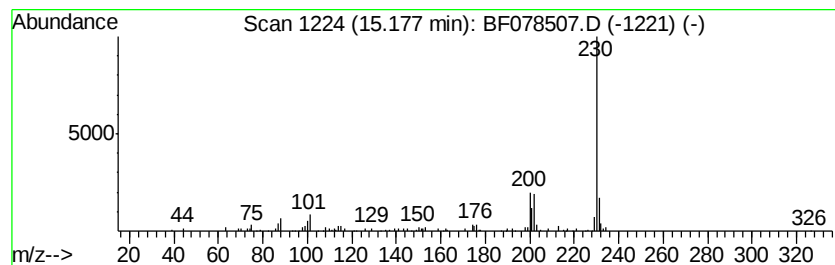
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 11H-Benzo[a]fluoren-11-one Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.18	2.29 ng	77597	Chrysene-d12	15.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	96
2		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	94
3		(E)-.alpha.,.alpha.'-Dicyanostil...	230	C16H10N2	002450-55-7	72
4		1-Pyrene-carboxaldehyde	230	C17H10O	003029-19-4	64
5		Dibenzo[c,h][2,6]naphthyridine	230	C16H10N2	000218-30-4	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078507.D
Acq On : 14 Apr 2015 19:02
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Sample : G1828-03 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

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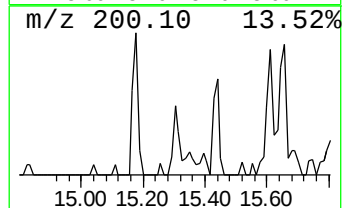
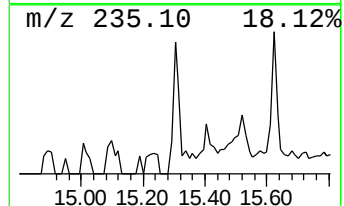
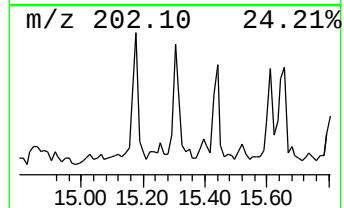
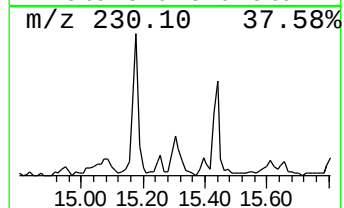
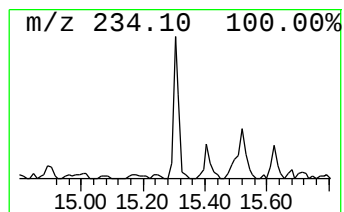
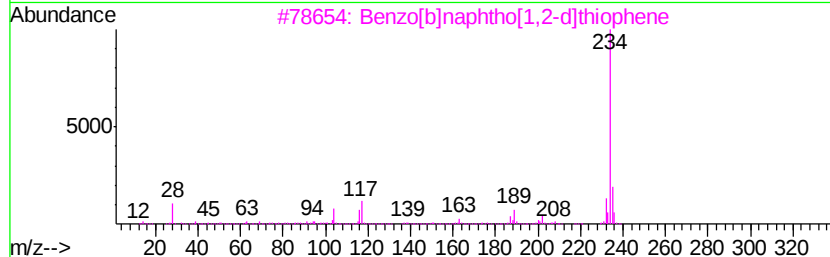
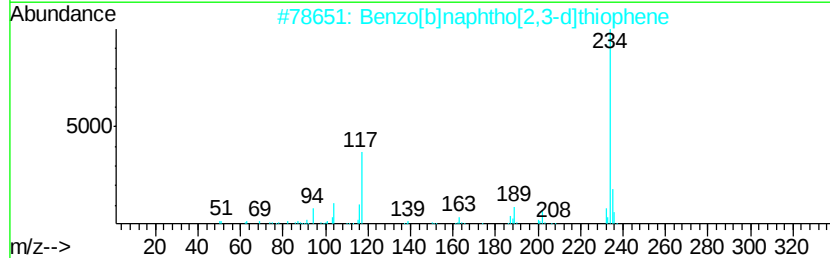
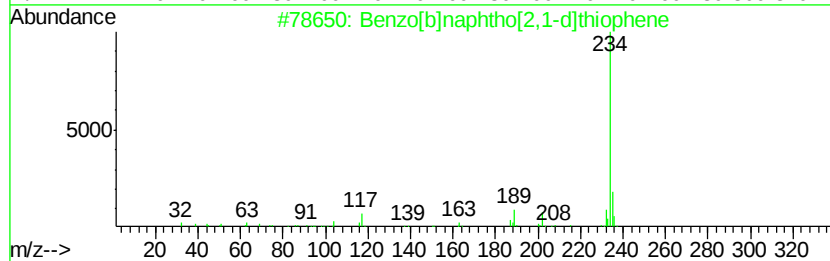
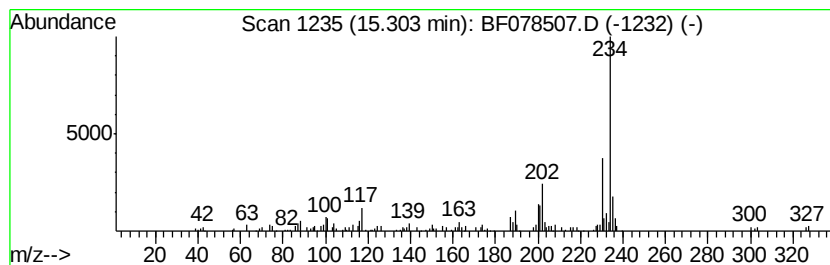
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.30	3.23 ng	109491	Chrysene-d12	15.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	70
2		Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	64
3		Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	49
4		Anthra[1,9a,9-cd;5,10a,10-c'd']d...	234	C14H6N2O2	000201-91-2	46
5		Imidazole, 4-pentafluorophenyl-	234	C9H3F5N2	081769-58-6	46



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
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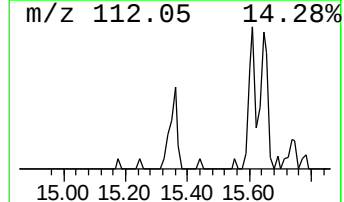
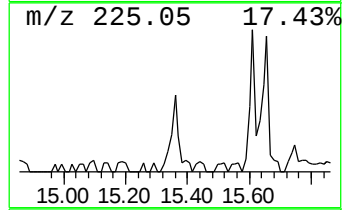
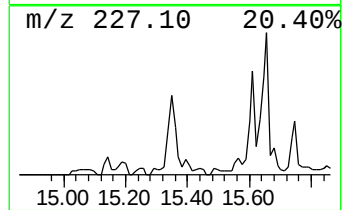
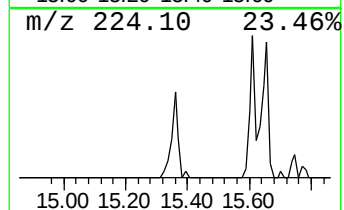
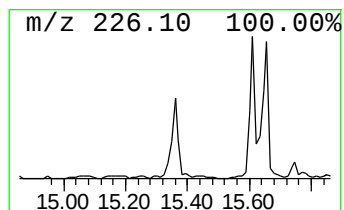
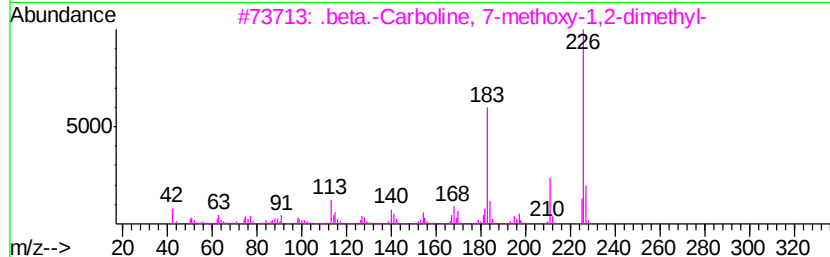
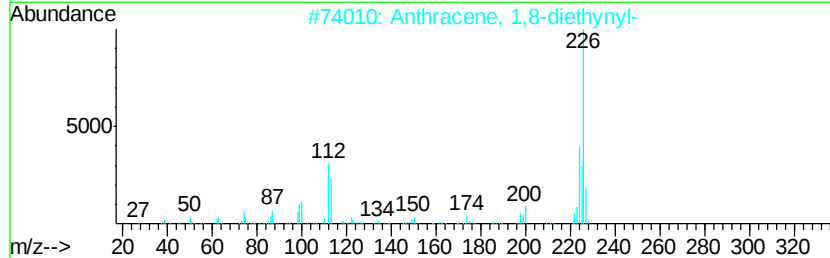
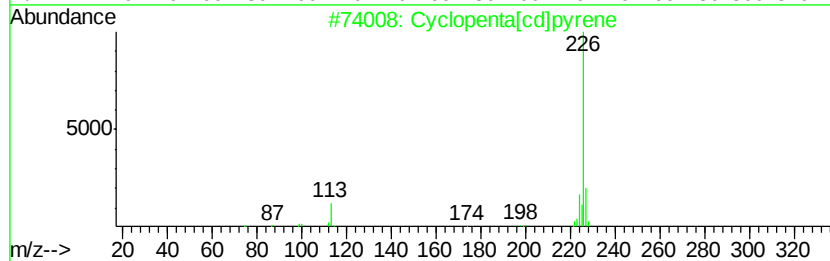
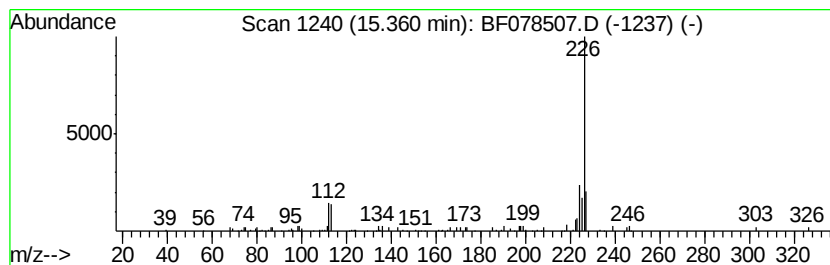
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Cyclopenta[cd]pyrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.		
15.36	3.23 ng	109485	Chrysene-d12	15.62		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	76
2		Anthracene, 1,8-diethynyl-	226	C18H10	078053-58-4	59
3		.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	58
4		Benzo[ghi]fluoranthene	226	C18H10	000203-12-3	43
5		2-Furfurylidene-7-hydroxy-1-inda...	226	C14H10O3	1000244-80-4	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078507.D
Acq On : 14 Apr 2015 19:02
Operator : TP/IZ
Sample : G1828-03 5X
Misc :
ALS Vial : 5 Sample Multiplier: 1

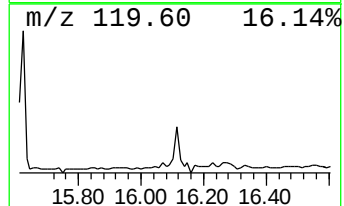
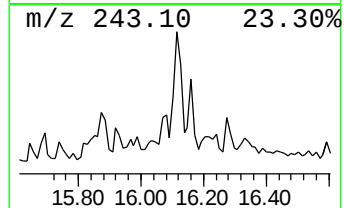
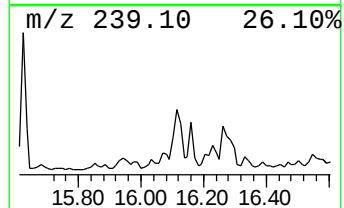
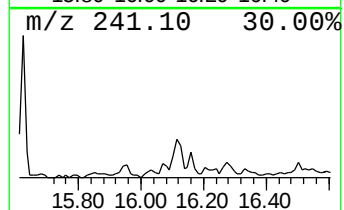
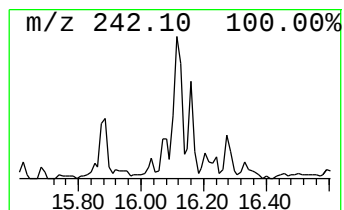
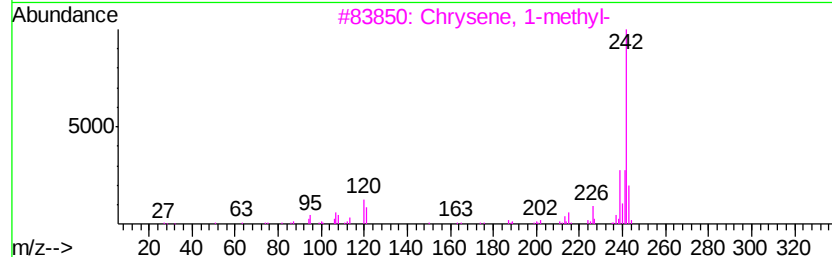
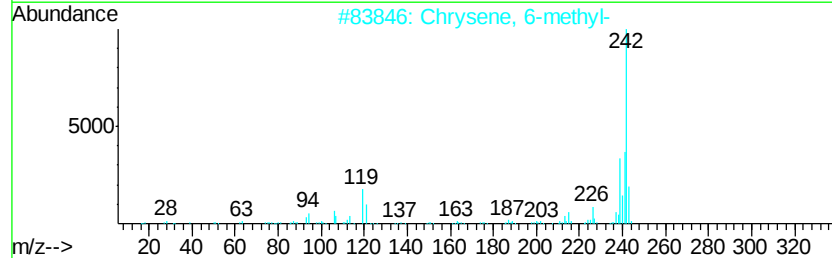
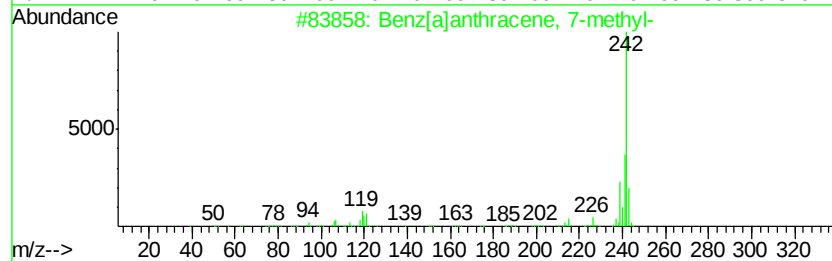
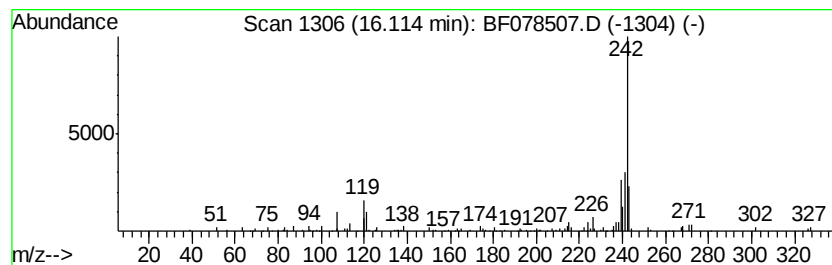
Instrument :
BNA_F
ClientSampleId :
HS-SB06A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Benz[a]anthracene, 7-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.11	2.49 ng	84451	Chrysene-d12		15.62	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	94
2		Chrysene, 6-methyl-	242	C19H14	001705-85-7	89
3		Chrysene, 1-methyl-	242	C19H14	003351-28-8	81
4		Benz[a]anthracene, 8-methyl-	242	C19H14	002381-31-9	81
5		Chrysene, 2-methyl-	242	C19H14	003351-32-4	81



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078507.D
 Acq On : 14 Apr 2015 19:02
 Operator : TP/IZ
 Sample : G1828-03 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB06A

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.52	6.52	14.9	ng	82924	1	6.80	111130	20.0
Phenanthrene, 2-m...	13.02	2.5	ng	29058	4	12.37	236653	20.0
5-Bromo-2-thiophe...	13.12	3.8	ng	45387	4	12.37	236653	20.0
5,16[1',2']:8,13[...	13.37	2.5	ng	30133	4	12.37	236653	20.0
di-p-Tolylacetylene	13.67	2.8	ng	33080	4	12.37	236653	20.0
Phenol, 4,4'-(1-m...	14.24	2.9	ng	98282	5	15.62	678024	20.0
Fluoranthene, 2-m...	14.51	4.0	ng	136284	5	15.62	678024	20.0
Pyrene, 1-methyl-	14.67	2.8	ng	96211	5	15.62	678024	20.0
11H-Benzo[a]fluor...	15.18	2.3	ng	77597	5	15.62	678024	20.0
Benzo[b]naphtho[2...	15.30	3.2	ng	109491	5	15.62	678024	20.0
Cyclopenta[cd]pyrene	15.36	3.2	ng	109485	5	15.62	678024	20.0
Benz[a]anthracene...	16.11	2.5	ng	84451	5	15.62	678024	20.0