

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078517.D
 Acq On : 14 Apr 2015 23:49
 Operator : TP/IZ
 Sample : G1828-14
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB04A

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	rVV	51328	87242	6.28%	0.840%
2	1.598	34	36	66	rVB	567073	1389212	100.00%	13.372%
3	4.456	283	286	293	rVB	27803	39404	2.84%	0.379%
4	5.050	336	338	350	rVB	238226	376557	27.11%	3.625%
5	6.410	455	457	464	rVB	362873	420409	30.26%	4.047%
6	6.525	464	467	469	rBV	374865	436609	31.43%	4.203%
7	6.799	489	491	494	rVB	93473	113933	8.20%	1.097%
8	6.993	505	508	510	rBV	261210	303632	21.86%	2.923%
9	7.508	550	553	555	rBV	286477	255298	18.38%	2.457%
10	8.388	627	630	632	rBV	138914	173584	12.50%	1.671%
11	9.736	745	748	750	rBV	443072	510194	36.73%	4.911%
12	10.251	791	793	794	rBV	18276	20233	1.46%	0.195%
13	10.365	801	803	806	rBV	21937	20721	1.49%	0.199%
14	10.536	816	818	821	rBV	175857	208020	14.97%	2.002%
15	11.439	894	897	900	rBV	21704	33745	2.43%	0.325%
16	11.519	901	904	906	rBV	319239	350527	25.23%	3.374%
17	12.365	976	978	979	rBV	242966	228520	16.45%	2.200%
18	12.388	979	980	983	rVB	106244	127535	9.18%	1.228%
19	12.457	983	986	988	rBV	50156	49518	3.56%	0.477%
20	12.994	1030	1033	1034	rBV	20958	22310	1.61%	0.215%
21	13.028	1034	1036	1038	rVB	27477	31044	2.23%	0.299%
22	13.085	1038	1041	1042	rBV	12271	14041	1.01%	0.135%
23	13.120	1042	1044	1046	rVV	51571	60539	4.36%	0.583%
24	13.371	1064	1066	1070	rVV2	16726	30808	2.22%	0.297%
25	13.668	1090	1092	1094	rBV	12997	21672	1.56%	0.209%
26	13.760	1099	1100	1104	rVV2	8534	15218	1.10%	0.146%
27	13.862	1106	1109	1111	rBV	277992	392431	28.25%	3.777%
28	14.034	1122	1124	1126	rBV2	16277	20889	1.50%	0.201%
29	14.125	1130	1132	1135	rBV	309173	398963	28.72%	3.840%
30	14.240	1140	1142	1144	rVB	57029	58674	4.22%	0.565%
31	14.297	1144	1147	1148	rBV	20357	26021	1.87%	0.250%
32	14.354	1149	1152	1154	rVB	612424	648816	46.70%	6.245%
33	14.422	1154	1158	1160	rBV3	17884	31826	2.29%	0.306%
34	14.525	1164	1167	1168	rBV	19127	34726	2.50%	0.334%

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35	14.548	1168	1169	1171	rVB	42457	43380	3.12%	0.418%
36	14.628	1174	1176	1178	rVB	21822	28644	2.06%	0.276%
37	14.674	1178	1180	1182	rBV2	34733	50213	3.61%	0.483%
38	14.788	1185	1190	1191	rBV2	13553	23240	1.67%	0.224%
39	15.085	1210	1216	1218	rBV3	10980	32466	2.34%	0.312%
40	15.177	1220	1224	1228	rBV	37946	70335	5.06%	0.677%
41	15.280	1231	1233	1234	rVV	33262	38270	2.75%	0.368%
42	15.314	1234	1236	1237	rVV	38083	51683	3.72%	0.497%
43	15.348	1237	1239	1242	rVB2	32857	54103	3.89%	0.521%
44	15.394	1242	1243	1245	rBV2	17038	24199	1.74%	0.233%
45	15.440	1245	1247	1250	rVB	36818	44704	3.22%	0.430%
46	15.543	1251	1256	1259	rBV4	71748	130653	9.40%	1.258%
47	15.623	1259	1263	1264	rVV2	336293	583387	41.99%	5.615%
48	15.657	1264	1266	1268	rVV	160471	211659	15.24%	2.037%
49	15.748	1271	1274	1275	rBV	37514	45509	3.28%	0.438%
50	15.794	1276	1278	1280	rBV3	25698	33845	2.44%	0.326%
51	15.874	1283	1285	1287	rVB2	23376	32643	2.35%	0.314%
52	15.920	1287	1289	1291	rBV	20911	36669	2.64%	0.353%
53	16.114	1303	1306	1308	rBV	39565	51344	3.70%	0.494%
54	16.160	1308	1310	1312	rVB	14985	22936	1.65%	0.221%
55	16.228	1313	1316	1317	rBV2	19042	32626	2.35%	0.314%
56	16.503	1337	1340	1343	rBV3	19571	37404	2.69%	0.360%
57	16.880	1371	1373	1378	rBV	228967	489795	35.26%	4.714%
58	16.994	1381	1383	1385	rVB	48424	62407	4.49%	0.601%
59	17.188	1398	1400	1403	rBV	138518	175532	12.64%	1.690%
60	17.257	1403	1406	1408	rBV	166366	239712	17.26%	2.307%
61	17.314	1408	1411	1413	rBV	261945	345724	24.89%	3.328%
62	17.486	1424	1426	1428	rBV3	23715	40749	2.93%	0.392%
63	18.537	1515	1518	1522	rVB2	114230	201302	14.49%	1.938%
64	18.686	1529	1531	1533	rBV	27285	41398	2.98%	0.398%
65	18.880	1545	1548	1552	rBV	90466	159738	11.50%	1.538%

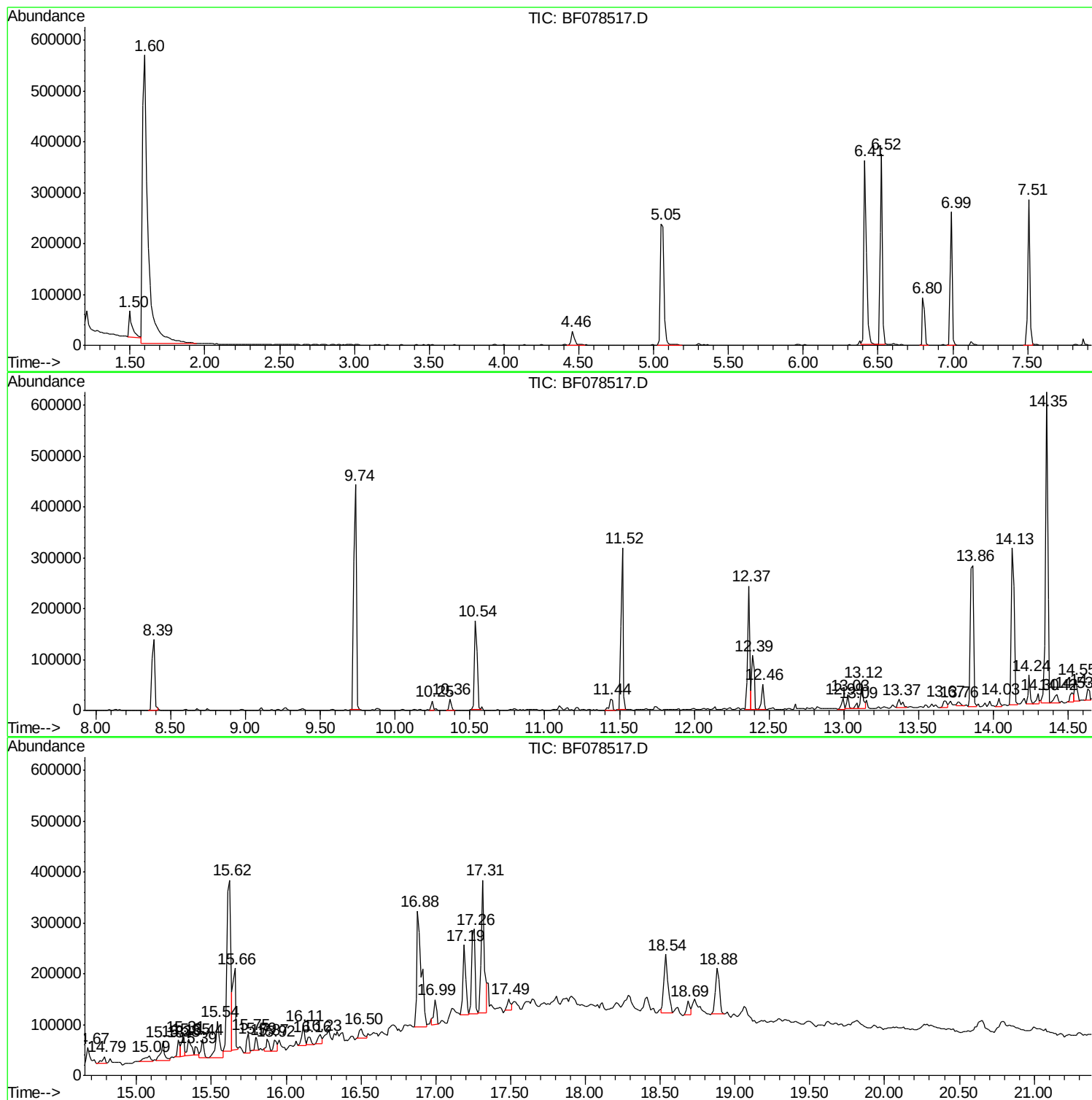
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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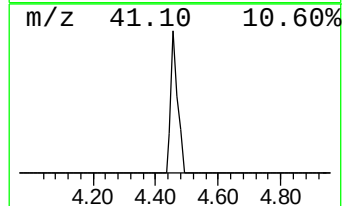
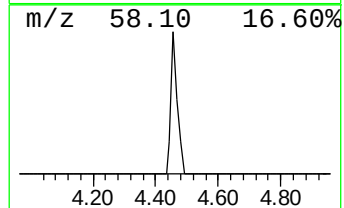
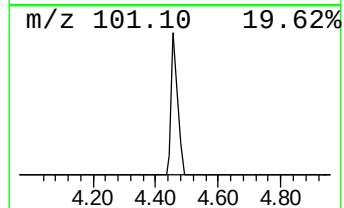
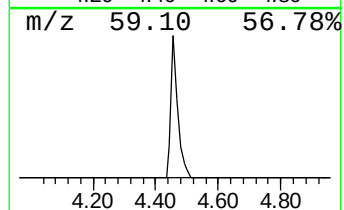
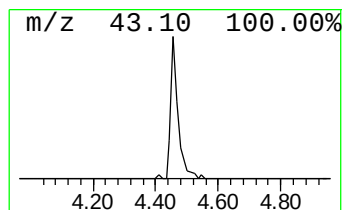
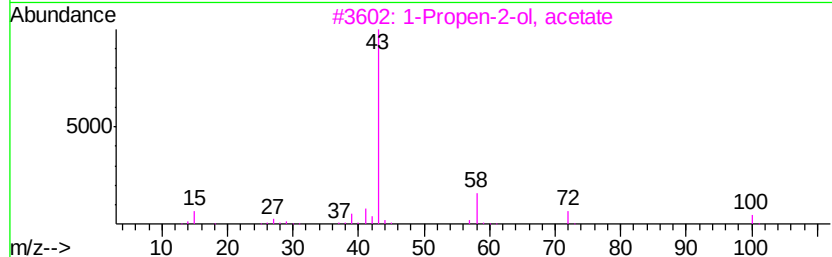
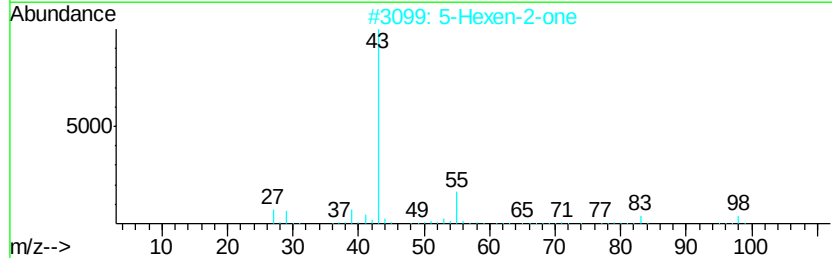
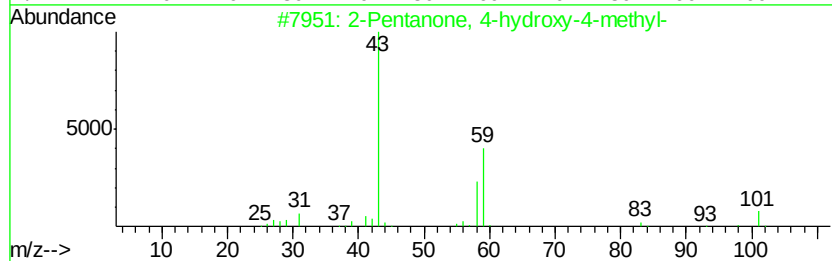
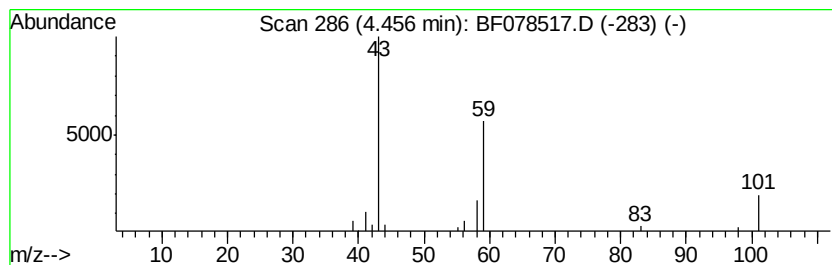
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.46	6.92 ng	39404	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	40
2		5-Hexen-2-one	98	C6H10O	000109-49-9	9
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	7
5		Ethanone, 1-(2-methylcyclopropyl)-	98	C6H10O	000930-56-3	5



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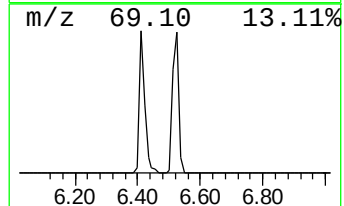
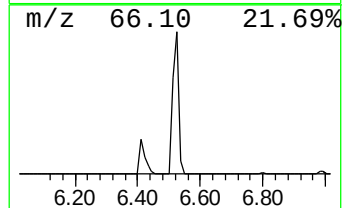
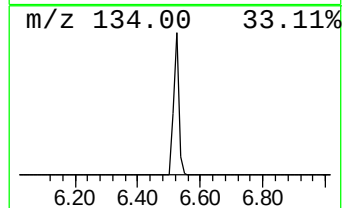
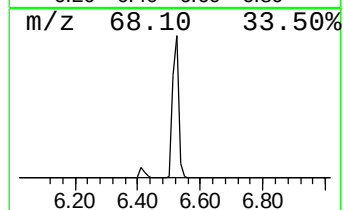
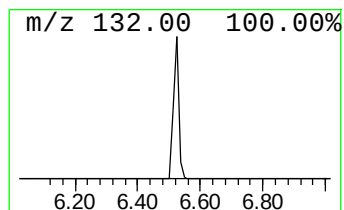
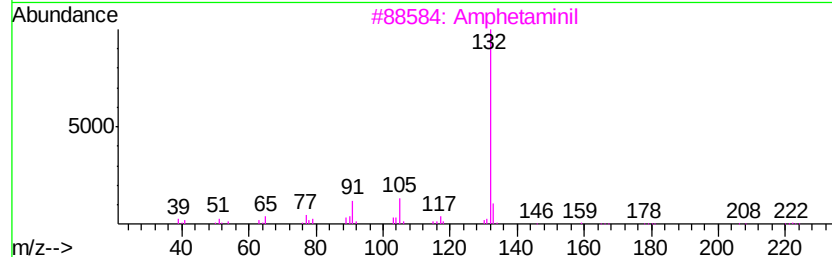
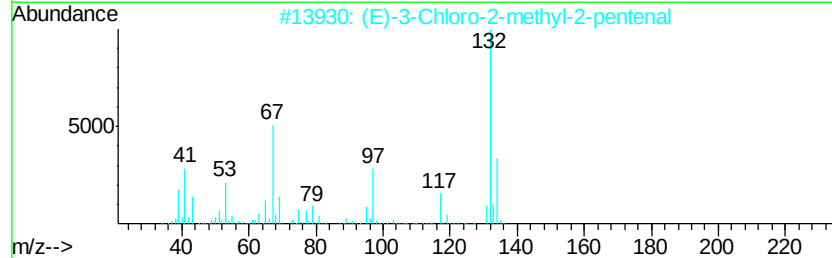
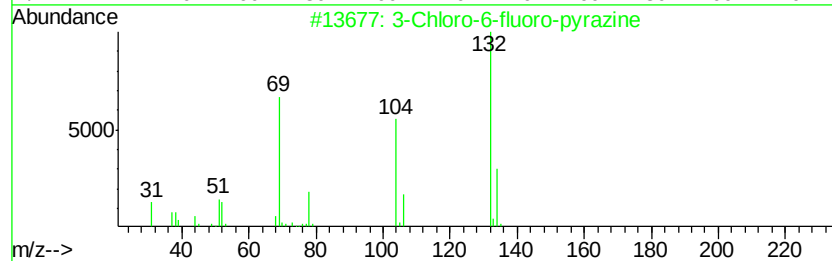
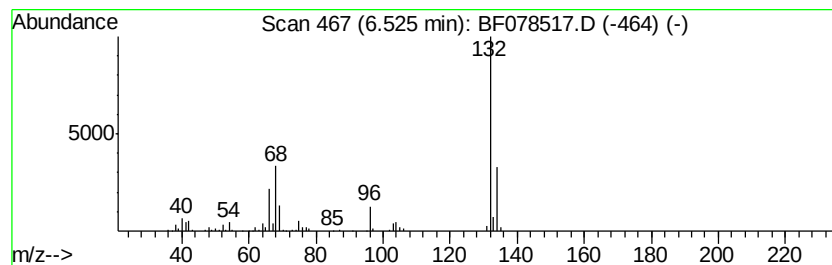
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.52 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	76.64 ng	436609	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	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Xenon	132	Xe	007440-63-3	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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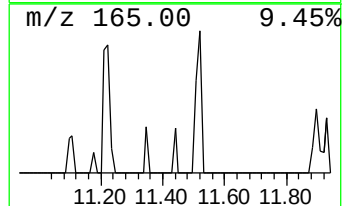
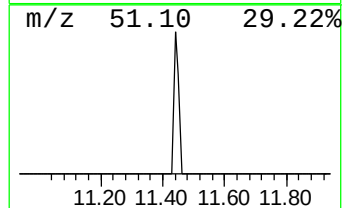
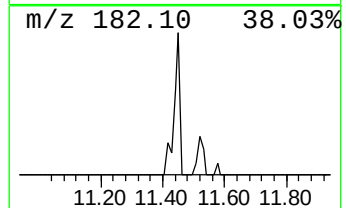
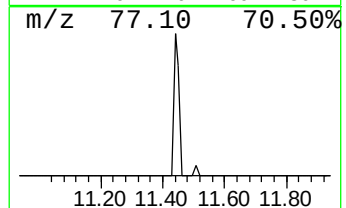
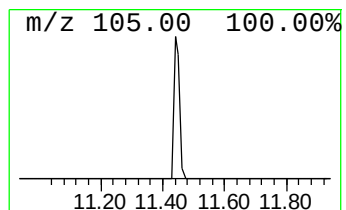
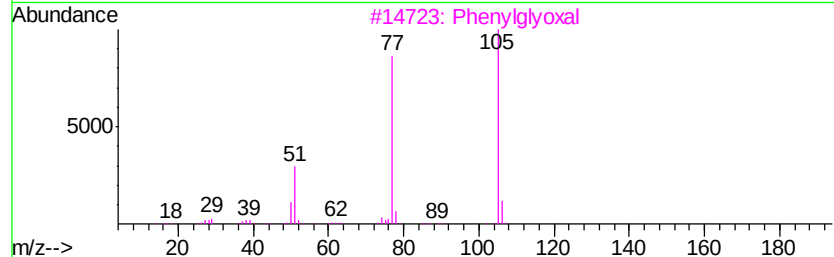
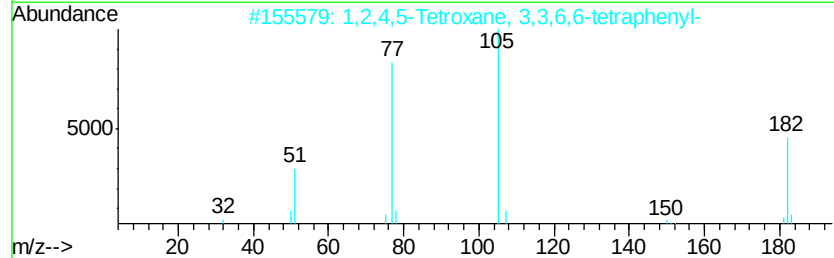
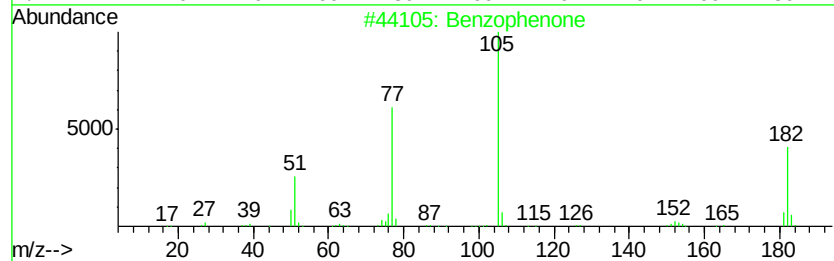
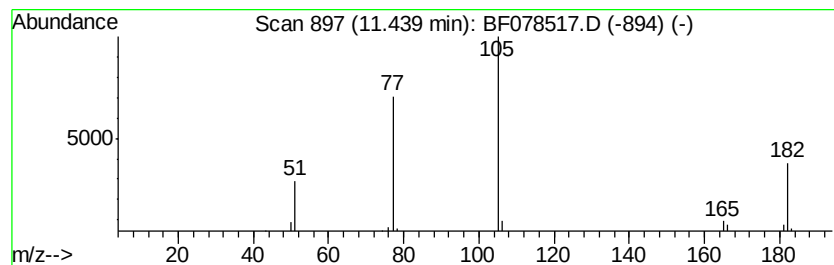
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzophenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	3.24 ng	33745	Acenaphthene-d10	10.54

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzophenone	182	C13H10O	000119-61-9	87
2		1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	59
3		Phenylglyoxal	134	C8H6O2	001074-12-0	50
4		Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	50
5		Benzoic acid, phenyl ester	198	C13H10O2	000093-99-2	50



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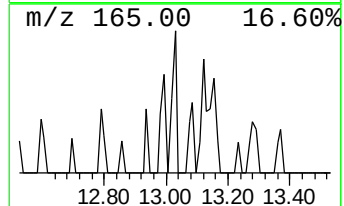
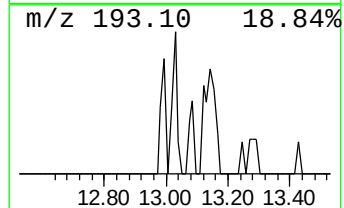
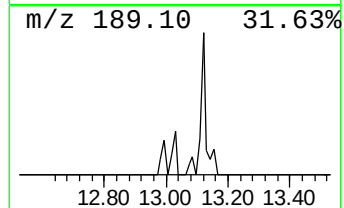
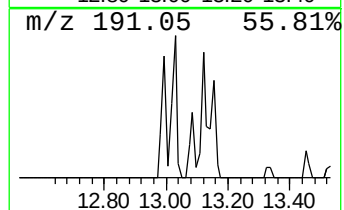
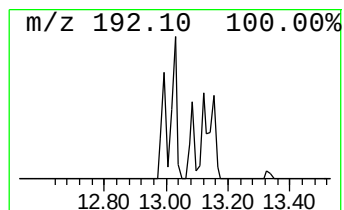
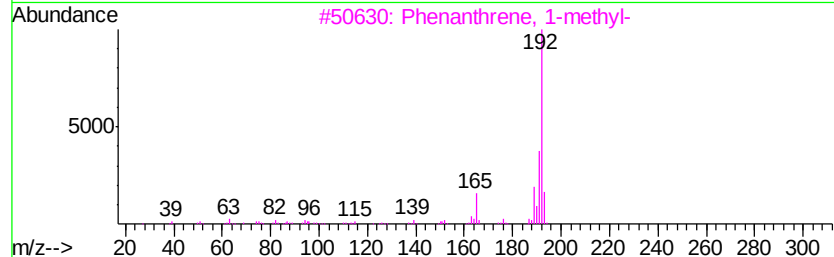
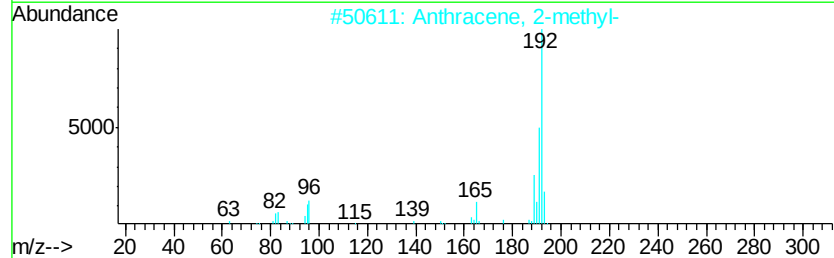
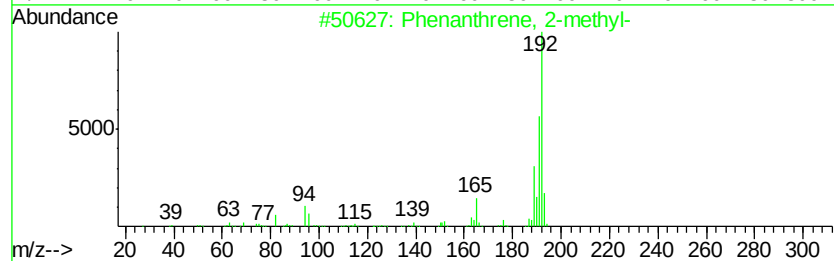
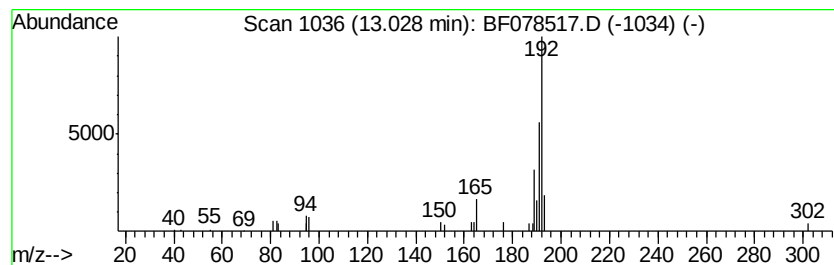
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phenanthrene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.03	2.72 ng	31044	Phenanthrene-d10	12.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	87
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	87
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	87
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	87



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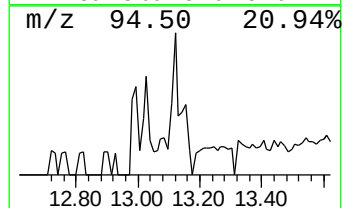
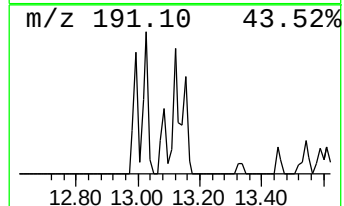
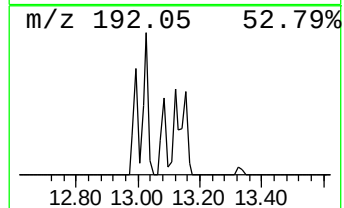
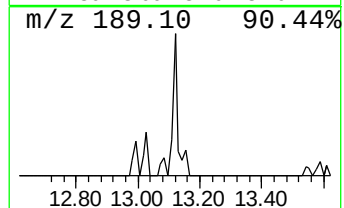
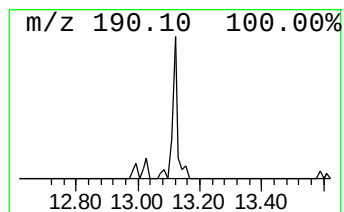
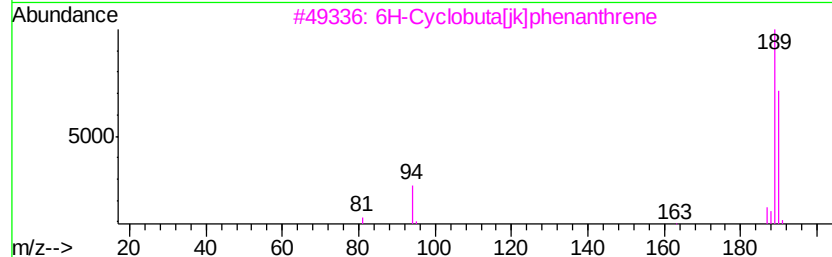
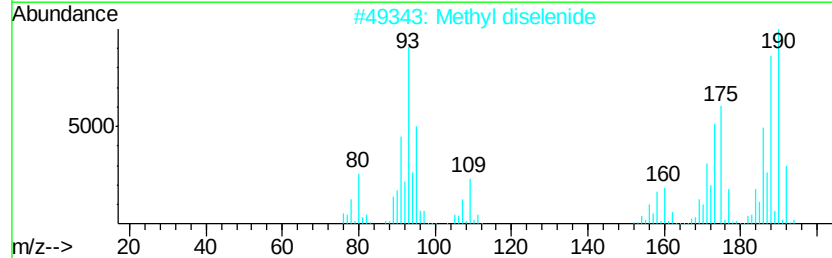
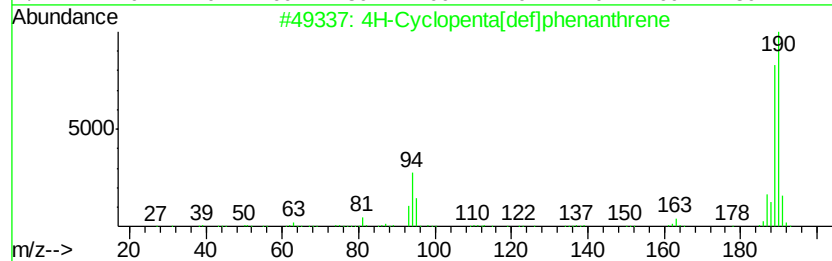
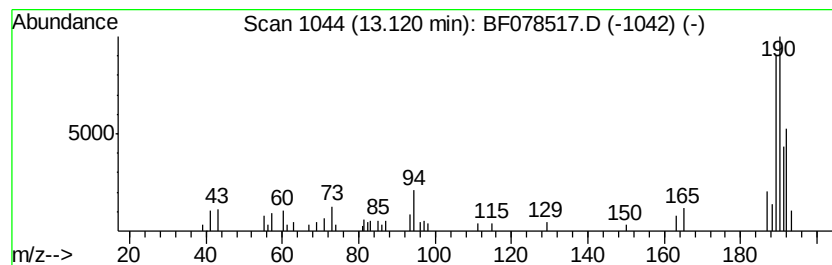
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	5.30 ng	60539	Phenanthrene-d10	12.37

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
2	Methyl diselenide	190	C2H6Se2	007101-31-7	41
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	40
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	38
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078517.D
Acq On : 14 Apr 2015 23:49
Operator : TP/IZ
Sample : G1828-14
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HS-SB04A

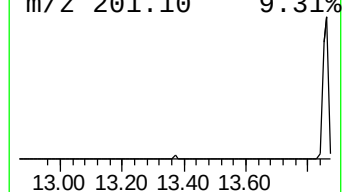
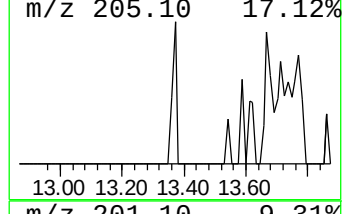
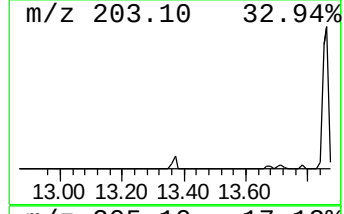
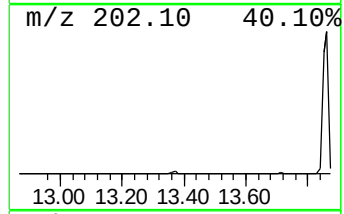
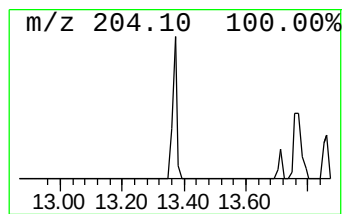
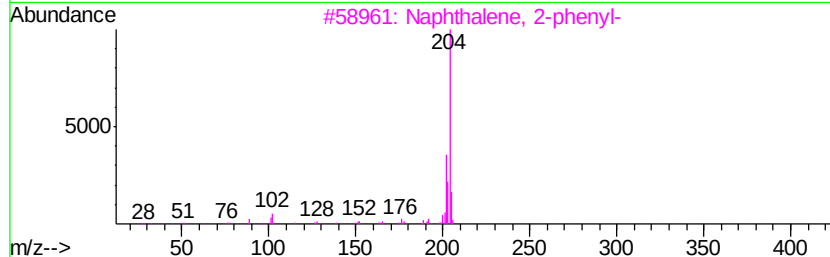
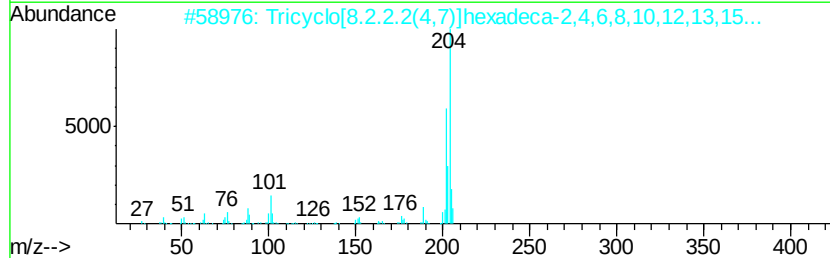
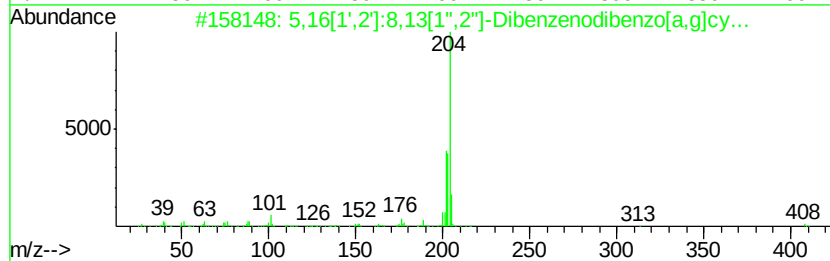
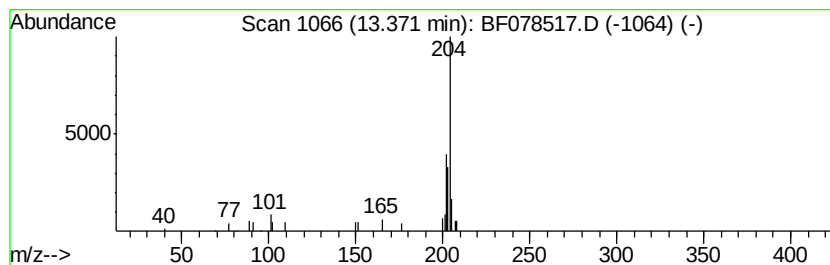
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.37	2.70 ng	30808	Phenanthrene-d10	12.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	78
2		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	58
3		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	53
4		9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	53
5		2-Phenylnaphthalene	204	C16H12	035465-71-5	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078517.D
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Sample : G1828-14
Misc :
ALS Vial : 15 Sample Multiplier: 1

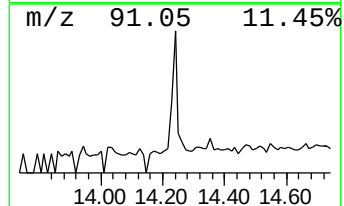
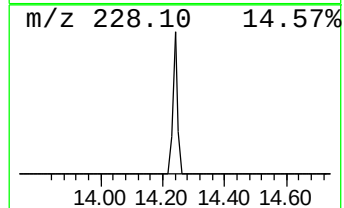
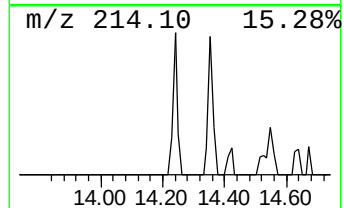
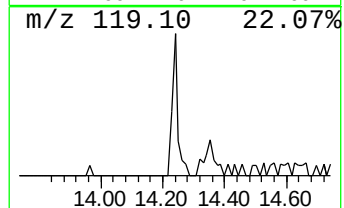
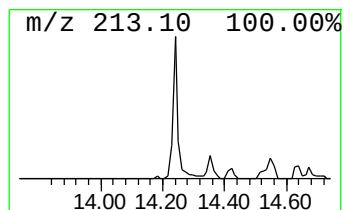
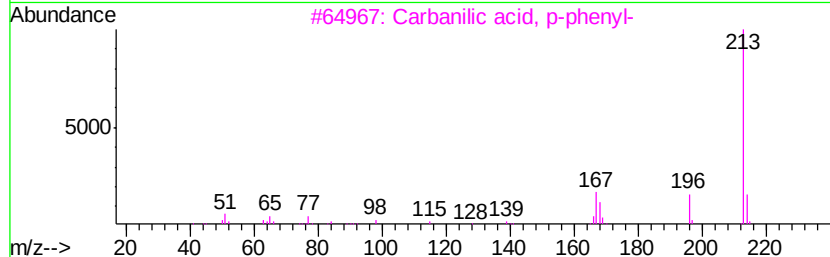
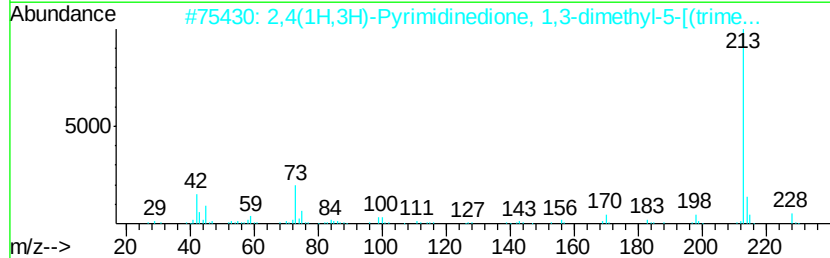
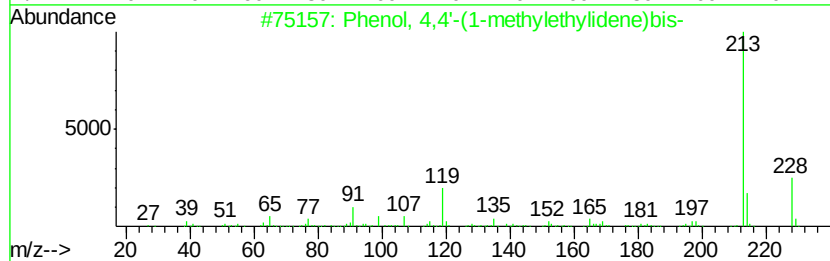
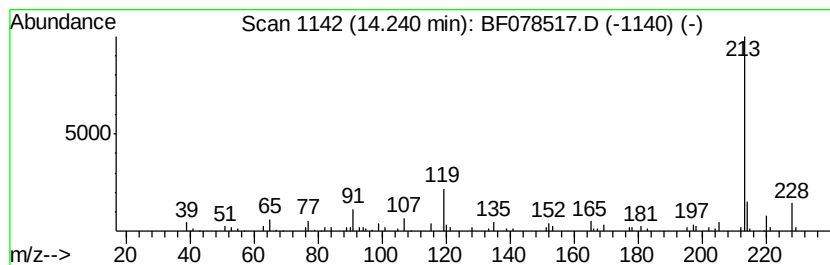
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenol, 4,4'-(1-methylethyl... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.24	2.01 ng	58674	Chrysene-d12	15.62		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	95	
2	2,4(1H,3H)-Pyrimidinedione, 1,3-...	228	C9H16N2O3Si	089447-84-7	53	
3	Carbanilic acid, p-phenyl-	213	C13H11NO2	004474-53-7	52	
4	2H,8H-Benzo[1,2-b:3,4-b']dipyran...	228	C14H12O3	000523-59-1	50	
5	2-Trifluoromethyl-5-hydroxyquino...	213	C10H6F3NO	041958-06-9	50	



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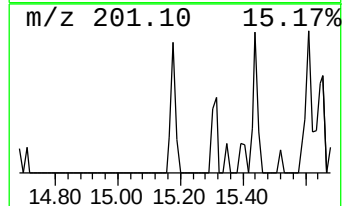
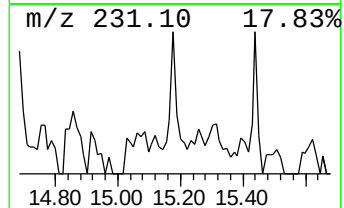
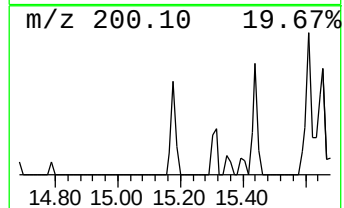
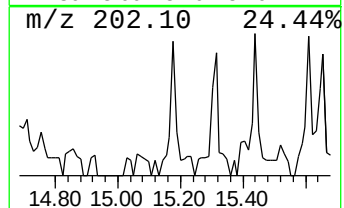
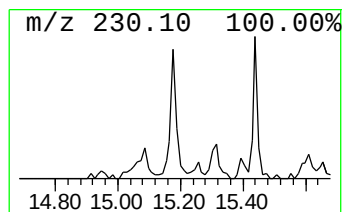
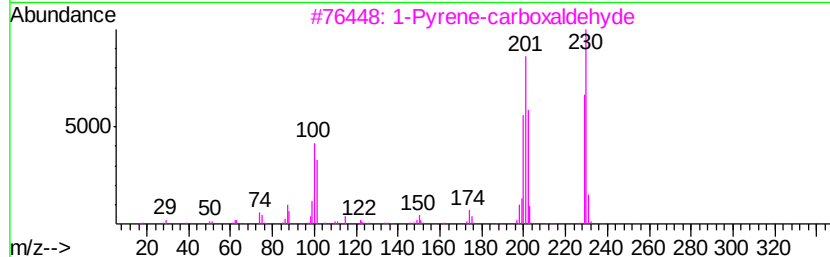
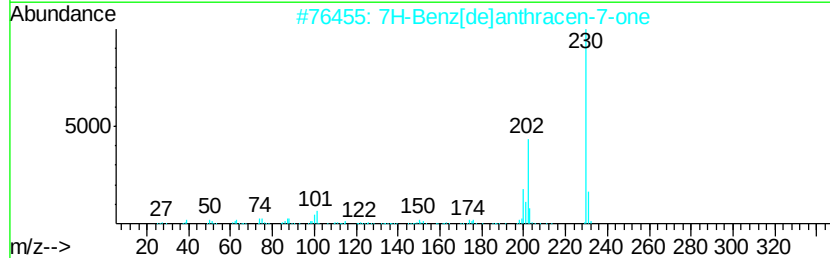
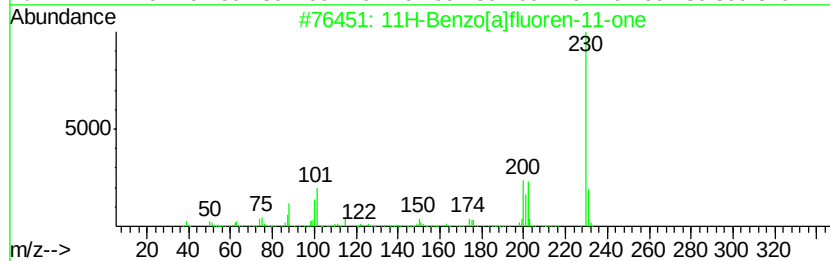
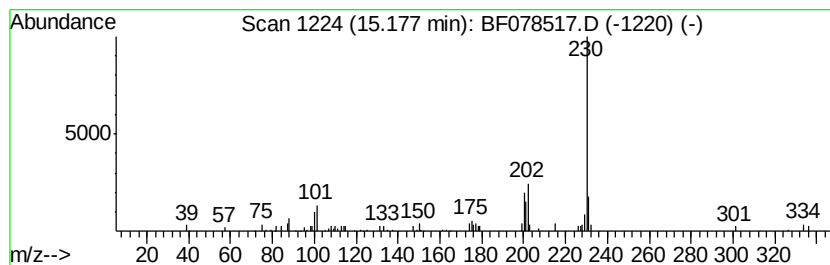
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 11H-Benzo[a]fluoren-11-one Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.		
15.18	2.41 ng	70335	Chrysene-d12	15.62		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	97
2		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	83
3		1-Pyrene-carboxaldehyde	230	C17H10O	003029-19-4	83
4		(E)-.alpha.,.alpha.'-Dicyanostil...	230	C16H10N2	002450-55-7	62
5		Pyrene, 1,3-dimethyl-	230	C18H14	064401-21-4	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078517.D
Acq On : 14 Apr 2015 23:49
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Misc :
ALS Vial : 15 Sample Multiplier: 1

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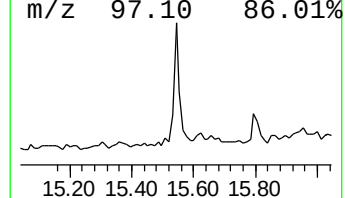
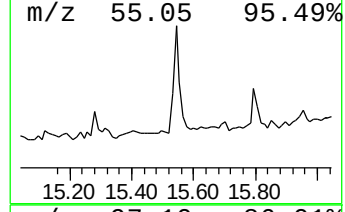
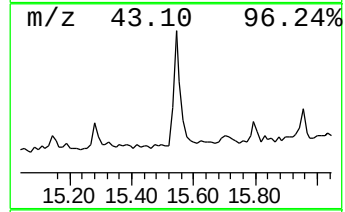
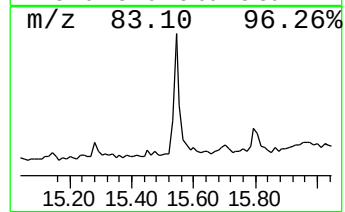
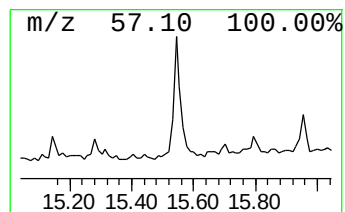
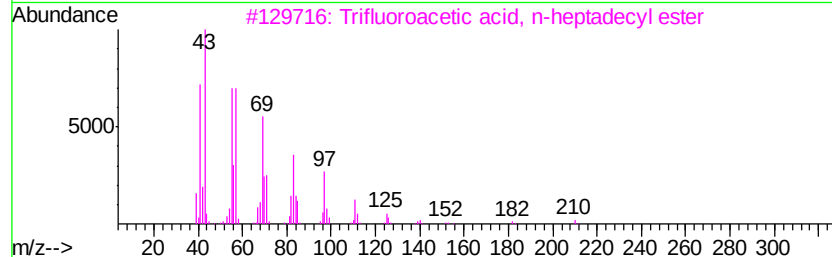
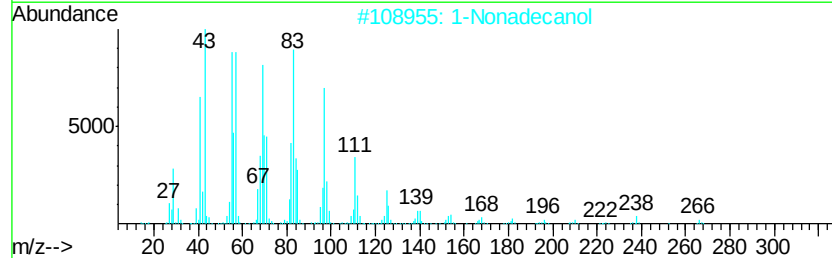
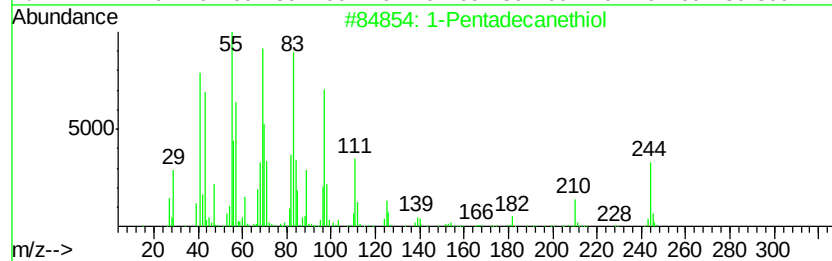
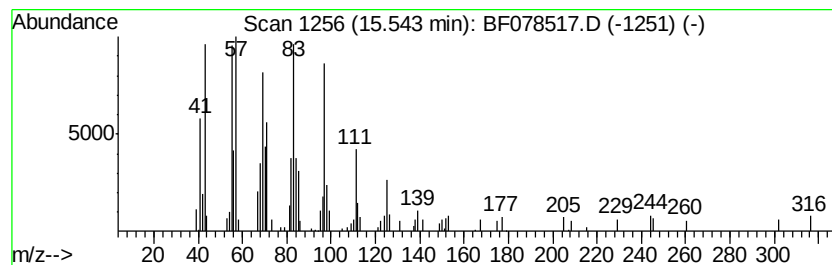
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Pentadecanethiol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.54	4.48 ng	130653	Chrysene-d12	15.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentadecanethiol	244	C15H32S	025276-70-4	96
2		1-Nonadecanol	284	C19H40O	001454-84-8	95
3		Trifluoroacetic acid, n-heptadec...	324	C17H31F3O2	1000216-79-2	91
4		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	91
5		2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078517.D
Acq On : 14 Apr 2015 23:49
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Misc :
ALS Vial : 15 Sample Multiplier: 1

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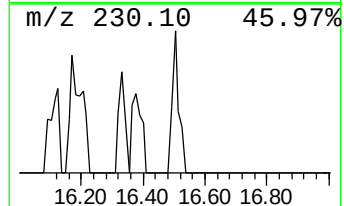
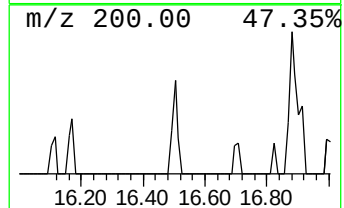
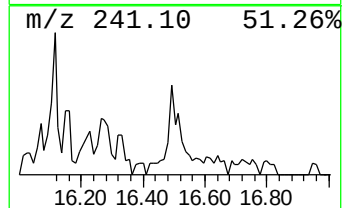
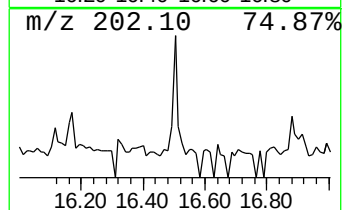
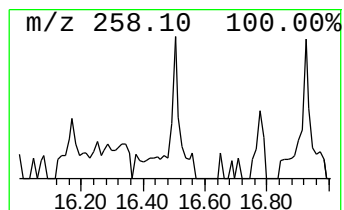
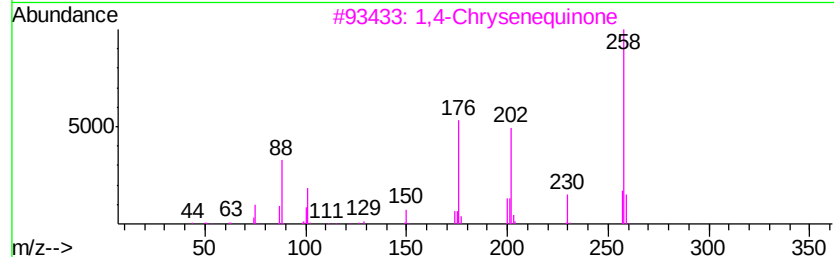
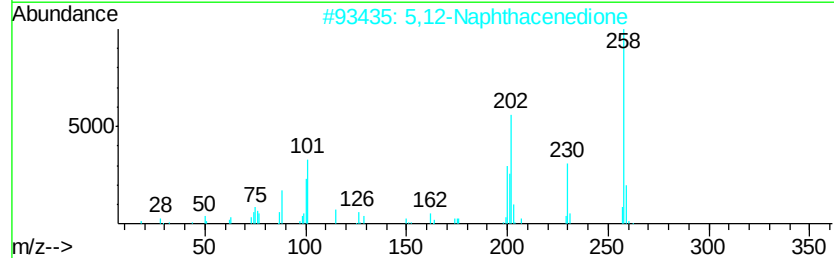
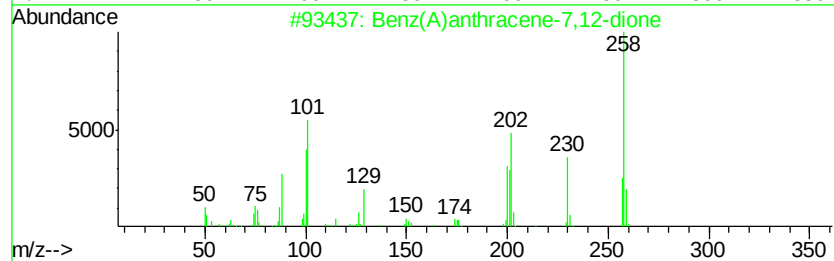
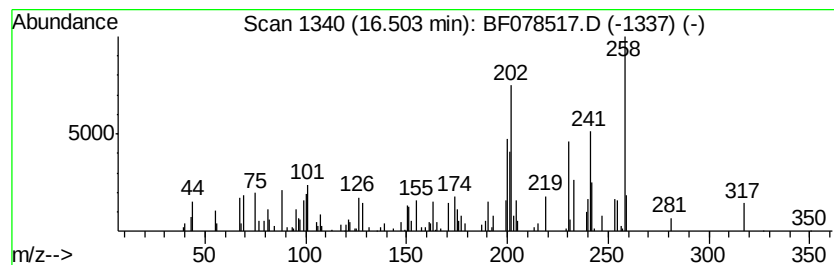
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benz(A)anthracene-7,12-dione Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	2.16 ng	37404	Perylene-d12	17.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benz(A)anthracene-7,12-dione	258	C18H10O2	002498-66-0	60
2		5,12-Naphthacenedione	258	C18H10O2	001090-13-7	38
3		1,4-Chrysenequinone	258	C18H10O2	100900-16-1	30
4		1,8-Anthracenedinitrile, 9,10-di...	258	C16H6N2O2	090866-50-5	27
5		4,8,9-Trihydroxy-6-methyl-3,4-di...	258	C15H14O4	1000210-20-3	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078517.D
Acq On : 14 Apr 2015 23:49
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Misc :
ALS Vial : 15 Sample Multiplier: 1

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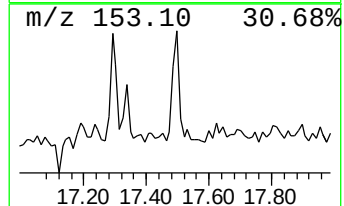
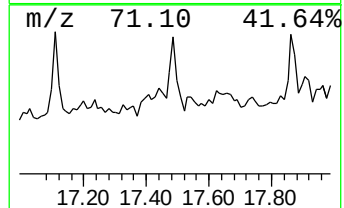
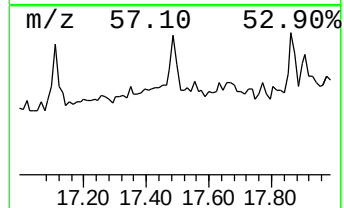
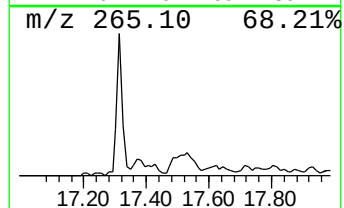
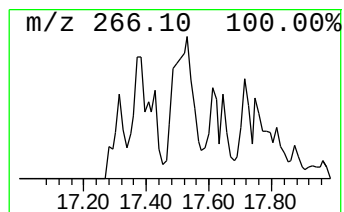
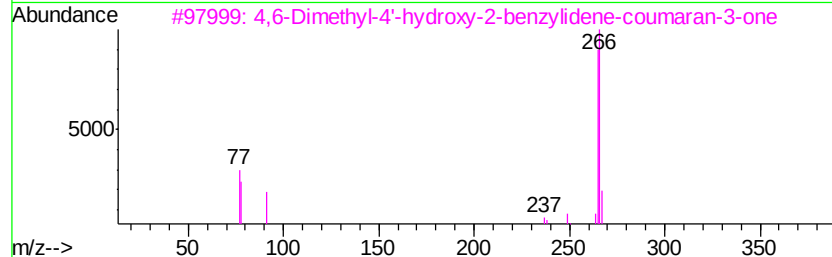
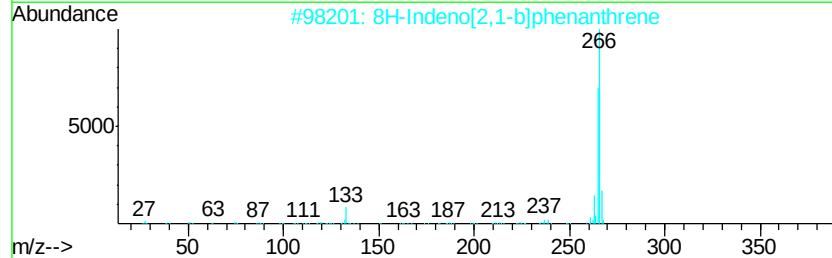
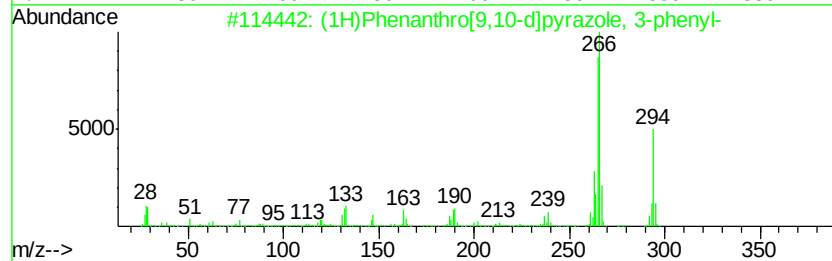
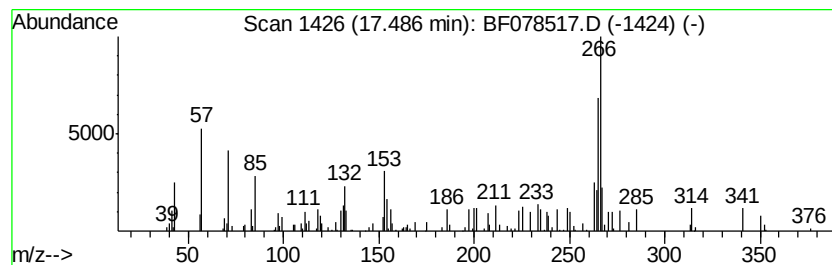
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 (1H)Phenanthro[9,10-d]pyraz... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.49	2.36 ng	40749	Perylene-d12	17.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(1H)Phenanthro[9,10-d]pyrazole, ...	294	C21H14N2	037944-54-0	53
2		8H-Indeno[2,1-b]phenanthrene	266	C21H14	000241-28-1	50
3		4,6-Dimethyl-4'-hydroxy-2-benzyl...	266	C17H14O3	002567-73-9	50
4		5-Hydroxy-4-methoxy-6H-indolo[3,...	266	C15H10N2O3	018110-86-6	49
5		4H-1-Benzopyran-4-one, 3-hydroxy...	266	C17H14O3	078396-37-9	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078517.D
Acq On : 14 Apr 2015 23:49
Operator : TP/IZ
Sample : G1828-14
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HS-SB04A

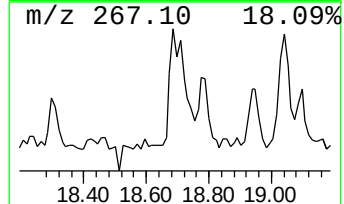
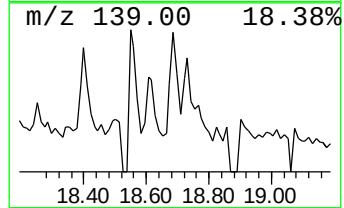
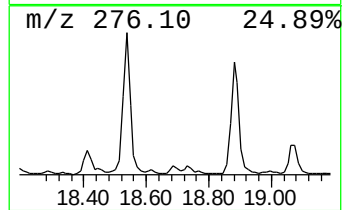
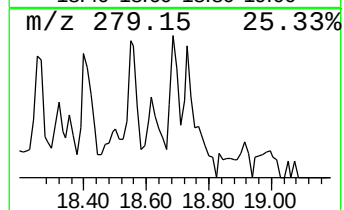
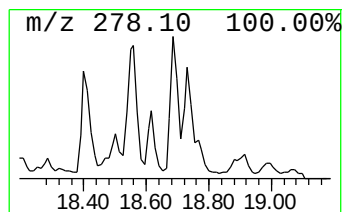
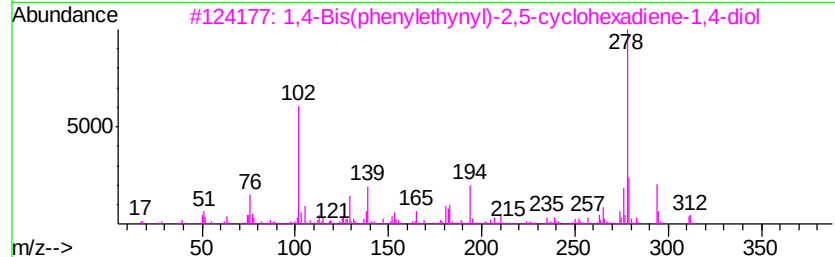
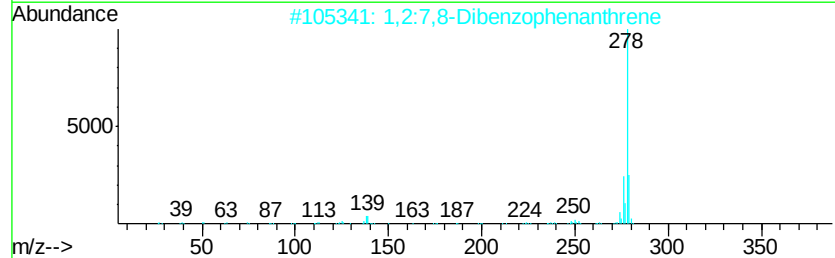
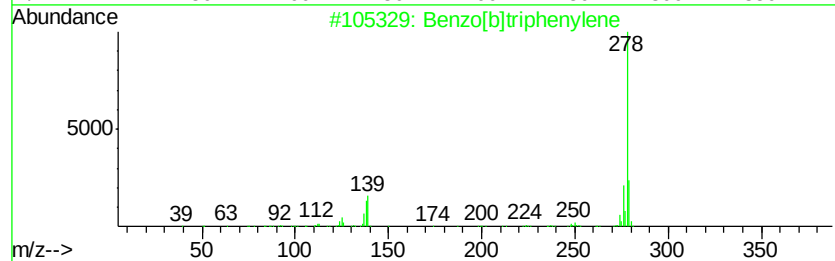
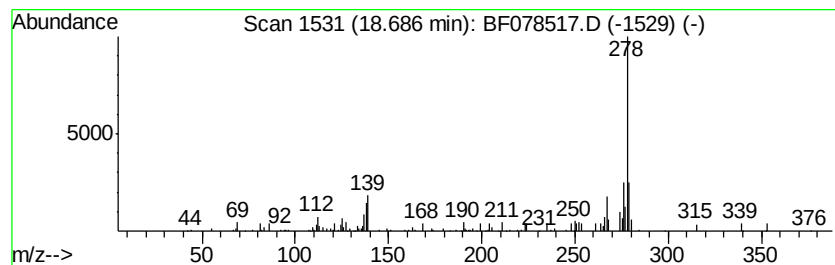
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Benzo[b]triphenylene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.69	2.39 ng	41398	Perylene-d12	17.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]triphenylene	278	C22H14	000215-58-7	92
2		1,2:7,8-Dibenzophenanthrene	278	C22H14	000213-46-7	86
3		1,4-Bis(phenylethynyl)-2,5-cyclo...	312	C22H16O2	055006-58-1	72
4		Benzo[b]chrysene	278	C22H14	000214-17-5	70
5		Pentaphene	278	C22H14	000222-93-5	70



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078517.D
 Acq On : 14 Apr 2015 23:49
 Operator : TP/IZ
 Sample : G1828-14
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB04A

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
2-Pentanone, 4-hy...	4.46	6.9 ng		39404	1	6.80	113933	20.0
unknown6.52	6.52	76.6 ng		436609	1	6.80	113933	20.0
Benzophenone	11.44	3.2 ng		33745	3	10.54	208020	20.0
Phenanthrene, 2-m...	13.03	2.7 ng		31044	4	12.37	228520	20.0
4H-Cyclopenta[def...	13.12	5.3 ng		60539	4	12.37	228520	20.0
5,16[1',2']:8,13[...	13.37	2.7 ng		30808	4	12.37	228520	20.0
Phenol, 4,4'-(1-m...	14.24	2.0 ng		58674	5	15.62	583387	20.0
11H-Benzo[a]fluor...	15.18	2.4 ng		70335	5	15.62	583387	20.0
1-Pentadecanethiol	15.54	4.5 ng		130653	5	15.62	583387	20.0
Benz(A)anthracene...	16.50	2.2 ng		37404	6	17.31	345724	20.0
(1H)Phenanthro[9,...	17.49	2.4 ng		40749	6	17.31	345724	20.0
Benzo[b]triphenylene	18.69	2.4 ng		41398	6	17.31	345724	20.0