

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
 Data File : BF079725.D
 Acq On : 5 Jun 2015 4:14
 Operator : TP/IZ
 Sample : G2474-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C-5

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-BF060415.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	2.261	92	94	108	rVB	129978	348443	10.94%	0.954%
2	2.455	108	111	125	rVV2	553021	1409465	44.26%	3.860%
3	2.673	126	130	142	rVB	81630	233080	7.32%	0.638%
4	2.935	150	153	157	rVB	288187	420616	13.21%	1.152%
5	6.124	430	432	435	rBV	2283251	2201775	69.14%	6.030%
6	7.107	515	518	520	rBV	2174359	2454153	77.07%	6.721%
7	7.279	530	533	535	rBV	2697796	2547838	80.01%	6.978%
8	7.507	551	553	555	rBV	705843	582434	18.29%	1.595%
9	7.667	565	567	570	rBV	2217761	1864511	58.55%	5.106%
10	8.067	599	602	604	rBV	1088930	1356606	42.60%	3.715%
11	8.799	664	666	669	rBV	926924	814898	25.59%	2.232%
12	9.873	758	760	763	rVB	3363568	2739080	86.01%	7.502%
13	10.262	792	794	795	rBV	282579	272441	8.56%	0.746%
14	10.582	820	822	824	rBV	825871	951379	29.88%	2.606%
15	11.371	889	891	894	rBV	2023110	1790370	56.22%	4.903%
16	12.091	951	954	955	rBV	1228497	1146974	36.02%	3.141%
17	12.605	996	999	1001	rBV2	97451	199381	6.26%	0.546%
18	12.731	1008	1010	1014	rBV2	144540	283168	8.89%	0.776%
19	13.382	1065	1067	1069	rBV	1042546	1091237	34.27%	2.989%
20	13.645	1088	1090	1093	rVB	997973	971327	30.50%	2.660%
21	13.794	1100	1103	1105	rBV	2243152	3184516	100.00%	8.722%
22	14.011	1118	1122	1125	rVB	114059	257258	8.08%	0.705%
23	14.765	1184	1188	1190	rBV2	96798	199634	6.27%	0.547%
24	14.834	1192	1194	1197	rBV3	374997	539648	16.95%	1.478%
25	15.074	1212	1215	1217	rBV2	989943	1779258	55.87%	4.873%
26	15.108	1217	1218	1224	rVB	366471	492001	15.45%	1.347%
27	15.680	1265	1268	1273	rBV3	154875	283575	8.90%	0.777%
28	16.468	1334	1337	1342	rBV	373202	1006522	31.61%	2.757%
29	16.571	1343	1346	1353	rVB4	141141	518688	16.29%	1.421%
30	16.880	1370	1373	1377	rVB	227191	364542	11.45%	0.998%
31	16.960	1377	1380	1385	rBV	261074	468081	14.70%	1.282%
32	17.051	1385	1388	1391	rBV	734978	1206103	37.87%	3.303%
33	17.714	1443	1446	1448	rBV	128585	240234	7.54%	0.658%
34	18.034	1471	1474	1481	rVB8	51324	178044	5.59%	0.488%

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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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060415.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	18.823	1540	1543	1548	rVB4	87203	235189	7.39%	0.644%
36	19.063	1560	1564	1569	rBV2	160667	451586	14.18%	1.237%
37	19.531	1601	1605	1608	rBV3	112767	293317	9.21%	0.803%
38	19.600	1608	1611	1614	rVV	193093	452096	14.20%	1.238%
39	19.669	1614	1617	1628	rVB	135868	395996	12.44%	1.085%
40	19.920	1635	1639	1642	rBV3	103101	287783	9.04%	0.788%

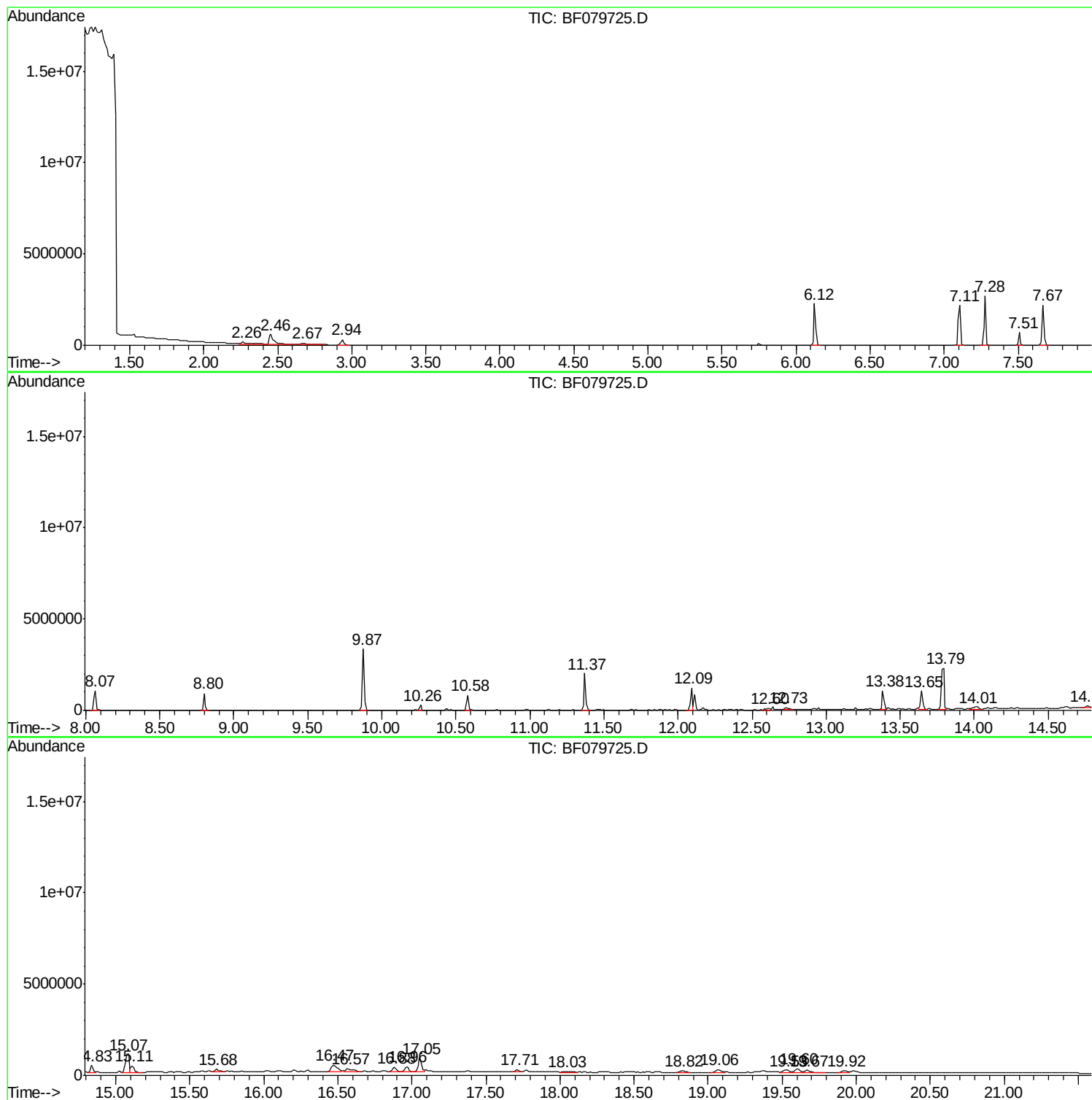
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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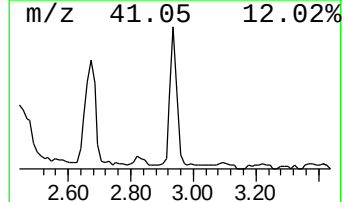
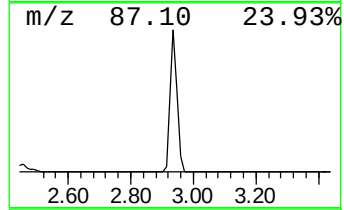
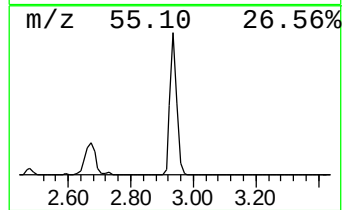
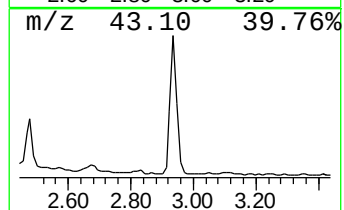
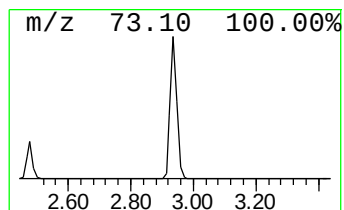
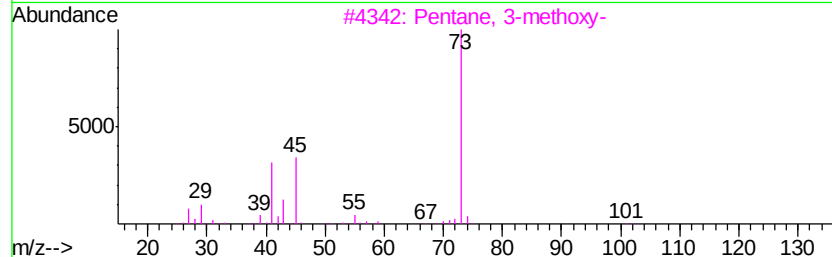
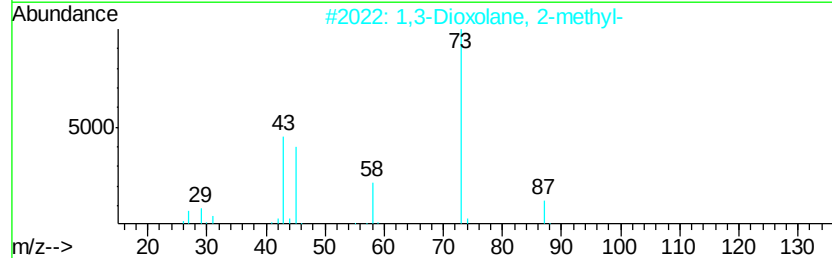
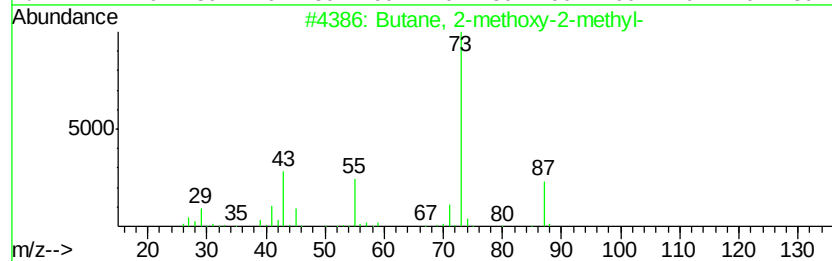
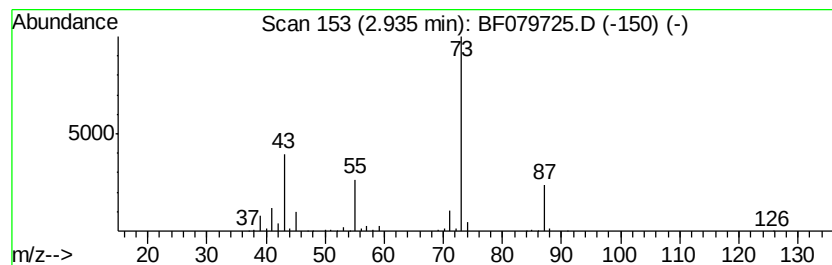
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.94	14.44 ng	420616	1,4-Dichlorobenzene-d4	7.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	10
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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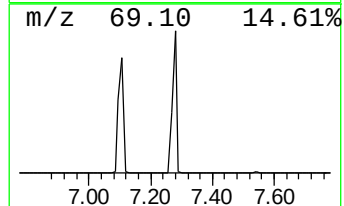
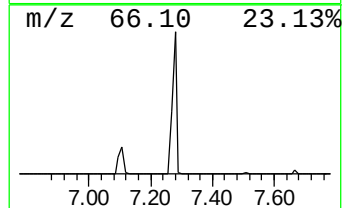
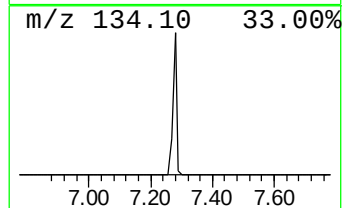
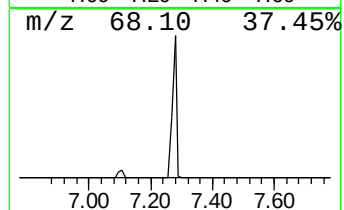
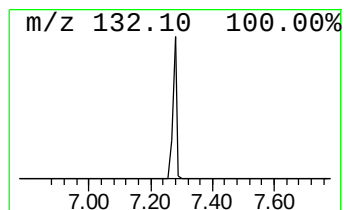
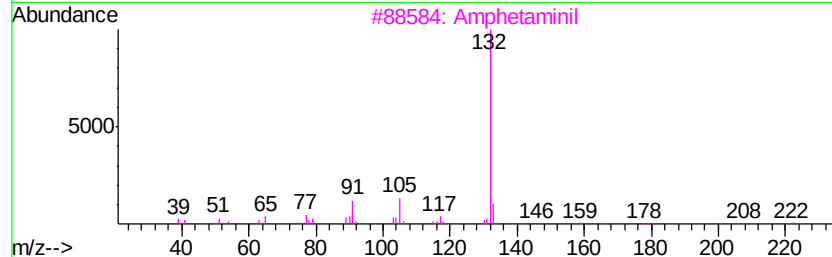
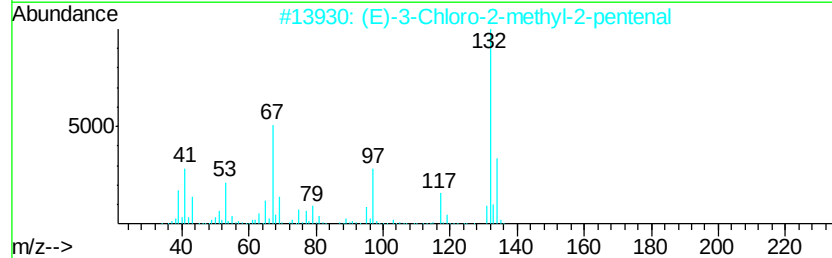
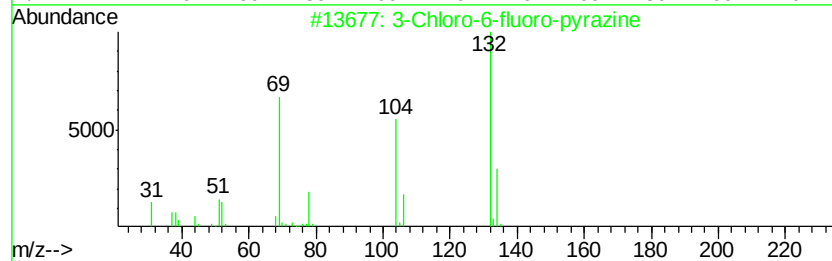
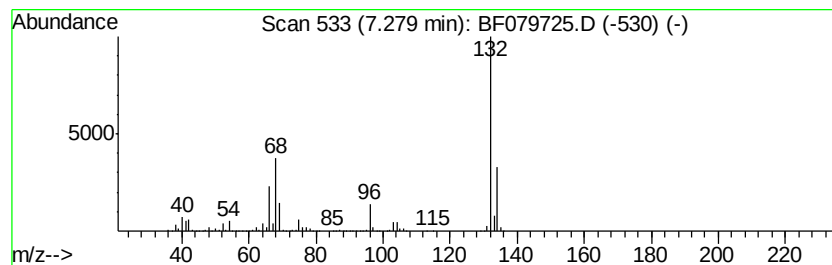
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown7.28 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.28	87.49 ng	2547840	1,4-Dichlorobenzene-d4	7.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	12



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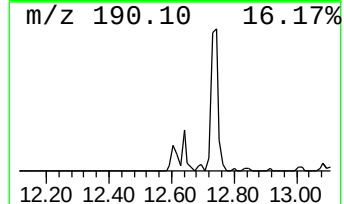
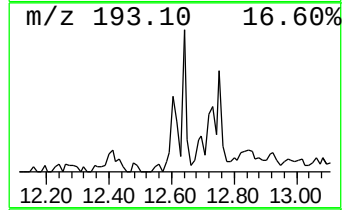
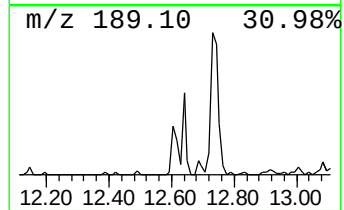
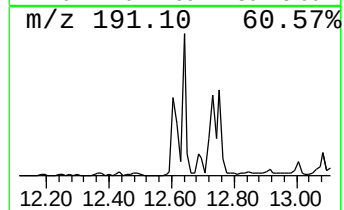
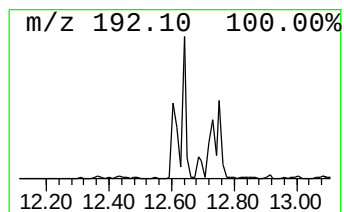
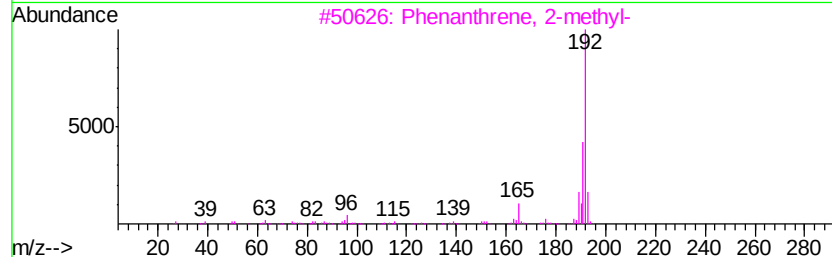
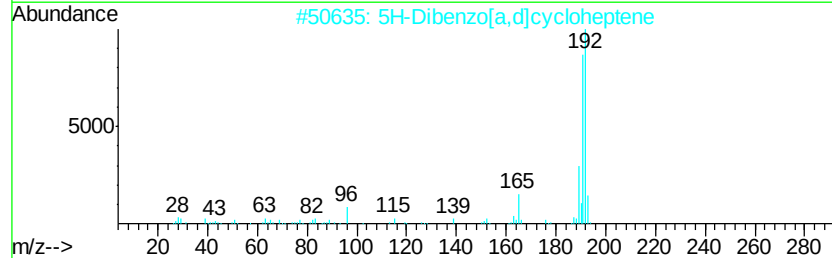
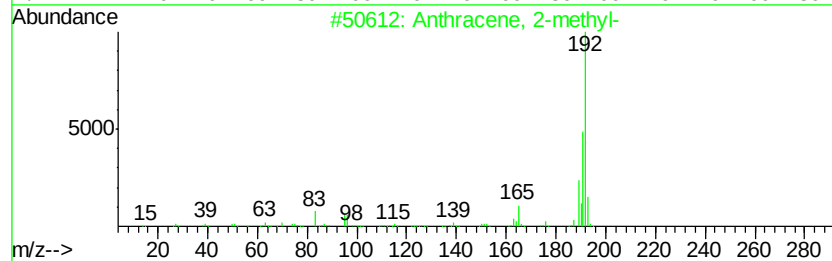
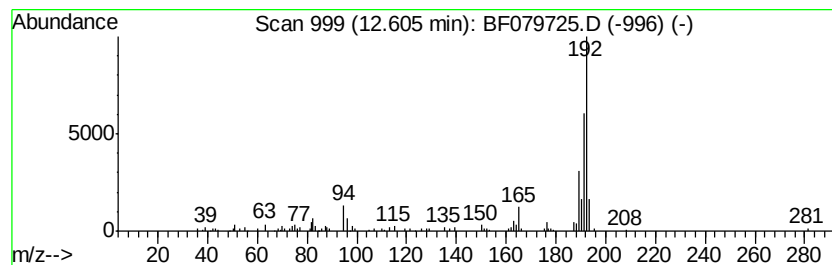
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Anthracene, 2-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.61	3.48 ng	199381	Phenanthrene-d10	12.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
2		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	94
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	90



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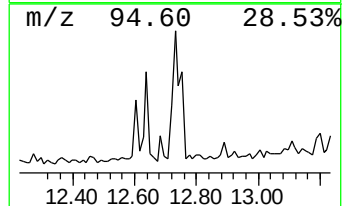
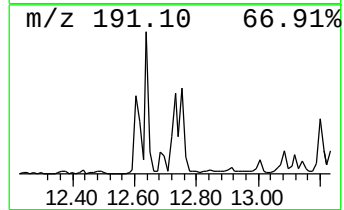
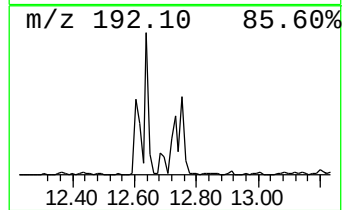
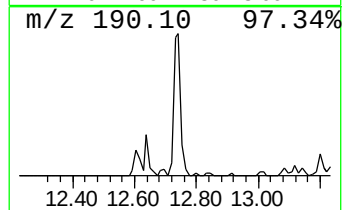
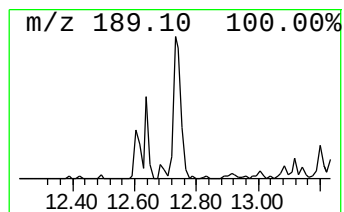
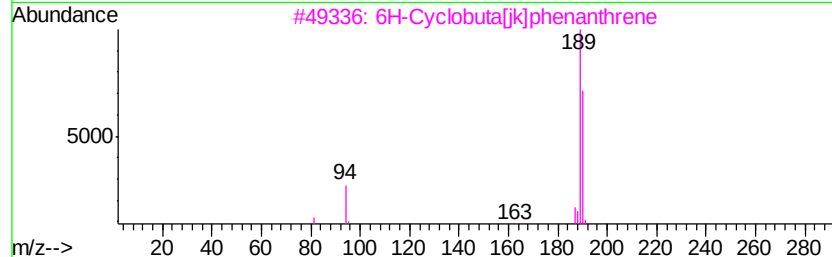
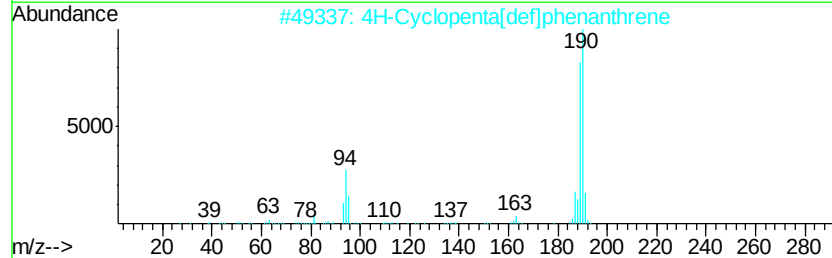
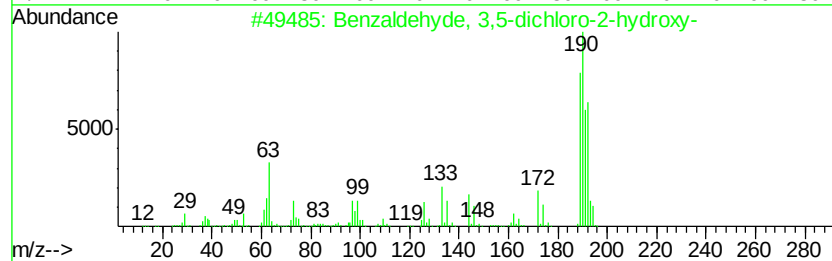
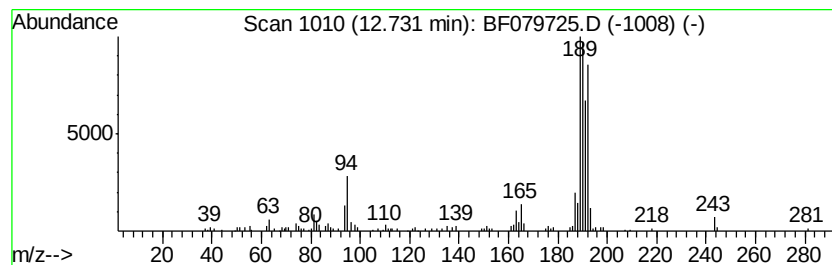
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzaldehyde, 3,5-dichloro-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.73	4.94 ng	283168	Phenanthrene-d10	12.09	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	72
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	49
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42
5	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	42



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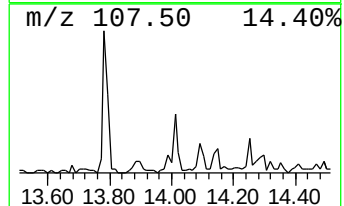
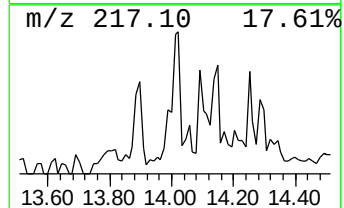
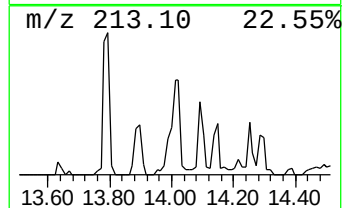
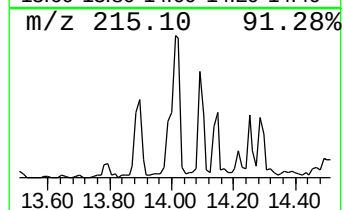
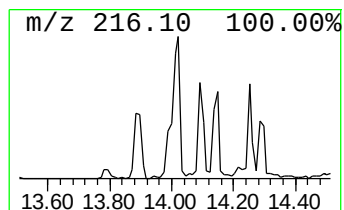
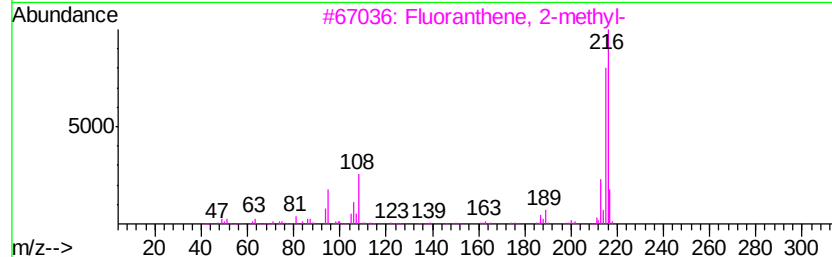
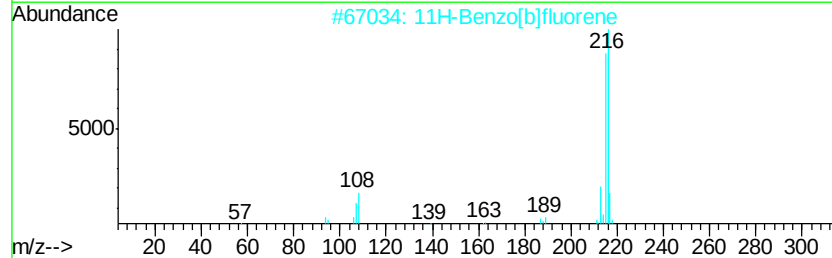
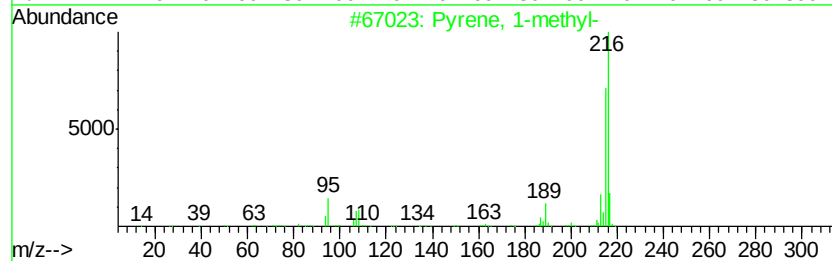
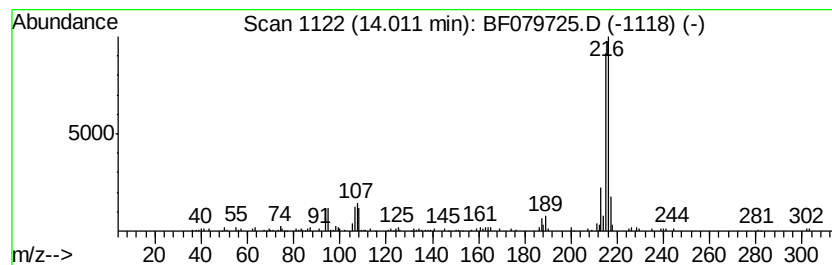
Instrument :
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ClientSampled :
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Pyrene, 1-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.01	2.89 ng	257258	Chrysene-d12	15.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Pyrene, 1-methyl-	216 C17H12	002381-21-7 94
2		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 90
3		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 90
4		11H-Benzo[a]fluorene	216 C17H12	000238-84-6 89
5		7H-Benzo[c]fluorene	216 C17H12	000205-12-9 68



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Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
C-5

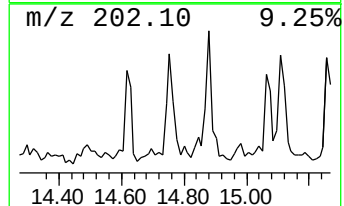
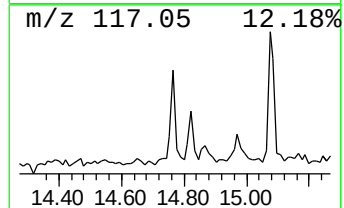
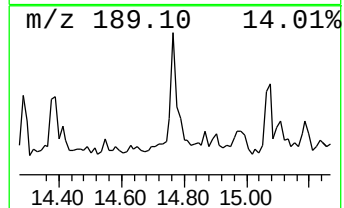
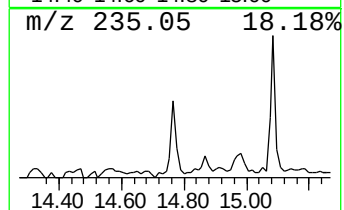
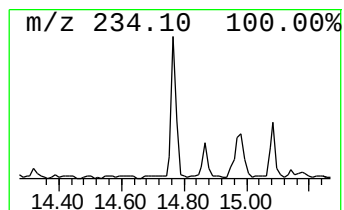
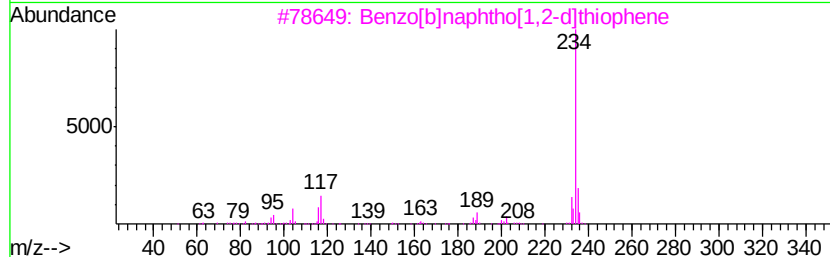
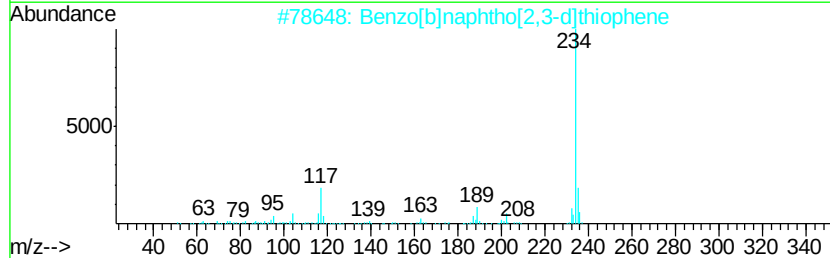
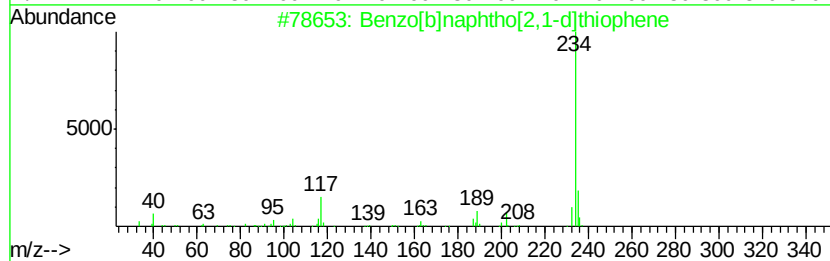
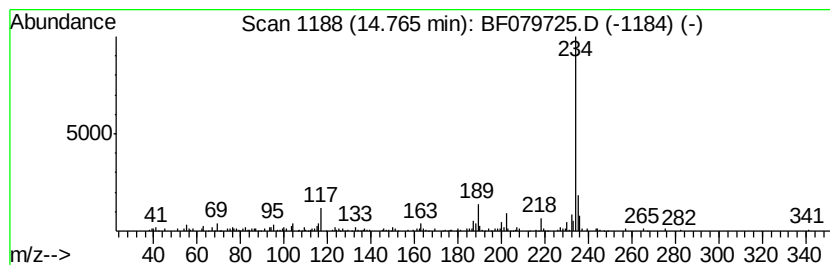
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.77	2.24 ng	199634	Chrysene-d12	15.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	96
2		Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	95
3		Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	90
4		Anthra[1,9a,9-cd;5,10a,10-c'd']d...	234	C14H6N2O2	000201-91-2	53
5		2-Amino-6-t-butyl-3-cyano-4,5,6,...	234	C13H18N2S	042159-76-2	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
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Acq On : 5 Jun 2015 4:14
Operator : TP/IZ
Sample : G2474-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

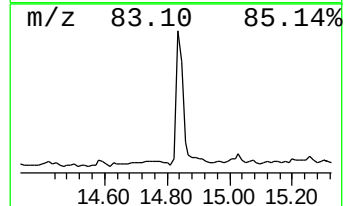
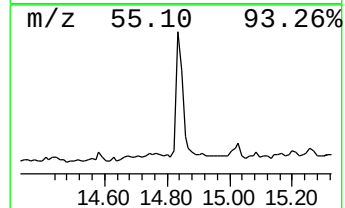
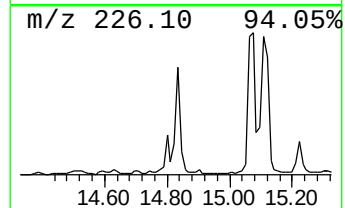
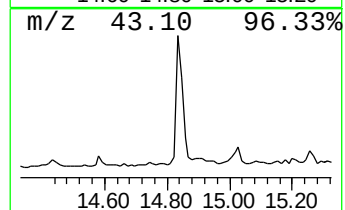
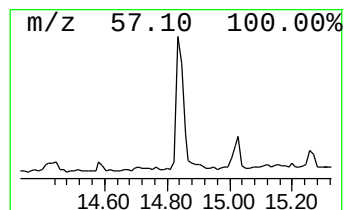
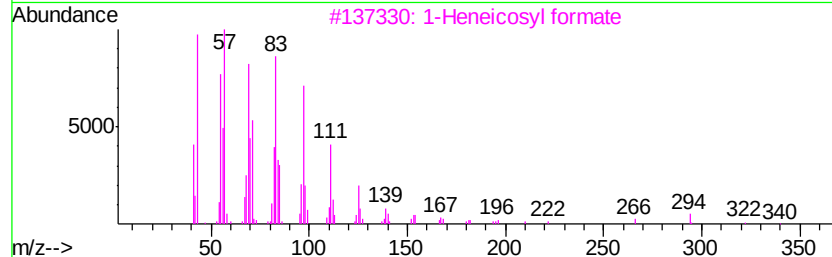
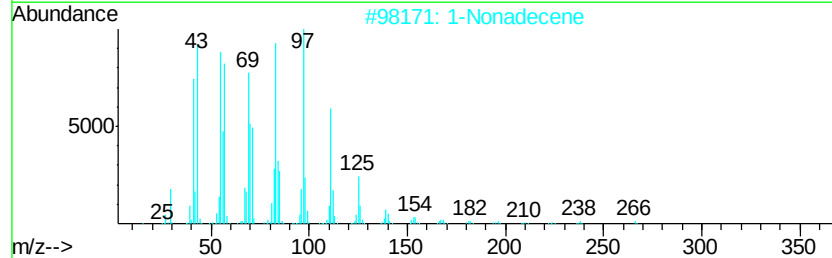
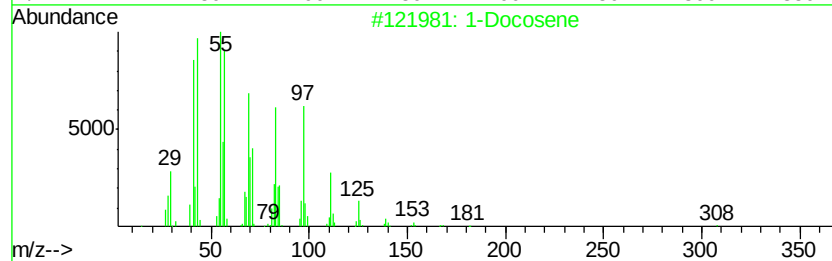
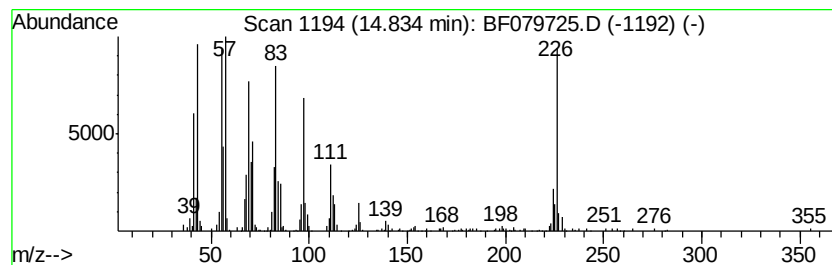
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Docosene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.83	6.07 ng	539648	Chrysene-d12	15.09	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	86
2	1-Nonadecene	266	C19H38	018435-45-5	76
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	76
4	Bromoacetic acid, hexadecyl ester	362	C18H35BrO2	005454-48-8	70
5	1-Hexadecanol	242	C16H34O	036653-82-4	70



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
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Sample : G2474-05
Misc :
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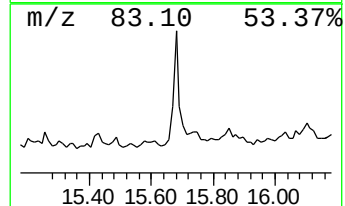
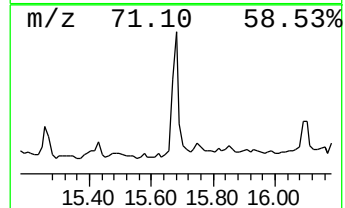
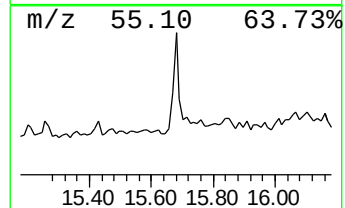
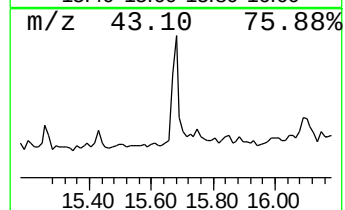
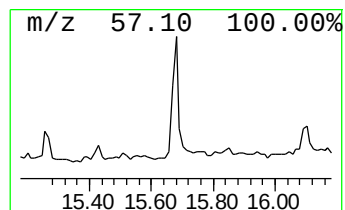
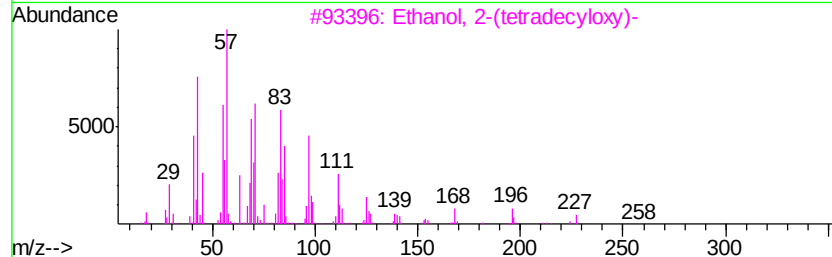
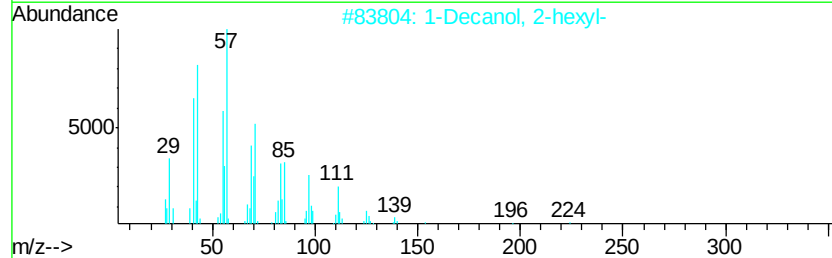
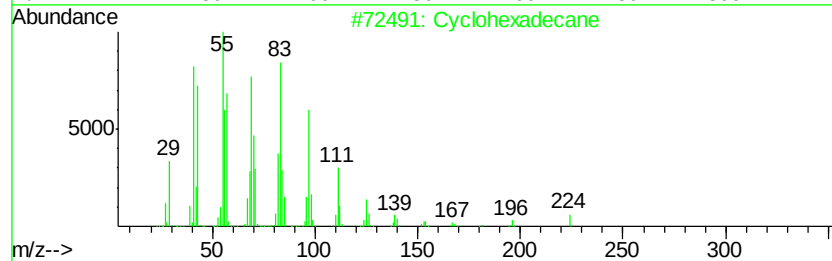
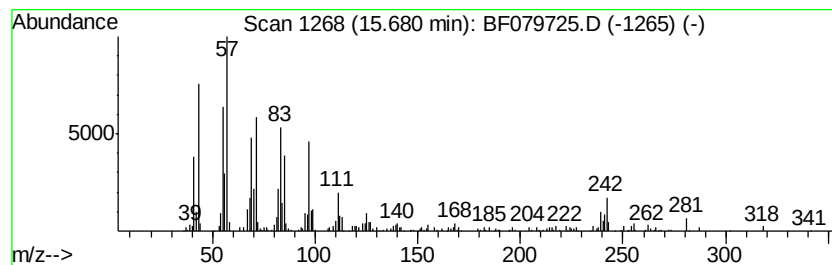
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Cyclohexadecane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.68	3.19 ng	283575	Chrysene-d12	15.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexadecane	224 C16H32	000295-65-8 83
2		1-Decanol, 2-hexyl-	242 C16H34O	002425-77-6 60
3		Ethanol, 2-(tetradecyloxy)-	258 C16H34O2	002136-70-1 58
4		10-Heneicosene (c,t)	294 C21H42	095008-11-0 53
5		Heptafluorobutanoic acid, heptad...	452 C21H35F7O2	1000282-97-3 53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
Data File : BF079725.D
Acq On : 5 Jun 2015 4:14
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Sample : G2474-05
Misc :
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Instrument :
BNA_F
ClientSampleId :
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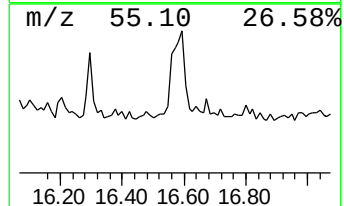
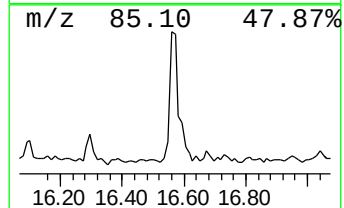
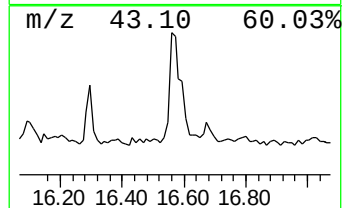
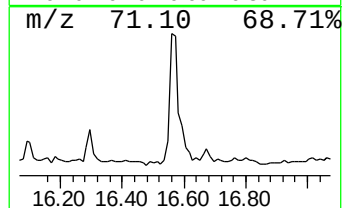
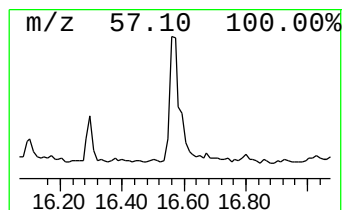
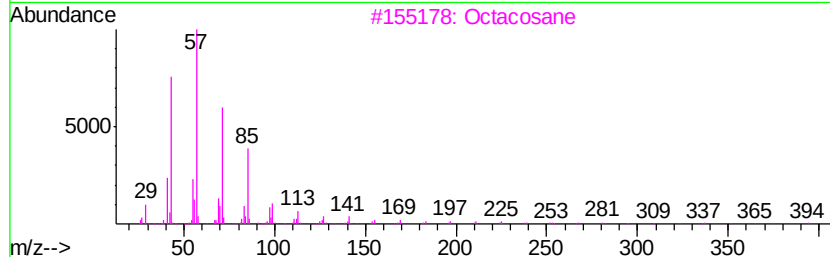
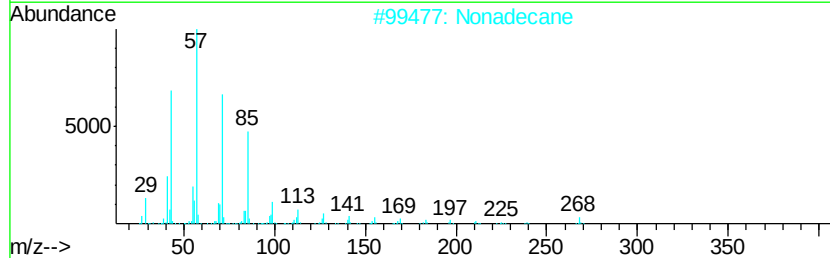
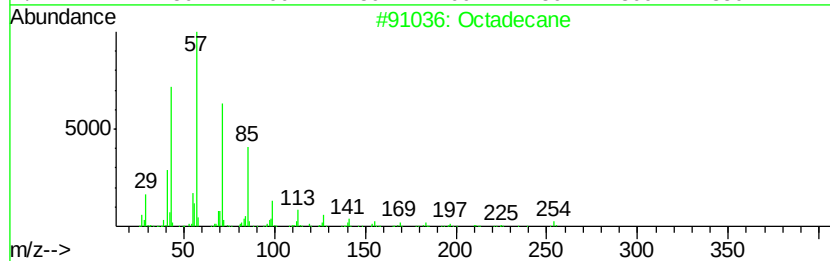
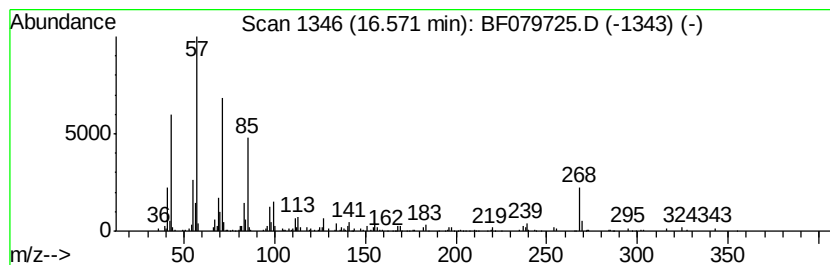
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Octadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.57	8.60 ng	518688	Perylene-d12	17.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	95
2		Nonadecane	268	C19H40	000629-92-5	87
3		Octacosane	394	C28H58	000630-02-4	81
4		Tetratetracontane	619	C44H90	007098-22-8	81
5		Tricosane	324	C23H48	000638-67-5	76



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
Data File : BF079725.D
Acq On : 5 Jun 2015 4:14
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Sample : G2474-05
Misc :
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Instrument :
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ClientSampleId :
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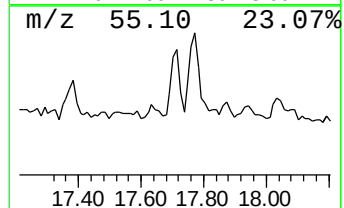
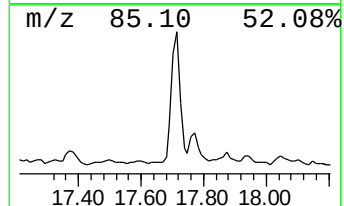
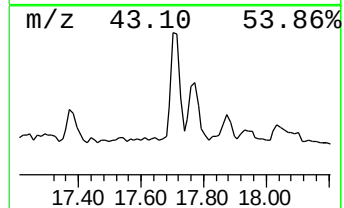
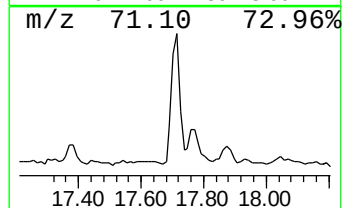
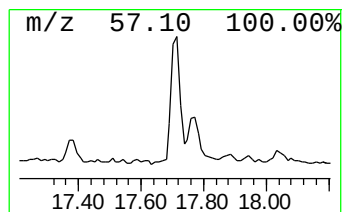
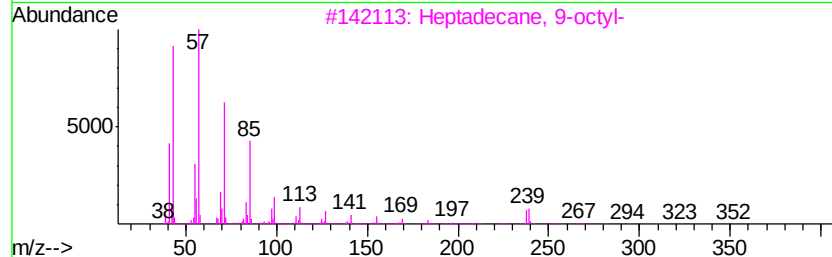
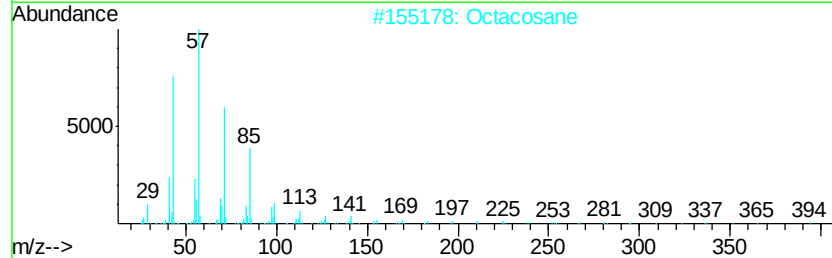
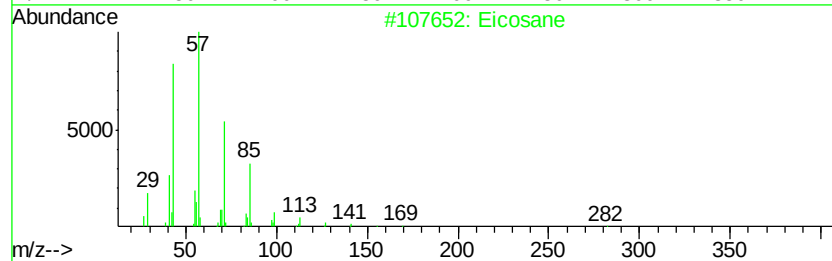
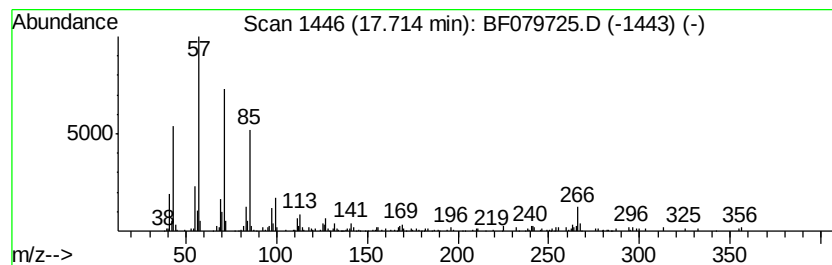
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Eicosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.71	3.98 ng	240234	Perylene-d12	17.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	87
2		Octacosane	394	C28H58	000630-02-4	83
3		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	80
4		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	80
5		Heneicosane	296	C21H44	000629-94-7	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
Data File : BF079725.D
Acq On : 5 Jun 2015 4:14
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Sample : G2474-05
Misc :
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Instrument :
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ClientSampleId :
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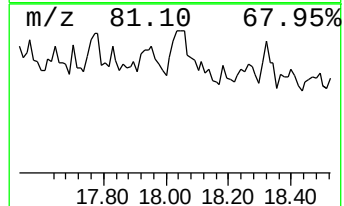
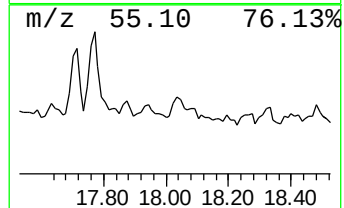
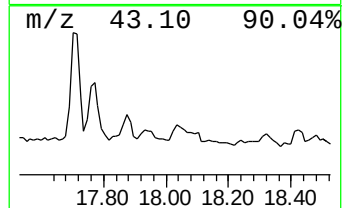
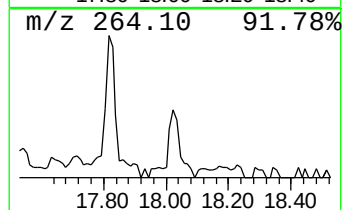
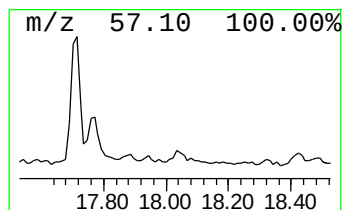
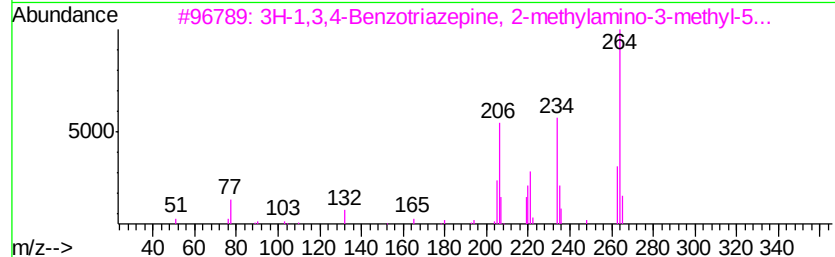
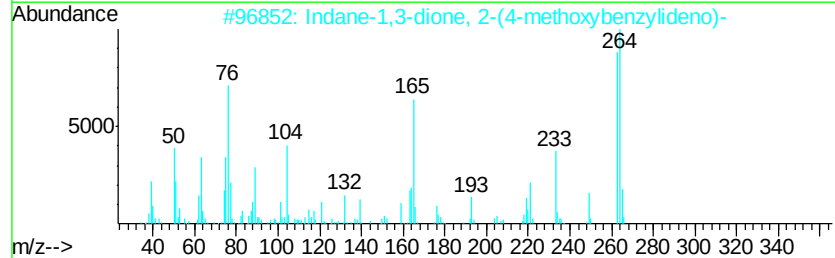
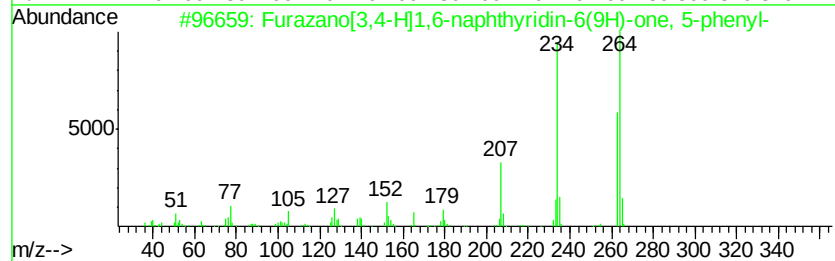
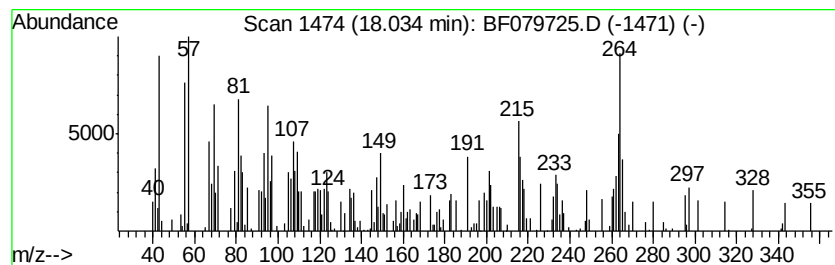
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown18.03 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.03	2.95 ng	178044	Perylene-d12	17.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Furazano[3,4-H]1,6-naphthyridin-...	264	C14H8N4O2	296245-19-7	38
2		Indane-1,3-dione, 2-(4-methoxybe...	264	C17H12O3	007421-76-3	38
3		3H-1,3,4-Benzotriazepine, 2-meth...	264	C16H16N4	121364-85-0	35
4		Benzo[b]naphtho[2,3-d]thiophene,...	264	C18H16S	024964-02-1	27
5		Pyrazolo[1,5-a]pyrimidin-7-amine...	264	C14H12N6	1000276-36-9	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
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Sample : G2474-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

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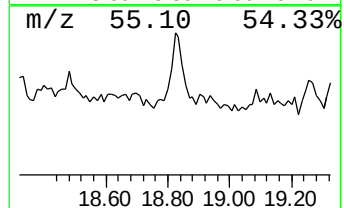
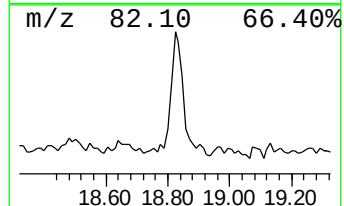
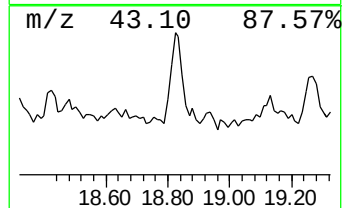
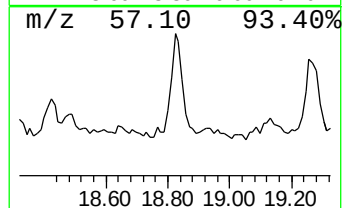
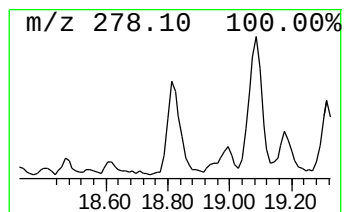
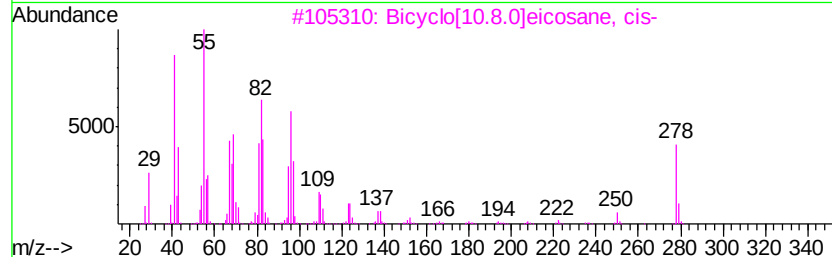
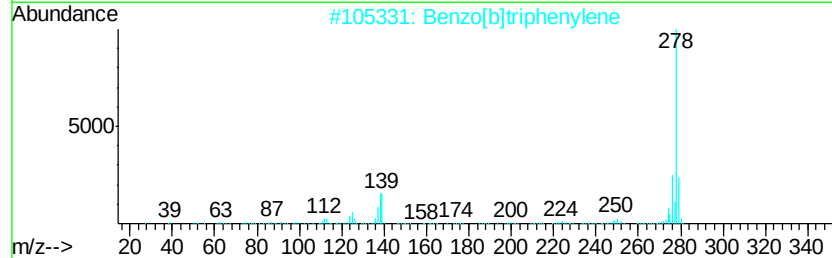
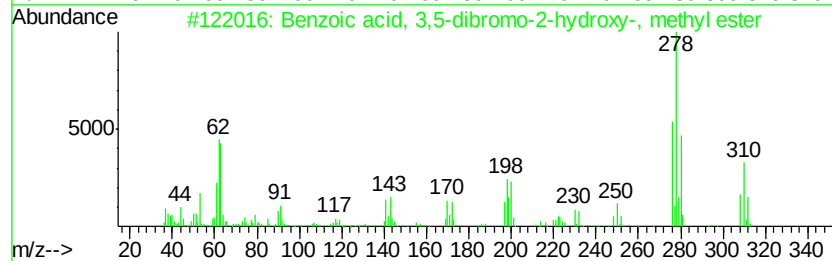
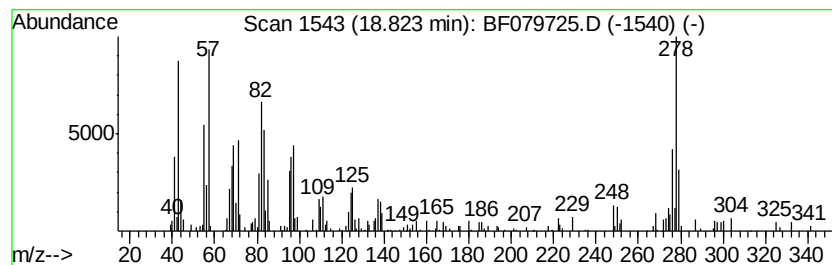
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 unknown18.82 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.82	3.90 ng	235189	Perylene-d12	17.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzoic acid, 3,5-dibromo-2-hydr...	308	C8H6Br2O3	021702-79-4	47
2		Benzo[b]triphenylene	278	C22H14	000215-58-7	46
3		Bicyclo[10.8.0]eicosane, cis-	278	C20H38	1000155-82-2	43
4		Pentaphene	278	C22H14	000222-93-5	38
5		Chromium, cycloheptatrienyl-(pen...	278	C17H22Cr	1000157-07-5	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
Data File : BF079725.D
Acq On : 5 Jun 2015 4:14
Operator : TP/IZ
Sample : G2474-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
C-5

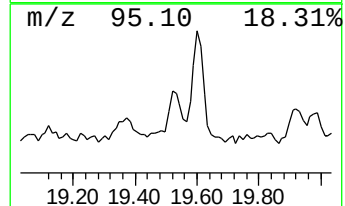
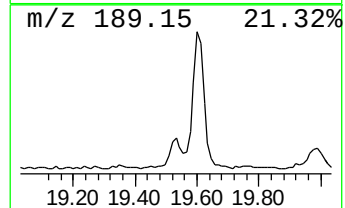
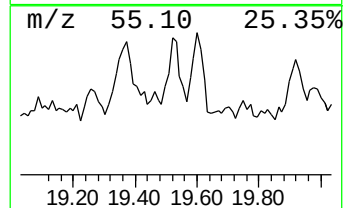
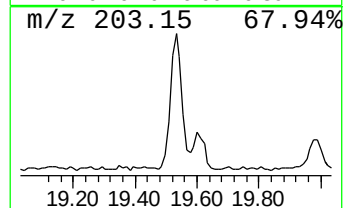
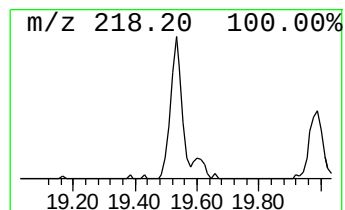
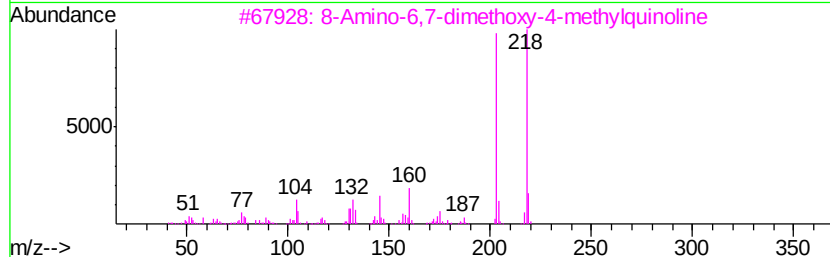
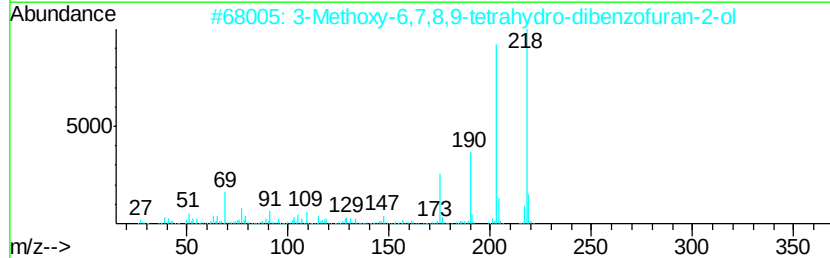
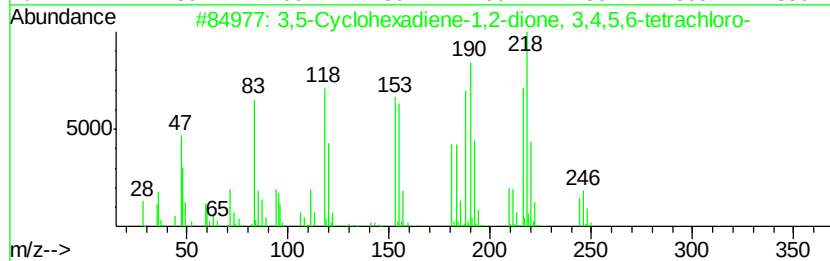
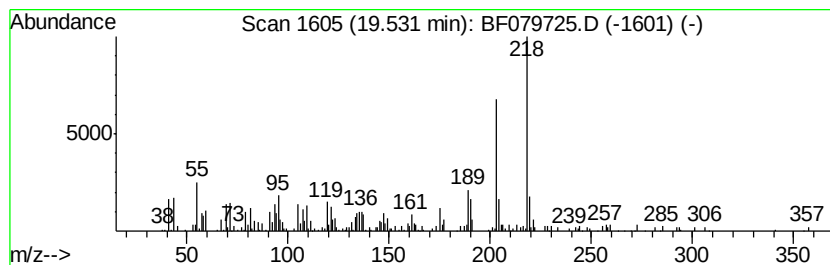
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 3,5-Cyclohexadiene-1,2-dione... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.53	4.86 ng	293317	Perylene-d12	17.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,5-Cyclohexadiene-1,2-dione, 3,...	244	C6Cl4O2	002435-53-2	86
2		3-Methoxy-6,7,8,9-tetrahydro-dib...	218	C13H14O3	1000186-57-9	74
3		8-Amino-6,7-dimethoxy-4-methylqu...	218	C12H14N2O2	1000213-07-2	58
4		.beta.-Amyrin trimethylsilyl ether	498	C33H58OSi	001721-67-1	58
5		2,9-Dimethyl-2,3,4,5,6,7-hexahyd...	203	C14H21N	077581-11-4	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
 Data File : BF079725.D
 Acq On : 5 Jun 2015 4:14
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 Sample : G2474-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C-5

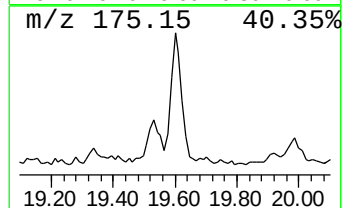
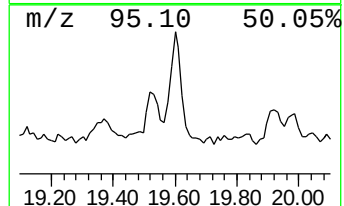
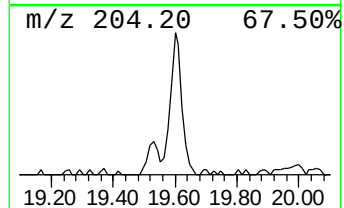
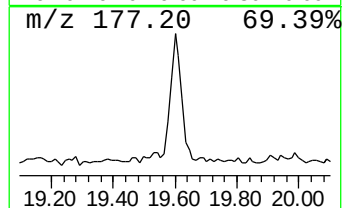
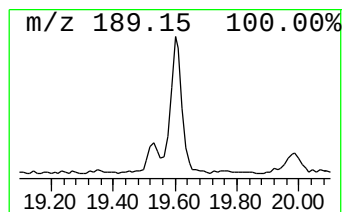
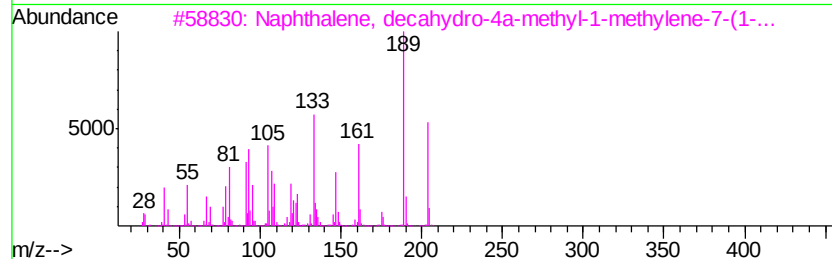
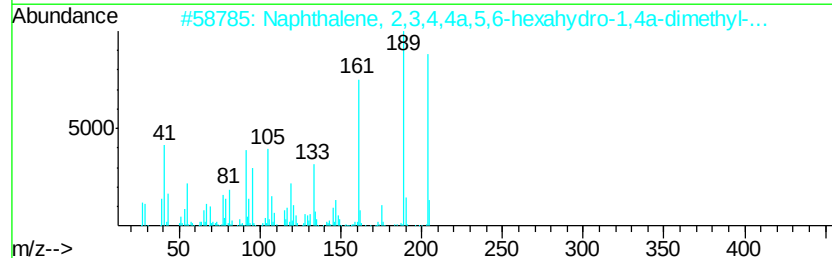
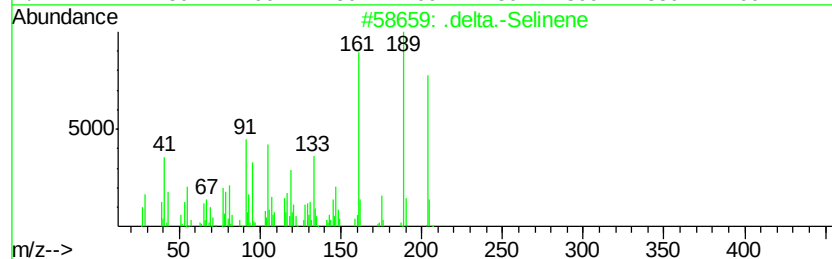
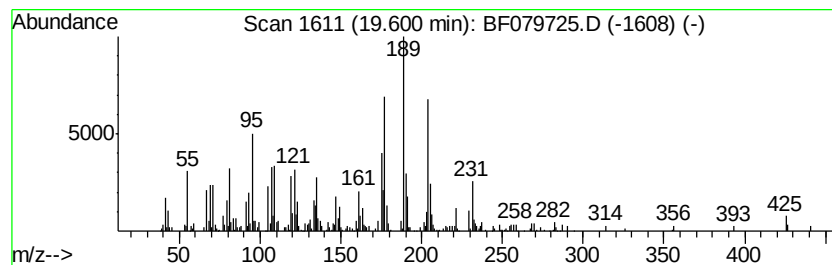
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 19 unknown19.60 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.60	7.50 ng	452096	Perylene-d12	17.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.delta.-Selinene	204	C15H24	028624-23-9	38
2		Naphthalene, 2,3,4,4a,5,6-hexahy...	204	C15H24	000473-14-3	38
3		Naphthalene, decahydro-4a-methyl...	204	C15H24	000515-17-3	38
4		2-Naphthalenemethanol, 1,2,3,4,4...	222	C15H26O	001209-71-8	35
5		Benzenepropanal, 3-(1,1-dimethyl...	204	C14H20O	062518-65-4	30



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Data File : BF079725.D
Acq On : 5 Jun 2015 4:14
Operator : TP/IZ
Sample : G2474-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

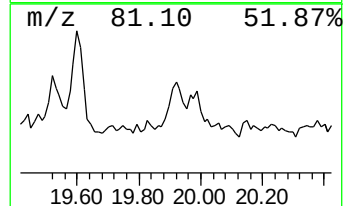
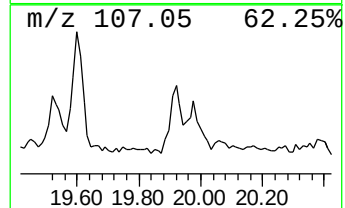
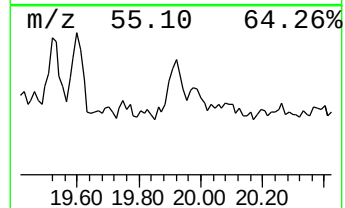
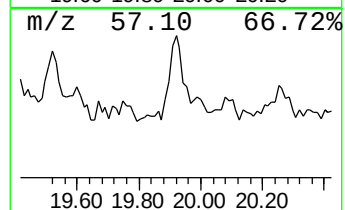
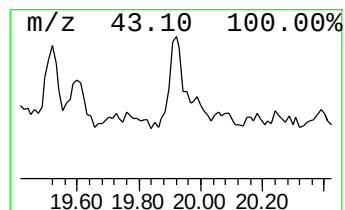
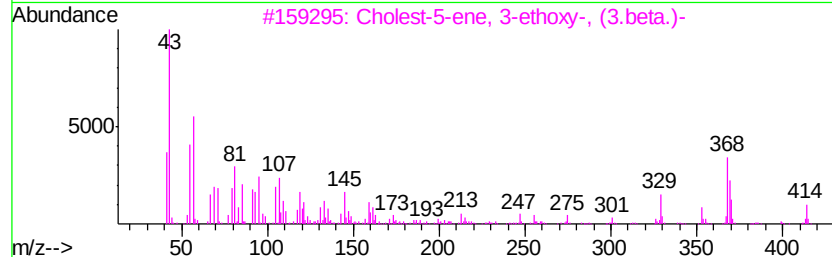
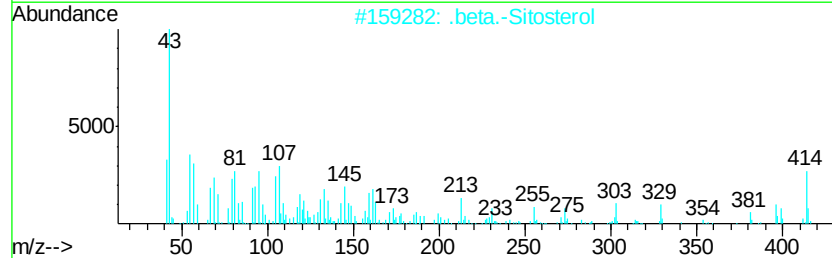
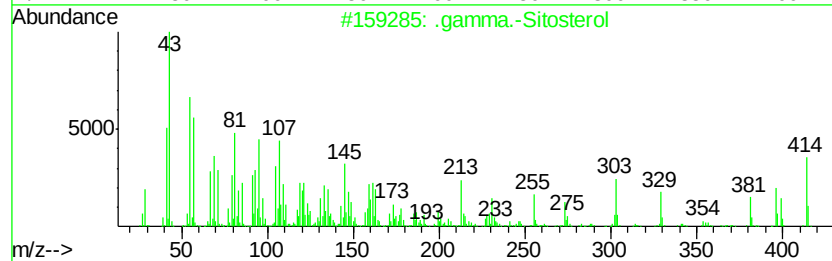
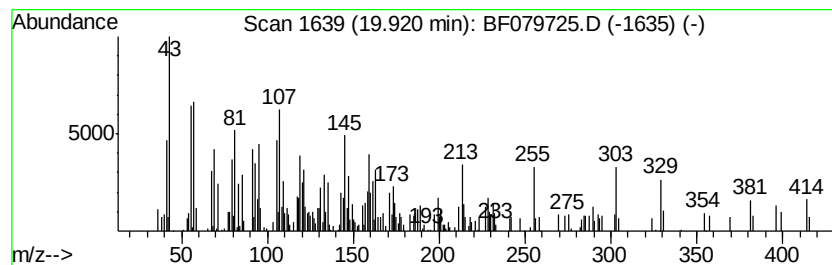
Instrument :
BNA_F
ClientSampleId :
C-5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 .gamma.-Sitosterol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.92	4.77 ng	287783	Perylene-d12	17.05	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.gamma.-Sitosterol	414	C29H50O	000083-47-6	94
2	.beta.-Sitosterol	414	C29H50O	000083-46-5	89
3	Cholest-5-ene, 3-ethoxy-, (3.beta...	414	C29H50O	000986-19-6	49
4	Stigmastan-3,5-diene	396	C29H48	1000214-16-4	35
5	17-(1,5-Dimethyl-hex-2-enyl)-10,...	382	C27H42O	1000194-79-4	33



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
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Sample : G2474-05
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
C-5

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060415.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