

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
 Data File : BF083630.D
 Acq On : 9 Dec 2015 9:09
 Operator : UM/IZ
 Sample : G4647-17
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BRIDGE14-9(0-2)

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-BF120815.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	3.538	160	162	175	rBV	426165	795622	23.75%	2.637%
2	3.984	199	201	216	rBV	2362746	3232121	96.47%	10.711%
3	5.276	311	314	322	rBV	2290796	3350375	100.00%	11.103%
4	5.402	322	325	336	rVB	2542899	3349035	99.96%	11.099%
5	5.699	349	351	360	rBV	713331	890400	26.58%	2.951%
6	5.916	367	370	382	rVB	1737170	2152379	64.24%	7.133%
7	6.339	404	407	415	rBV	19289	47523	1.42%	0.157%
8	6.533	421	424	438	rBV	1251231	2048228	61.13%	6.788%
9	7.607	515	518	528	rVB	656422	1085961	32.41%	3.599%
10	9.162	651	654	662	rVB	21305	42429	1.27%	0.141%
11	9.345	667	670	673	rVV	1873213	2344608	69.98%	7.770%
12	9.436	676	678	681	rVV	38939	55747	1.66%	0.185%
13	9.528	683	686	687	rVV	25087	35780	1.07%	0.119%
14	9.848	711	714	718	rVB4	20377	39389	1.18%	0.131%
15	10.031	727	730	733	rBV	477265	579686	17.30%	1.921%
16	10.213	744	746	751	rVB2	41921	61714	1.84%	0.205%
17	10.385	759	761	766	rBV2	653305	1150189	34.33%	3.812%
18	10.499	768	771	774	rVB	28505	53117	1.59%	0.176%
19	10.648	780	784	786	rBV3	25929	57093	1.70%	0.189%
20	10.808	792	798	799	rBV2	13986	37825	1.13%	0.125%
21	11.231	832	835	838	rBV	116513	150139	4.48%	0.498%
22	11.585	861	866	869	rBV2	30652	85951	2.57%	0.285%
23	11.688	872	875	881	rBV	886638	1473691	43.99%	4.884%
24	11.779	881	883	886	rVB	43361	53286	1.59%	0.177%
25	12.008	900	903	907	rBV	88550	186186	5.56%	0.617%
26	12.076	907	909	911	rVB	28115	40238	1.20%	0.133%
27	12.351	932	933	937	rVB4	18424	33722	1.01%	0.112%
28	12.739	965	967	969	rBV	45715	60599	1.81%	0.201%
29	12.785	969	971	978	rVB	633955	866031	25.85%	2.870%
30	13.711	1047	1052	1055	rBV3	18382	48486	1.45%	0.161%
31	13.768	1055	1057	1060	rVV3	14700	35380	1.06%	0.117%
32	13.837	1060	1063	1071	rVV	99164	173559	5.18%	0.575%
33	14.214	1092	1096	1101	rBV	21976	52109	1.56%	0.173%
34	14.351	1105	1108	1111	rBV2	24049	42680	1.27%	0.141%

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Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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35	14.557	1124	1126	1129	rBV	44740	69819	2.08%	0.231%
36	14.762	1140	1144	1147	rBV2	18229	33870	1.01%	0.112%
37	14.865	1150	1153	1155	rVB	92512	121838	3.64%	0.404%
38	15.117	1172	1175	1178	rBV	88979	134355	4.01%	0.445%
39	15.254	1185	1187	1189	rVB2	29949	39418	1.18%	0.131%
40	15.414	1198	1201	1204	rBV	1137553	1518834	45.33%	5.033%
41	15.700	1223	1226	1230	rVB2	30457	59079	1.76%	0.196%
42	16.225	1266	1272	1273	rBV2	32500	90766	2.71%	0.301%
43	16.294	1277	1278	1283	rVB4	28327	42560	1.27%	0.141%
44	16.397	1285	1287	1289	rBV2	24592	44089	1.32%	0.146%
45	16.500	1294	1296	1299	rVB	37367	59147	1.77%	0.196%
46	16.557	1299	1301	1305	rBV3	19759	37805	1.13%	0.125%
47	16.774	1318	1320	1325	rVB	49599	80804	2.41%	0.268%
48	16.923	1331	1333	1337	rBV2	187220	344613	10.29%	1.142%
49	17.003	1337	1340	1344	rVV	494984	793512	23.68%	2.630%
50	17.357	1369	1371	1374	rBV2	37881	67829	2.02%	0.225%
51	17.746	1402	1405	1408	rBV	182032	277210	8.27%	0.919%
52	18.409	1461	1463	1468	rBV2	89761	239961	7.16%	0.795%
53	18.637	1482	1483	1489	rVB	300827	451698	13.48%	1.497%
54	19.357	1544	1546	1555	rVB6	123287	287479	8.58%	0.953%
55	20.877	1677	1679	1686	rVB4	145626	342780	10.23%	1.136%
56	21.860	1762	1765	1769	rVB2	128938	326059	9.73%	1.081%

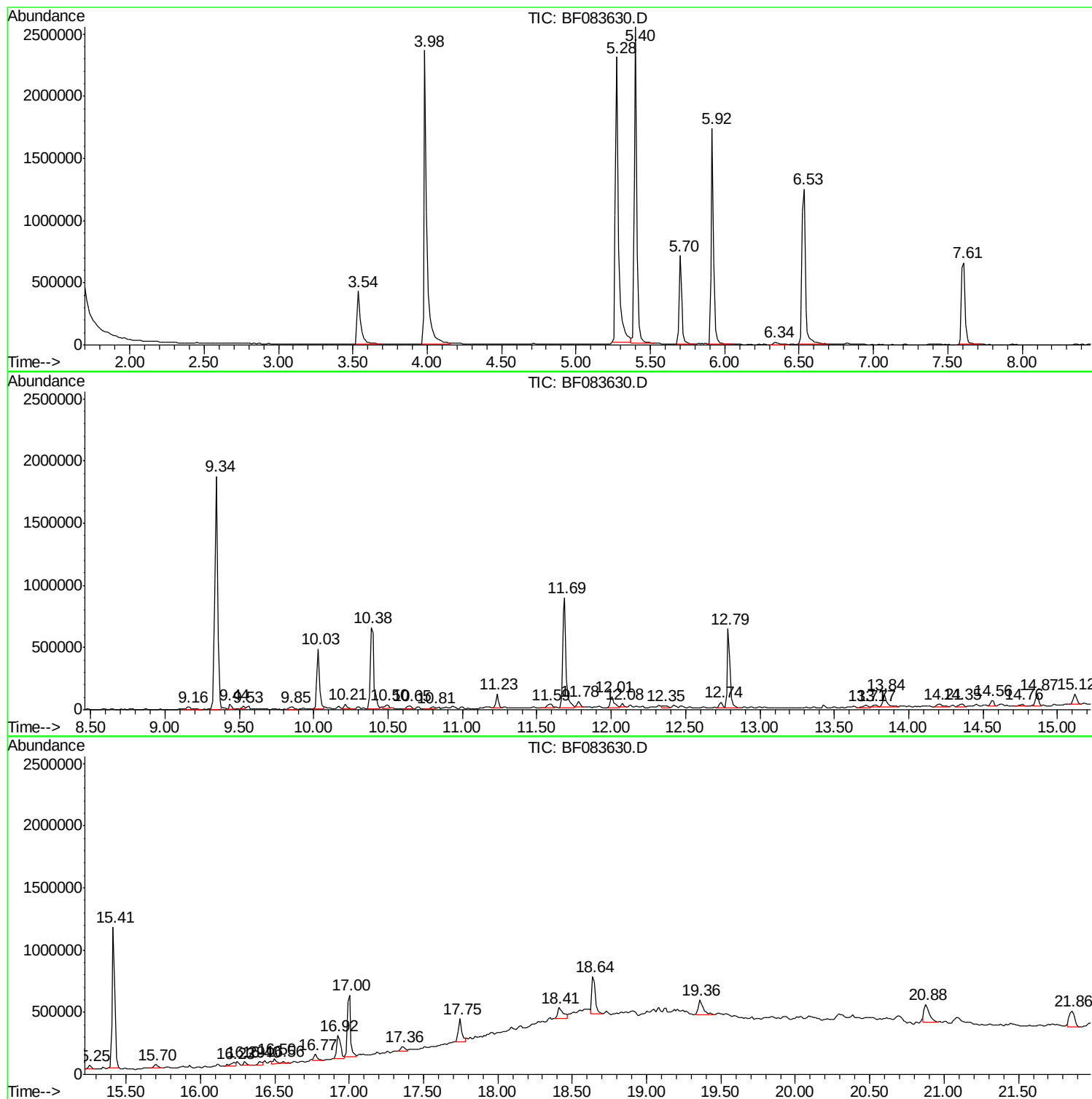
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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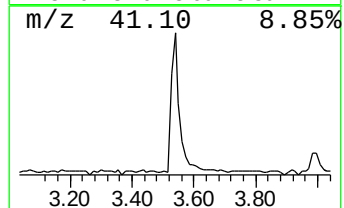
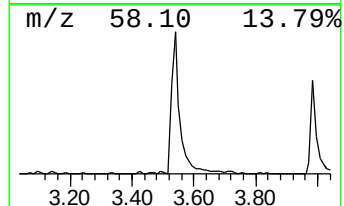
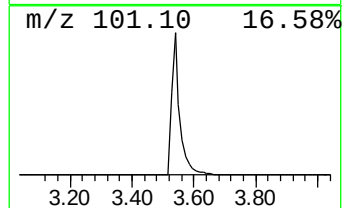
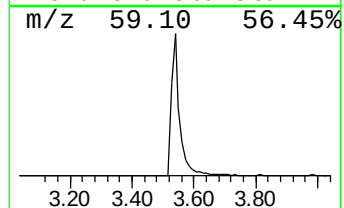
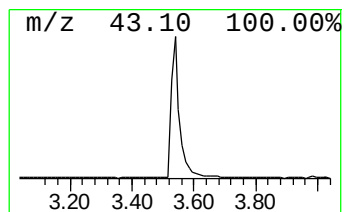
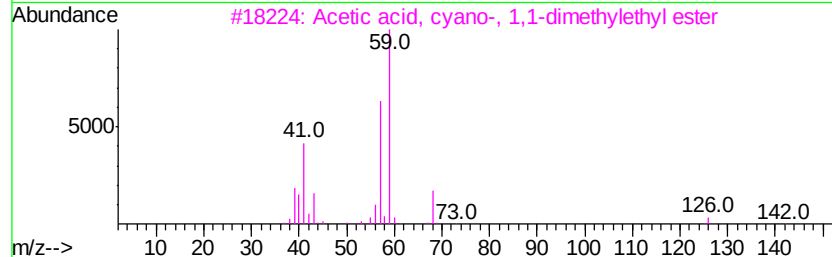
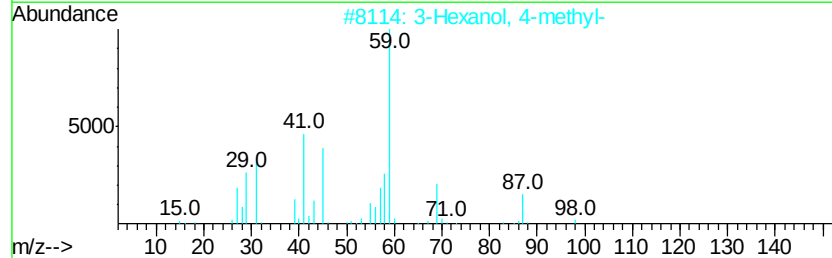
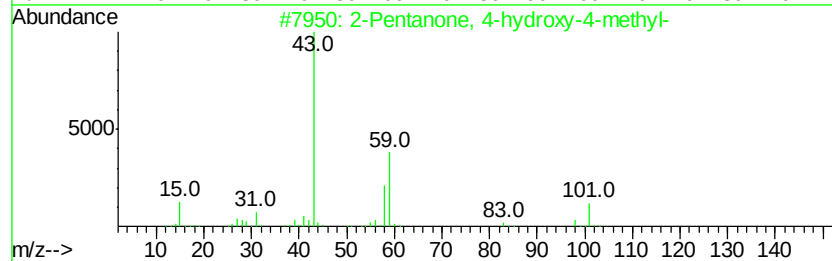
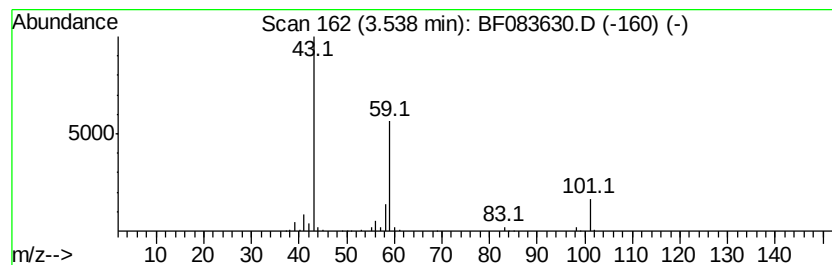
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.54	17.87 ng	795622	1,4-Dichlorobenzene-d4	5.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Ethanol, pentamethyl-	116	C7H16O	000594-83-2	9



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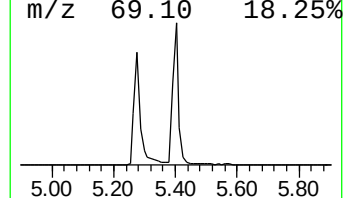
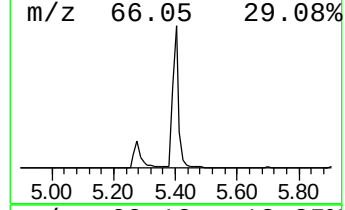
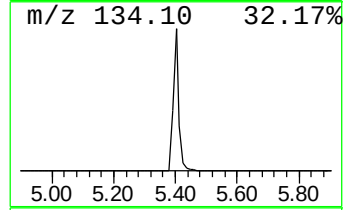
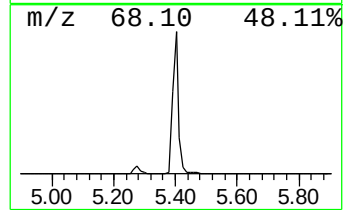
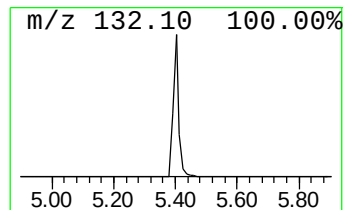
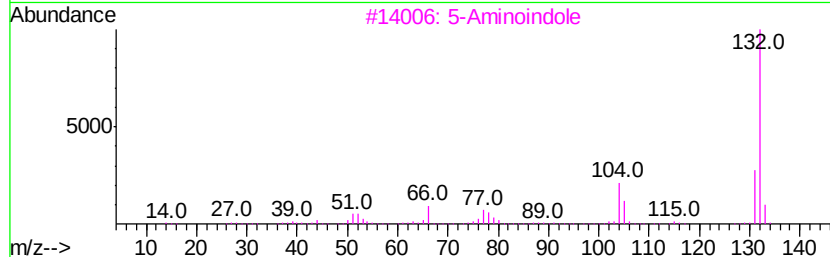
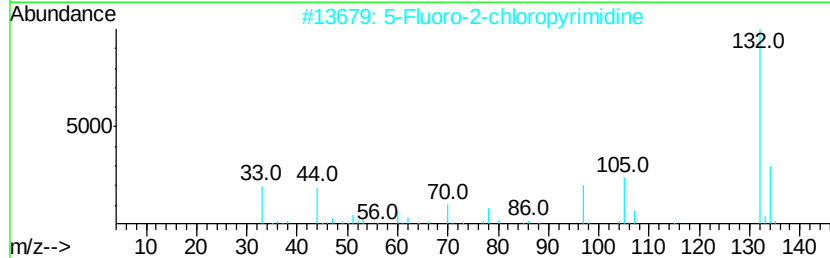
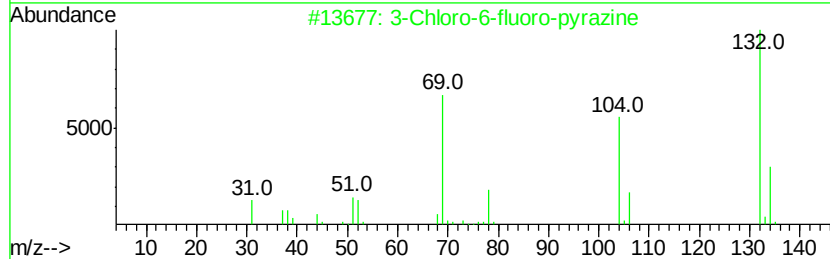
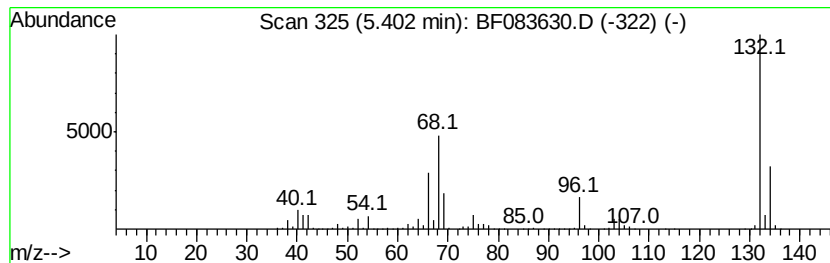
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown5.40 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.40	75.23 ng	3349040	1,4-Dichlorobenzene-d4	5.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Xenon	132	Xe	007440-63-3	9



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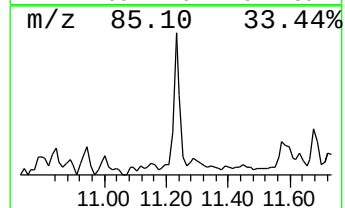
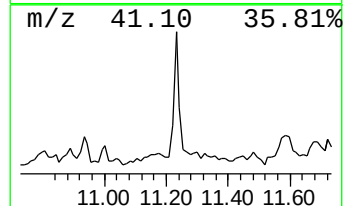
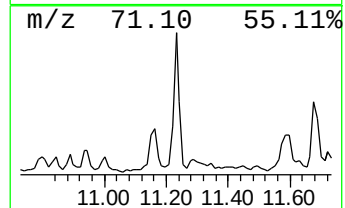
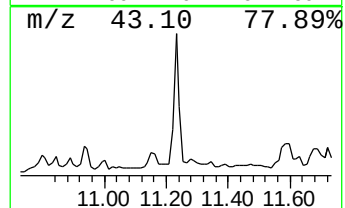
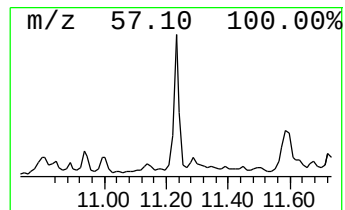
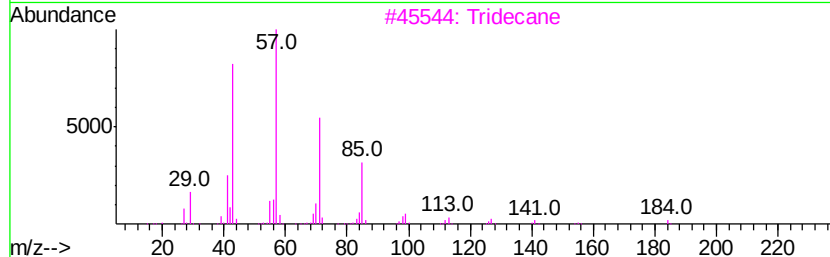
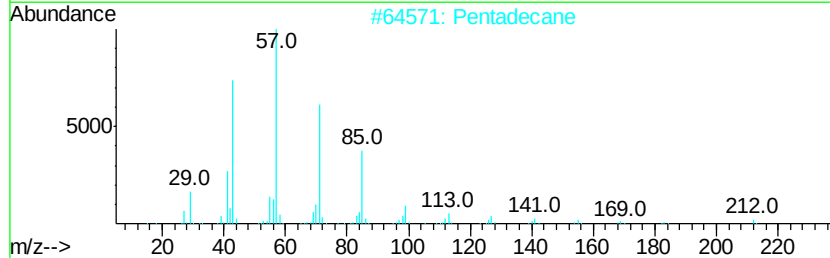
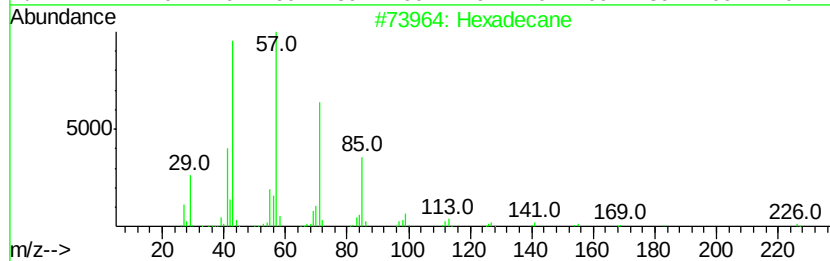
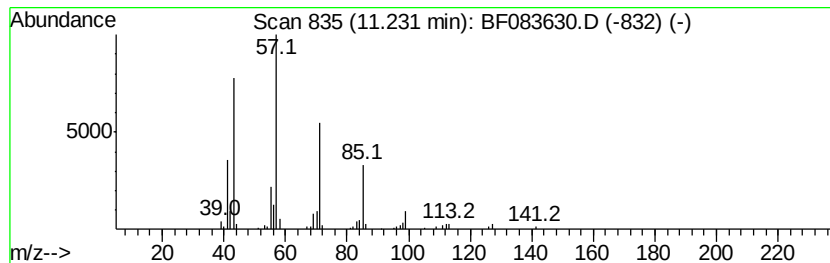
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.23	2.61 ng	150139	Acenaphthene-d10	10.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	91
2		Pentadecane	212	C15H32	000629-62-9	86
3		Tridecane	184	C13H28	000629-50-5	86
4		Tetradecane	198	C14H30	000629-59-4	80
5		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	59



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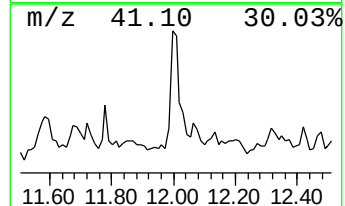
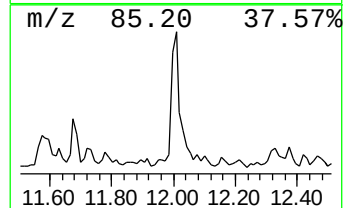
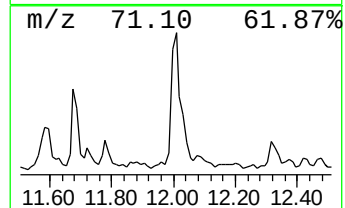
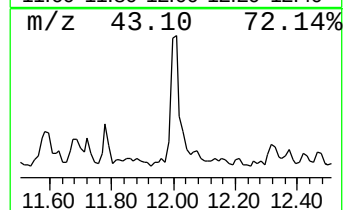
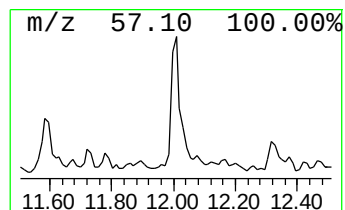
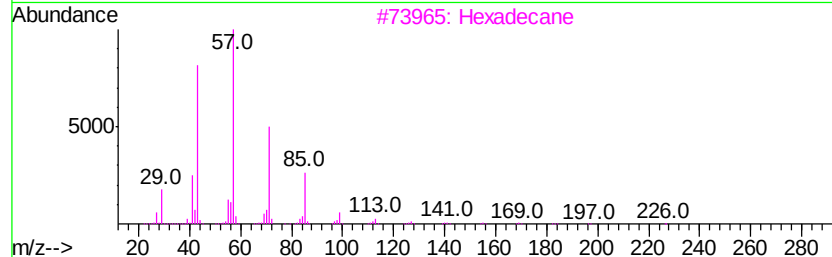
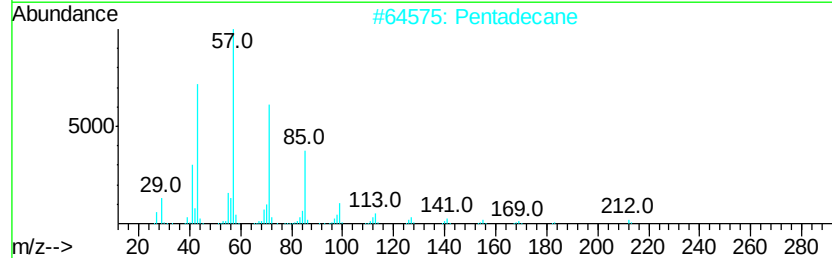
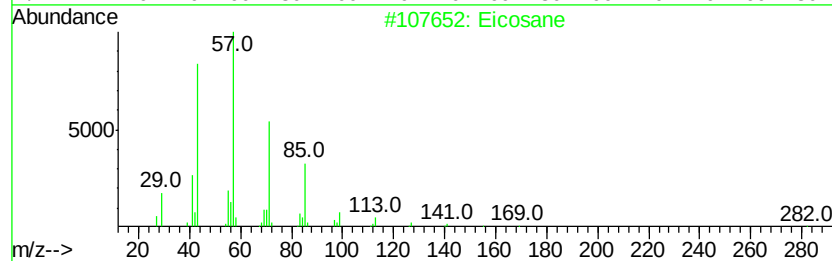
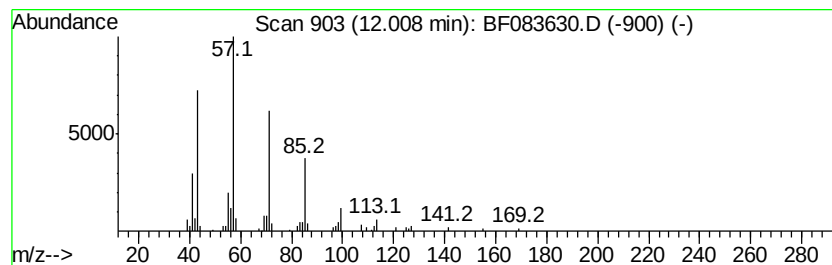
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Eicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	4.30 ng	186186	Phenanthrene-d10	12.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Pentadecane	212	C15H32	000629-62-9	86
3		Hexadecane	226	C16H34	000544-76-3	83
4		Octadecane	254	C18H38	000593-45-3	78
5		Docosane	310	C22H46	000629-97-0	78



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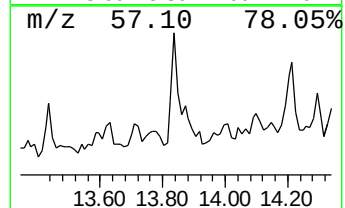
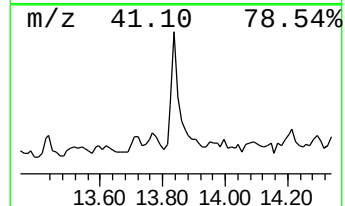
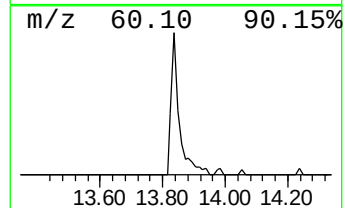
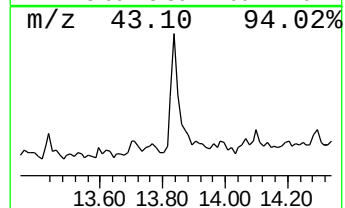
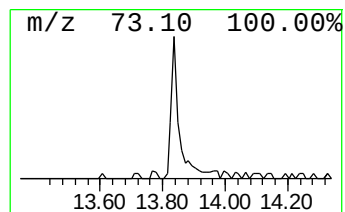
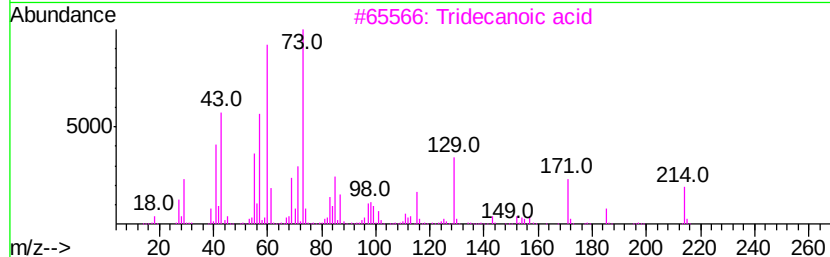
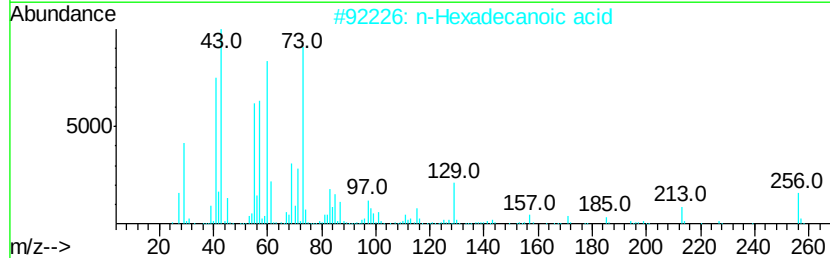
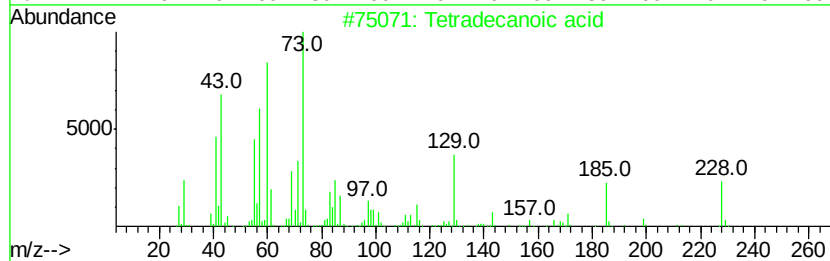
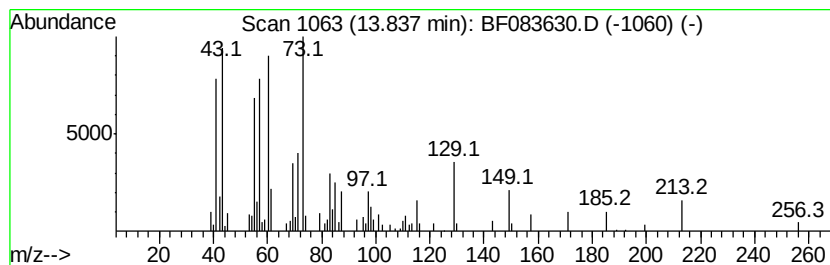
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tetradecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.84	4.01 ng	173559	Phenanthrene-d10	12.79	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanoic acid	228	C14H28O2	000544-63-8	80
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	80
3	Tridecanoic acid	214	C13H26O2	000638-53-9	80
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	80
5	Undecanoic acid	186	C11H22O2	000112-37-8	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
Data File : BF083630.D
Acq On : 9 Dec 2015 9:09
Operator : UM/IZ
Sample : G4647-17
Misc :
ALS Vial : 38 Sample Multiplier: 1

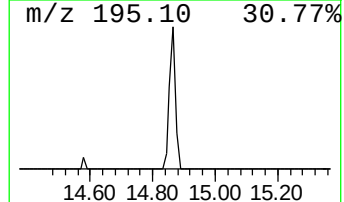
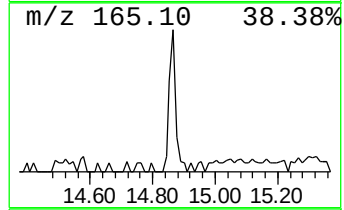
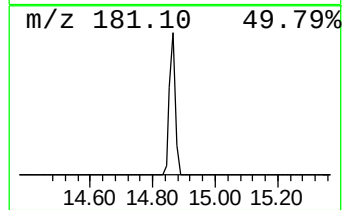
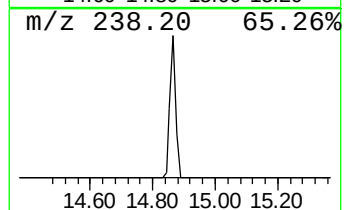
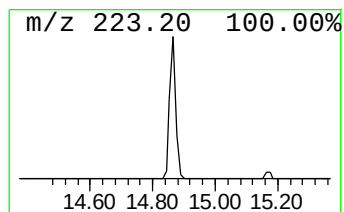
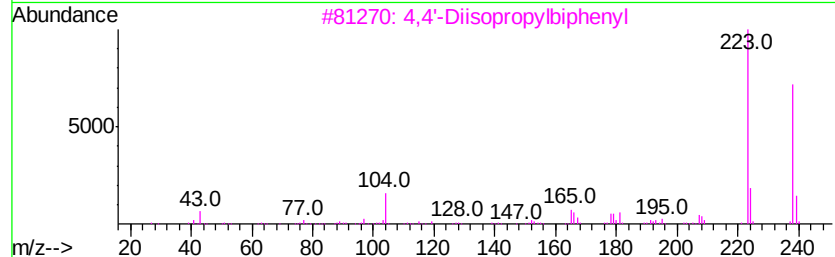
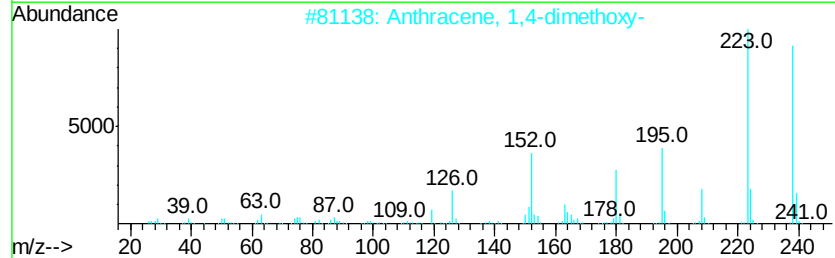
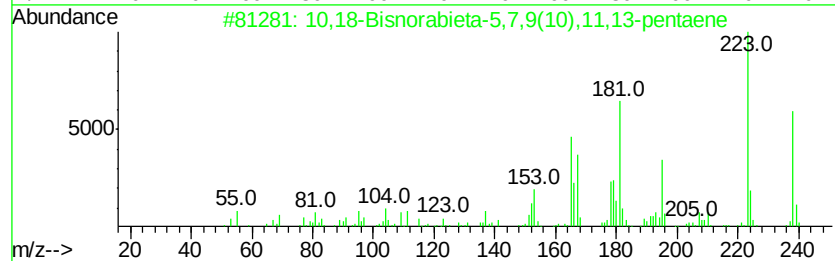
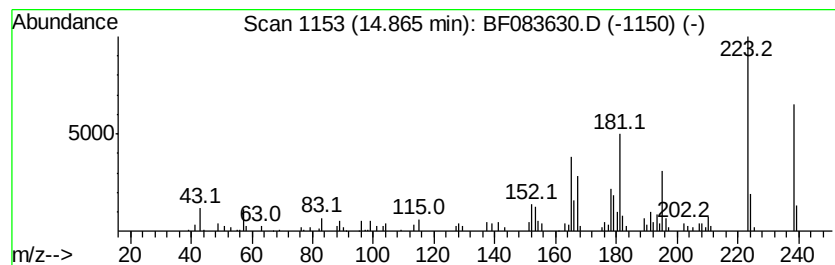
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
14.87	2.81 ng	121838	Phenanthrene-d10		12.79
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2	Anthracene, 1,4-dimethoxy-	238	C16H14O2	013076-29-4	53
3	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	49
4	3,3'-Diacetylbiphenyl	238	C16H14O2	094113-07-2	46
5	4,4'-Diacetyl biphenyl	238	C16H14O2	000787-69-9	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
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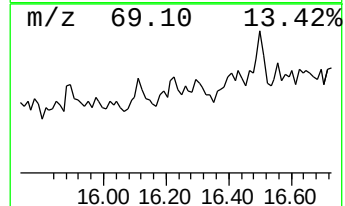
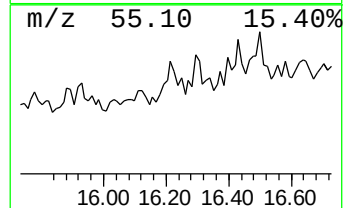
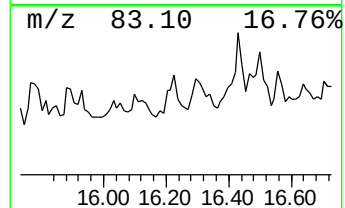
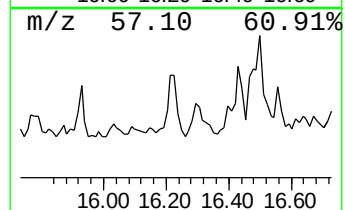
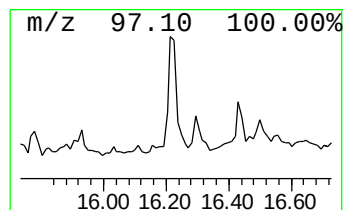
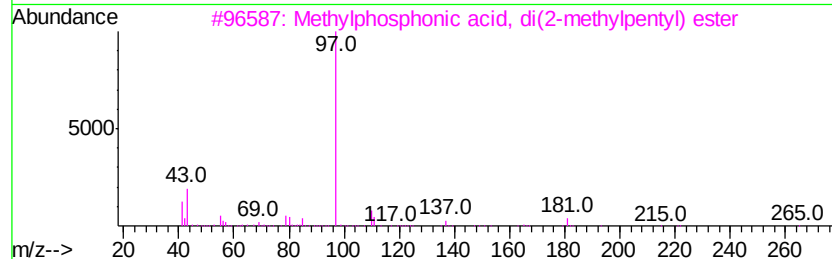
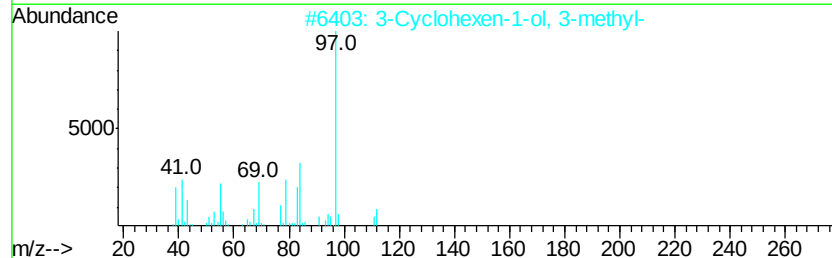
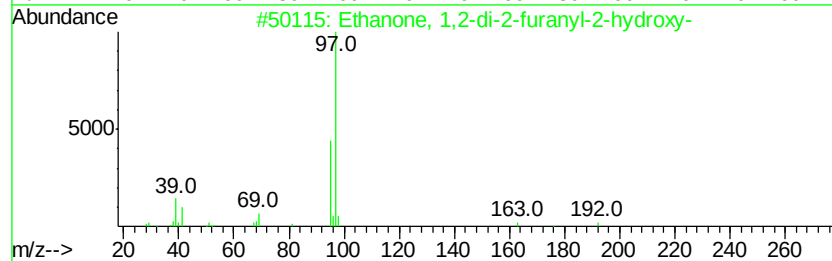
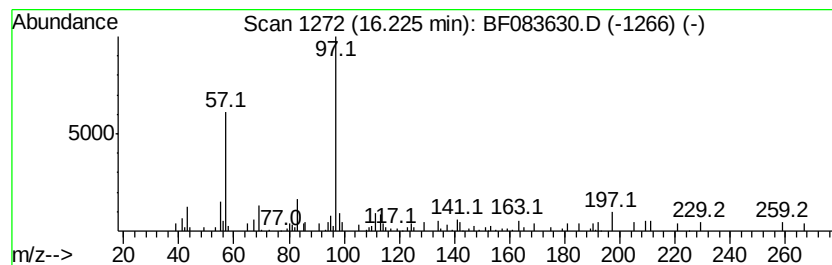
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Ethanone, 1,2-di-2-furanyl-... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.23	2.29 ng	90766	Chrysene-d12	17.00
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Ethanone, 1,2-di-2-furanyl-2-hyd...	192 C10H8O4	000552-86-3	50
2	3-Cyclohexen-1-ol, 3-methyl-	112 C7H12O	053783-91-8	42
3	Methylphosphonic acid, di(2-meth...	264 C13H29O3P	1000273-18-3	40
4	Isoxazole, 3,5-dimethyl-	97 C5H7NO	000300-87-8	40
5	Cyclohexanone, 3-(3,3-dimethylbu...	182 C12H22O	040564-94-1	37



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
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Operator : UM/IZ
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Instrument :
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ClientSampleId :
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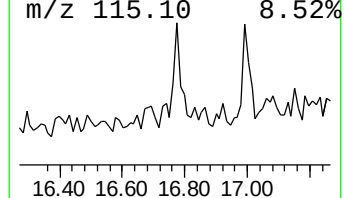
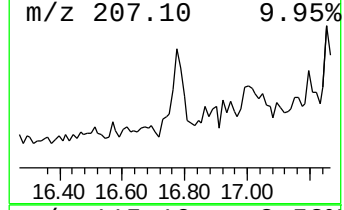
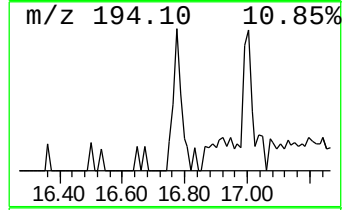
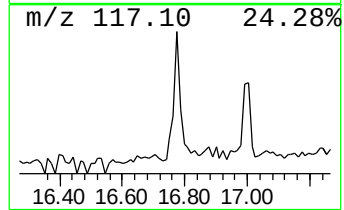
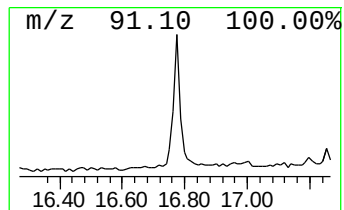
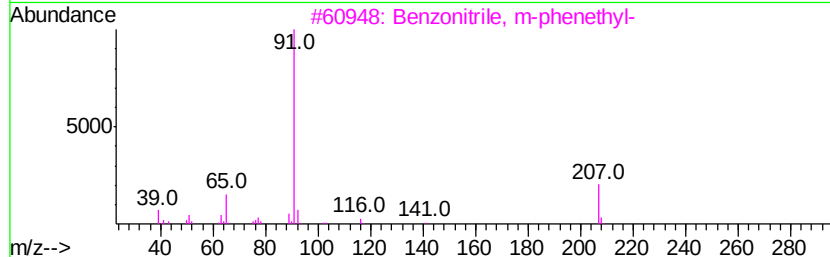
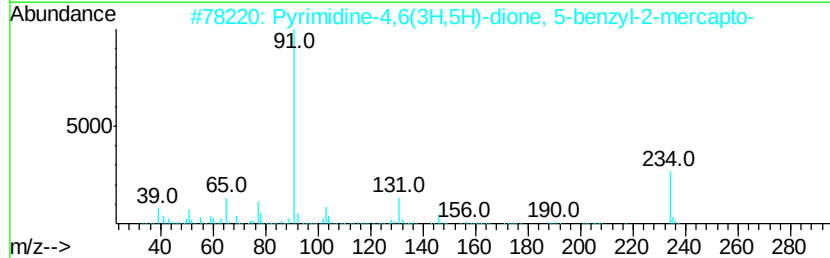
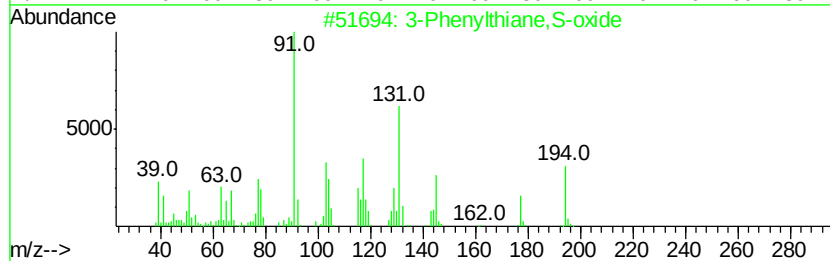
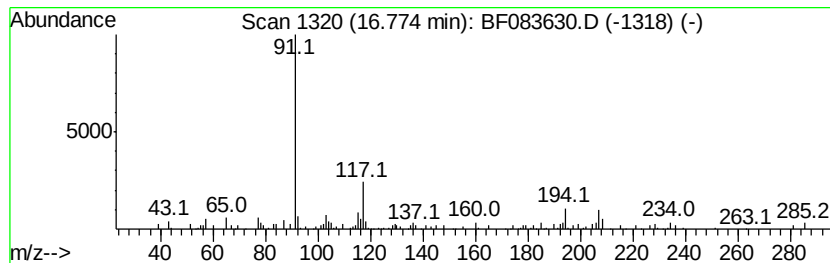
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown16.77 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.77	2.04 ng	80804	Chrysene-d12	17.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Phenylthiane,S-oxide	194	C11H14OS	058121-26-9	17
2		Pyrimidine-4,6(3H,5H)-dione, 5-b...	234	C11H10N2O2S	015018-53-8	14
3		Benzonitrile, m-phenethyl-	207	C15H13N	034176-91-5	12
4		2-Pentenal, 5-phenyl-	160	C11H12O	033046-84-3	10
5		Benzene, 2-benzyloxy-1-methoxy-4...	285	C16H15NO4	063909-29-5	10



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
Data File : BF083630.D
Acq On : 9 Dec 2015 9:09
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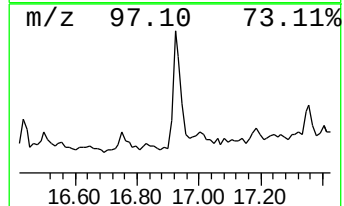
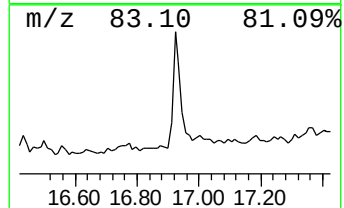
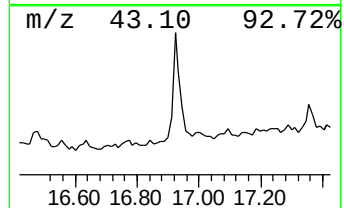
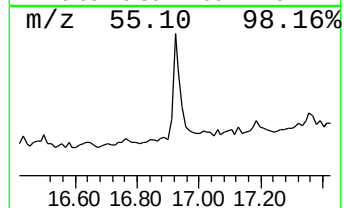
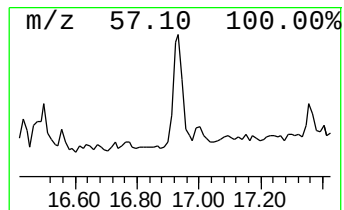
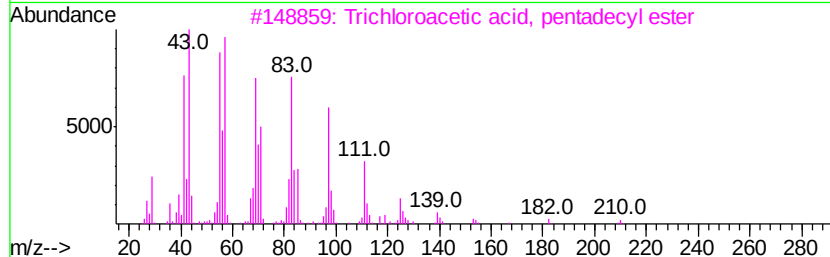
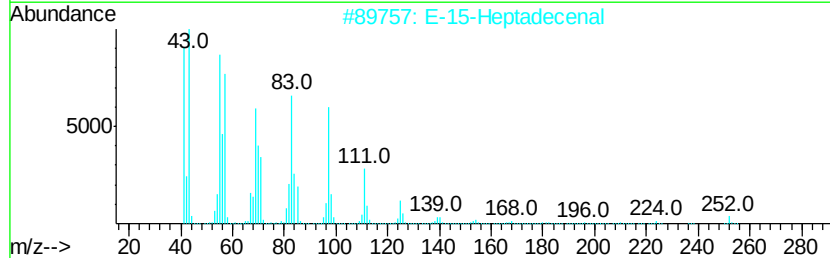
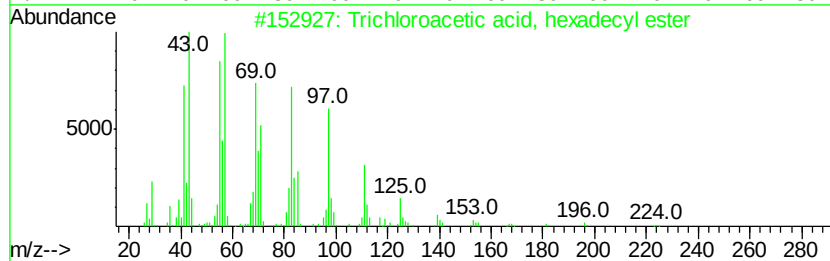
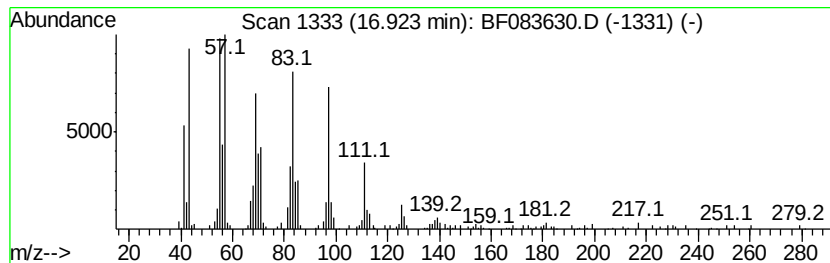
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Trichloroacetic acid, hexad... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.92	8.69 ng	344613	Chrysene-d12	17.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
2			E-15-Heptadecenal	252	C17H32O	1000130-97-9	91
3			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
4			1-Nonadecene	266	C19H38	018435-45-5	91
5			2- Chloropropionic acid, hexadec...	332	C19H37ClO2	086711-81-1	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
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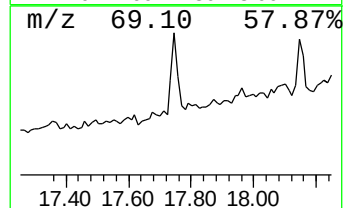
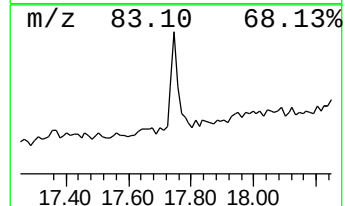
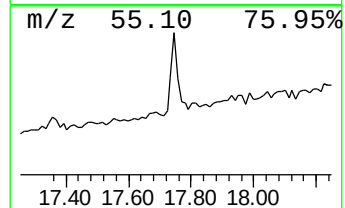
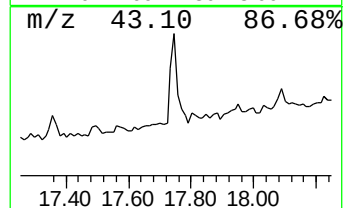
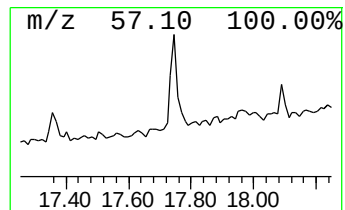
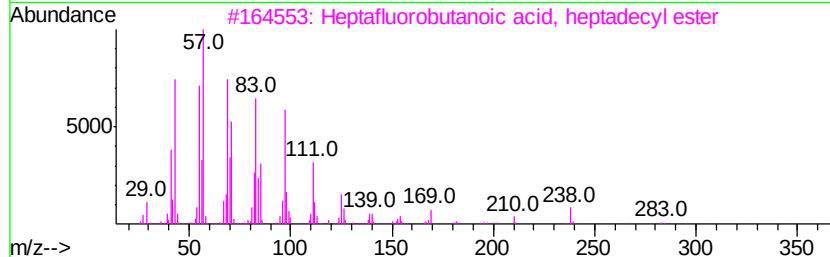
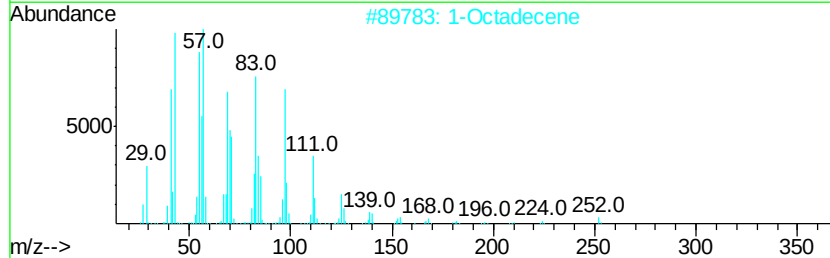
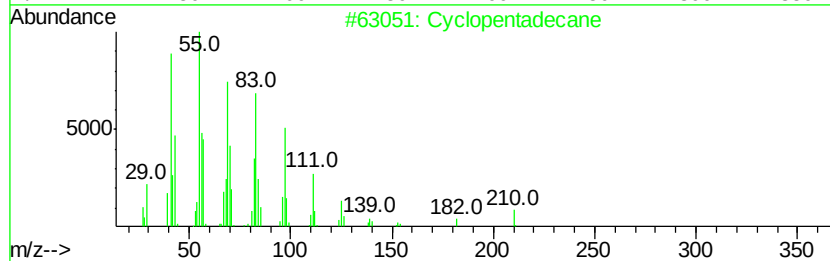
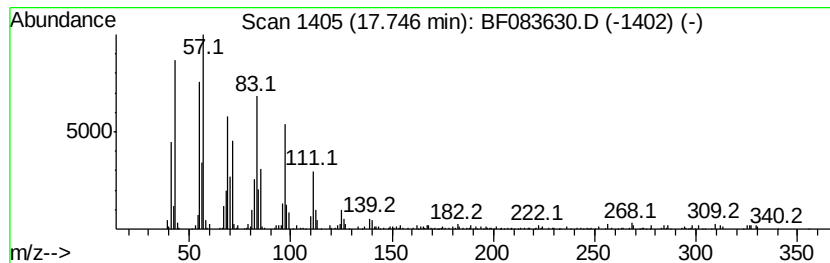
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Cyclopentadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.75	6.99 ng	277210	Chrysene-d12	17.00		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentadecane	210	C15H30	000295-48-7	97
2		1-Octadecene	252	C18H36	000112-88-9	95
3		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94
4		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	94
5		Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
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Acq On : 9 Dec 2015 9:09
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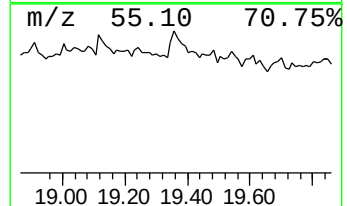
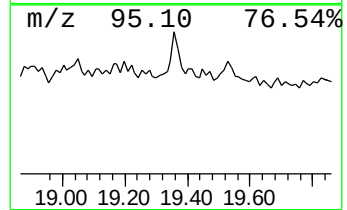
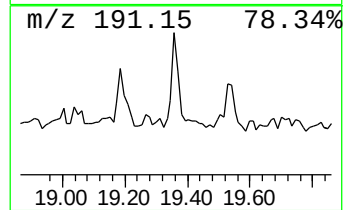
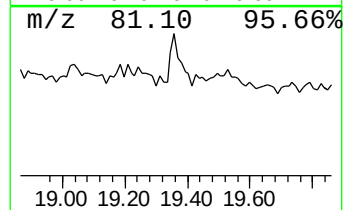
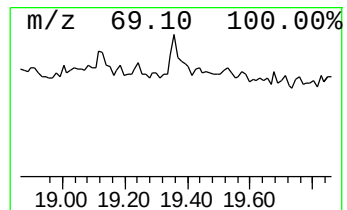
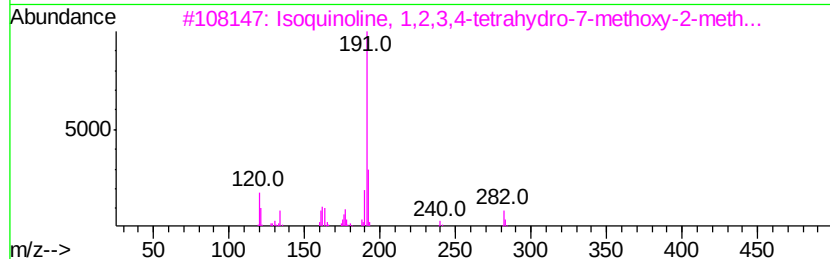
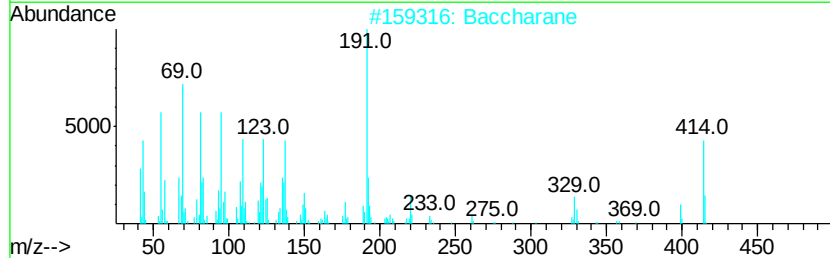
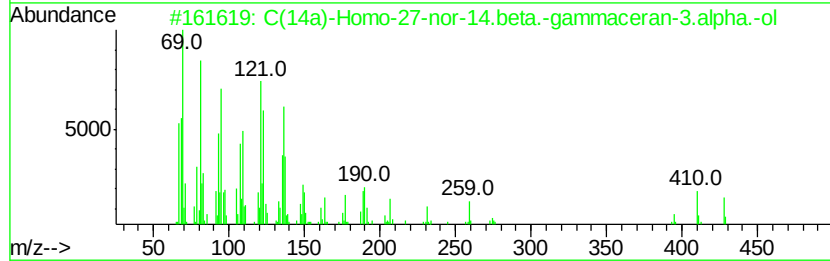
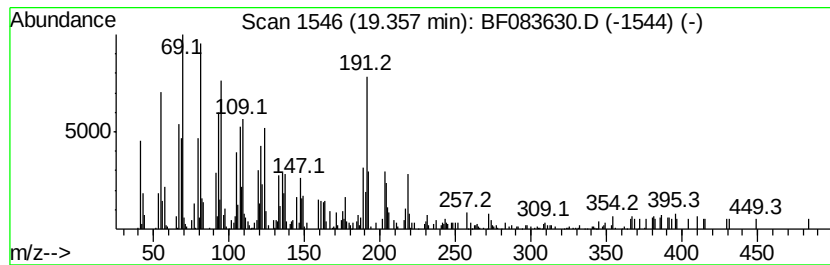
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 C(14a)-Homo-27-nor-14.beta.... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.36	12.73 ng	287479	Perylene-d12	18.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	C(14a)-Homo-27-nor-14.beta.-gamm...	428	C30H52O	024739-08-0	59
2		Baccharane	414	C30H54	036441-74-4	49
3		Isoquinoline, 1,2,3,4-tetrahydro...	283	C18H21N02	036646-87-4	49
4		A'-Neogammacer-22(29)-ene	410	C30H50	001615-91-4	46
5		Cyclopentane, 1-methyl-1-(2-meth...	138	C10H18	074764-47-9	41



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
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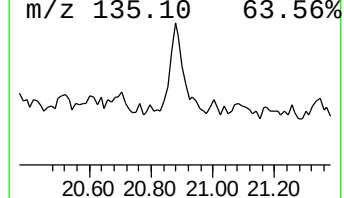
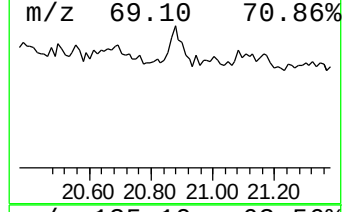
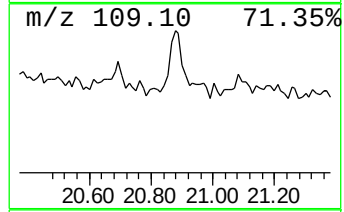
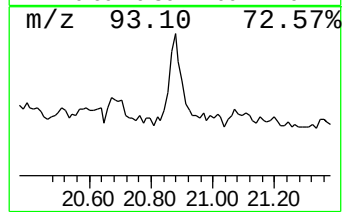
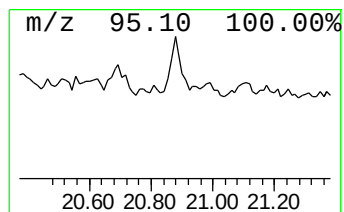
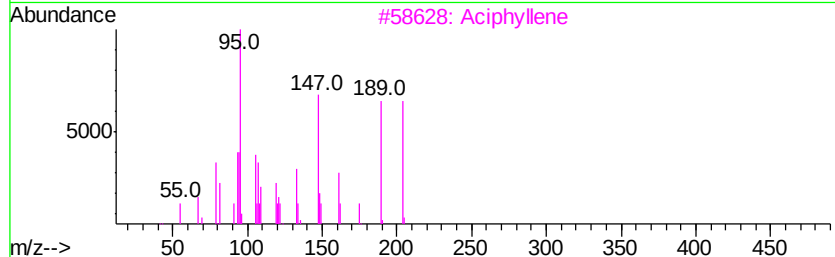
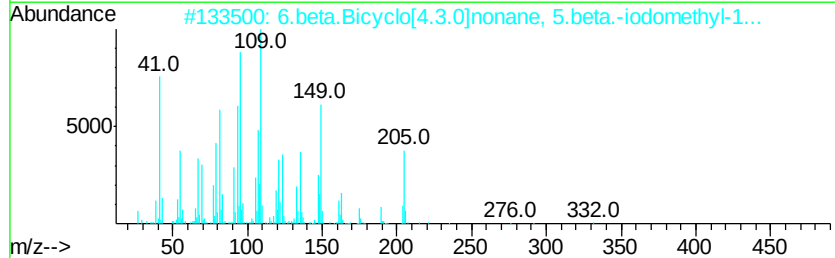
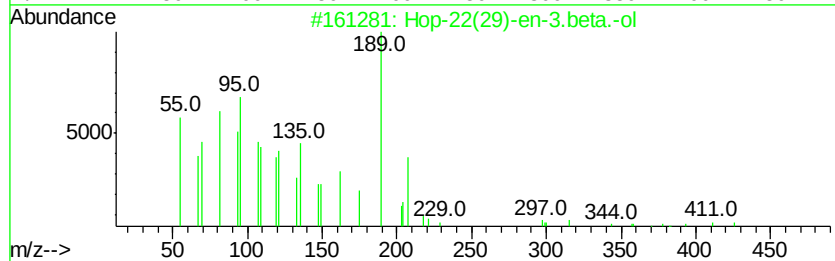
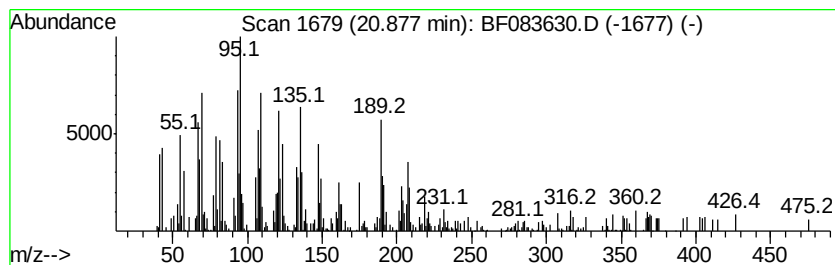
Instrument :
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ClientSampleId :
BRIDGE14-9(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Hop-22(29)-en-3.beta.-ol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.88	15.18 ng	342780	Perylene-d12	18.64
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Hop-22(29)-en-3.beta.-ol	426 C30H50O	058801-23-3 83
2		6.beta.Bicyclo[4.3.0]nonane, 5.b...	332 C15H25I	1000195-85-9 46
3		Aciphyllene	204 C15H24	087745-31-1 46
4		Cholest-8(14)-en-3-ol, 2,2-dimet...	414 C29H50O	064519-14-8 43
5		1-Naphthalenecarboxylic acid, de...	316 C21H32O2	001235-39-8 40



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
Data File : BF083630.D
Acq On : 9 Dec 2015 9:09
Operator : UM/IZ
Sample : G4647-17
Misc :
ALS Vial : 38 Sample Multiplier: 1

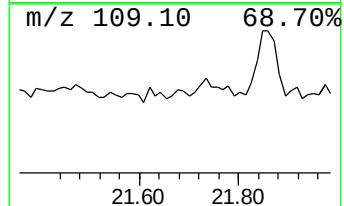
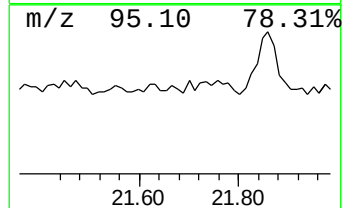
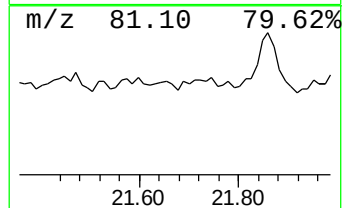
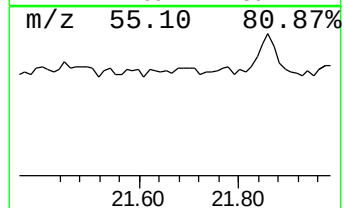
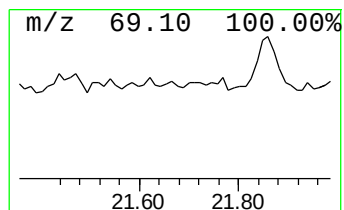
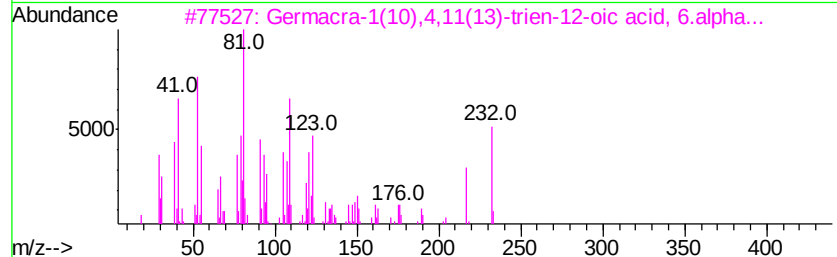
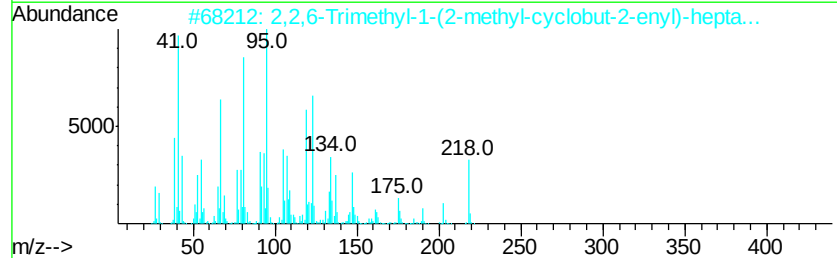
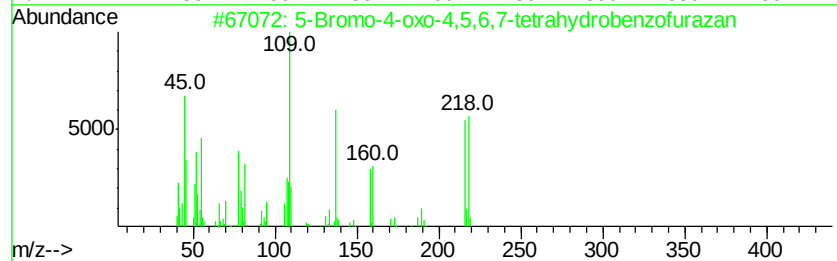
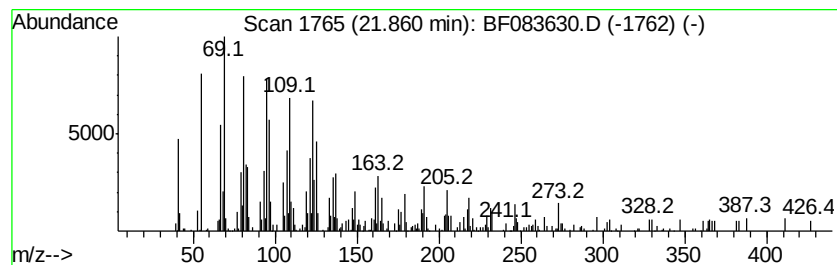
Instrument :
BNA_F
ClientSampled :
BRIDGE14-9(0-2)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.86	14.44 ng	326059	Perylene-d12	18.64	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	90
2	2,2,6-Trimethyl-1-(2-methyl-cycl...	218	C15H22O	1000188-72-8	46
3	Germacra-1(10),4,11(13)-trien-12...	232	C15H20O2	000553-21-9	46
4	1-Ethynylcyclopentanol	110	C7H10O	017356-19-3	38
5	Sesquirosefuran	218	C15H22O	039007-93-7	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
Data File : BF083630.D
Acq On : 9 Dec 2015 9:09
Operator : UM/IZ
Sample : G4647-17
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
BRIDGE14-9(0-2)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.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...	3.54	17.9	ng	795622	1	5.70	890400	20.0
unknown5.40	5.40	75.2	ng	3349040	1	5.70	890400	20.0
Hexadecane	11.23	2.6	ng	150139	3	10.38	1150190	20.0
Eicosane	12.01	4.3	ng	186186	4	12.79	866031	20.0
Tetradecanoic acid	13.84	4.0	ng	173559	4	12.79	866031	20.0
10,18-Bisnorabiet...	14.87	2.8	ng	121838	4	12.79	866031	20.0
Ethanone, 1,2-di-...	16.23	2.3	ng	90766	5	17.00	793512	20.0
unknown16.77	16.77	2.0	ng	80804	5	17.00	793512	20.0
Trichloroacetic a...	16.92	8.7	ng	344613	5	17.00	793512	20.0
Cyclopentadecane	17.75	7.0	ng	277210	5	17.00	793512	20.0
C(14a)-Homo-27-no...	19.36	12.7	ng	287479	6	18.64	451698	20.0
Hop-22(29)-en-3.b...	20.88	15.2	ng	342780	6	18.64	451698	20.0
5-Bromo-4-oxo-4,5...	21.86	14.4	ng	326059	6	18.64	451698	20.0