

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121315\
 Data File : BF083766.D
 Acq On : 14 Dec 2015 8:49
 Operator : UM/IZ
 Sample : G4648-13
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-07(0-0.16)

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-BF121315.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.116	262	265	269	rBV	457358	617442	21.47%	1.928%
2	5.527	298	301	304	rVB	2368052	2548131	88.59%	7.957%
3	6.544	387	390	393	rBV	2070761	2389404	83.07%	7.462%
4	6.693	400	403	405	rBV	2223333	2876307	100.00%	8.982%
5	6.921	420	423	425	rBV	678505	592054	20.58%	1.849%
6	7.081	434	437	439	rVB	2189845	2364747	82.21%	7.385%
7	7.482	469	472	474	rBV	1729807	1864683	64.83%	5.823%
8	8.202	533	535	537	rBV	841131	720020	25.03%	2.248%
9	9.276	627	629	632	rBV	2695027	2498286	86.86%	7.802%
10	9.665	661	663	666	rVB	408596	403956	14.04%	1.261%
11	9.893	681	683	686	rBV	47824	38140	1.33%	0.119%
12	9.950	686	688	691	rBV	597360	642122	22.32%	2.005%
13	10.396	723	727	728	rBV	53539	66651	2.32%	0.208%
14	10.613	743	746	749	rBV	19746	30259	1.05%	0.094%
15	10.739	754	757	759	rBV	1645472	1544834	53.71%	4.824%
16	10.865	766	768	773	rBV	67915	77849	2.71%	0.243%
17	11.082	782	787	792	rBV5	9362	28793	1.00%	0.090%
18	11.436	816	818	819	rVV	649882	650698	22.62%	2.032%
19	11.459	819	820	822	rVV	62940	61827	2.15%	0.193%
20	11.493	822	823	826	rVB2	42264	56808	1.98%	0.177%
21	11.962	863	864	867	rBV2	141582	171644	5.97%	0.536%
22	12.042	869	871	875	rVB2	21169	38089	1.32%	0.119%
23	12.145	876	880	885	rVB3	25514	57748	2.01%	0.180%
24	12.248	885	889	892	rBV3	14727	32525	1.13%	0.102%
25	12.579	916	918	921	rVV	152703	155486	5.41%	0.486%
26	12.636	921	923	925	rVB	138446	154519	5.37%	0.483%
27	12.762	930	934	937	rBV4	16385	36241	1.26%	0.113%
28	12.842	939	941	945	rBV2	215924	342410	11.90%	1.069%
29	12.922	945	948	950	rVB2	67120	76277	2.65%	0.238%
30	13.014	954	956	959	rVB	2180142	2246641	78.11%	7.016%
31	13.196	969	972	973	rBV	26276	31895	1.11%	0.100%
32	13.254	976	977	979	rVB2	31641	37275	1.30%	0.116%
33	13.494	994	998	1000	rBV2	97861	142136	4.94%	0.444%
34	13.551	1000	1003	1005	rVV	166296	173247	6.02%	0.541%

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35	13.596	1005	1007	1010	rVV2	57928	82948	2.88%	0.259%
36	13.711	1014	1017	1021	rVV2	22959	41676	1.45%	0.130%
37	13.916	1032	1035	1039	rBV2	273072	396920	13.80%	1.239%
38	14.065	1045	1048	1049	rBV2	689613	907282	31.54%	2.833%
39	14.236	1061	1063	1065	rBV2	106195	121287	4.22%	0.379%
40	14.271	1065	1066	1069	rVV2	37869	48888	1.70%	0.153%
41	14.431	1078	1080	1083	rBV2	24186	36214	1.26%	0.113%
42	14.545	1087	1090	1092	rBV	254068	276677	9.62%	0.864%
43	14.842	1113	1116	1120	rBV	94447	135976	4.73%	0.425%
44	14.922	1120	1123	1126	rBV	141292	160069	5.57%	0.500%
45	14.991	1126	1129	1131	rVB2	41836	63152	2.20%	0.197%
46	15.082	1135	1137	1143	rBV	90955	225740	7.85%	0.705%
47	15.174	1143	1145	1149	rBV	208235	354179	12.31%	1.106%
48	15.288	1153	1155	1157	rVV2	61910	77005	2.68%	0.240%
49	15.379	1161	1163	1166	rVB	51940	69052	2.40%	0.216%
50	15.437	1166	1168	1171	rBV	51345	78571	2.73%	0.245%
51	15.505	1171	1174	1176	rBV	443905	558900	19.43%	1.745%
52	15.551	1176	1178	1180	rVB	57049	79973	2.78%	0.250%
53	15.745	1193	1195	1200	rBV	101041	161550	5.62%	0.504%
54	15.917	1208	1210	1213	rBV	81463	126729	4.41%	0.396%
55	15.985	1213	1216	1218	rVV	393480	585585	20.36%	1.829%
56	16.031	1218	1220	1225	rVV3	66730	136601	4.75%	0.427%
57	16.134	1227	1229	1234	rVB2	50066	96351	3.35%	0.301%
58	16.237	1234	1238	1245	rVB3	62679	215792	7.50%	0.674%
59	16.351	1245	1248	1255	rVB2	36556	105926	3.68%	0.331%
60	16.477	1257	1259	1261	rBV	26444	44755	1.56%	0.140%
61	16.762	1281	1284	1290	rBV	156613	322691	11.22%	1.008%
62	16.934	1296	1299	1301	rBV2	25544	47590	1.65%	0.149%
63	16.991	1301	1304	1307	rVV2	37947	75489	2.62%	0.236%
64	17.071	1307	1311	1314	rVV	69926	159048	5.53%	0.497%
65	17.140	1314	1317	1321	rVB4	41319	100117	3.48%	0.313%
66	17.528	1348	1351	1357	rBV3	73196	189190	6.58%	0.591%
67	17.757	1367	1371	1376	rVB	48291	122867	4.27%	0.384%
68	17.848	1376	1379	1382	rBV5	22946	60532	2.10%	0.189%
69	17.928	1382	1386	1388	rVV3	46915	131522	4.57%	0.411%
70	17.985	1388	1391	1398	rVB	118008	328917	11.44%	1.027%
71	18.168	1404	1407	1416	rVB5	42096	158689	5.52%	0.496%

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72	18.443	1427	1431	1436	rBV2	88750	257080	8.94%	0.803%
73	18.603	1436	1445	1454	rVB	327977	1112233	38.67%	3.473%

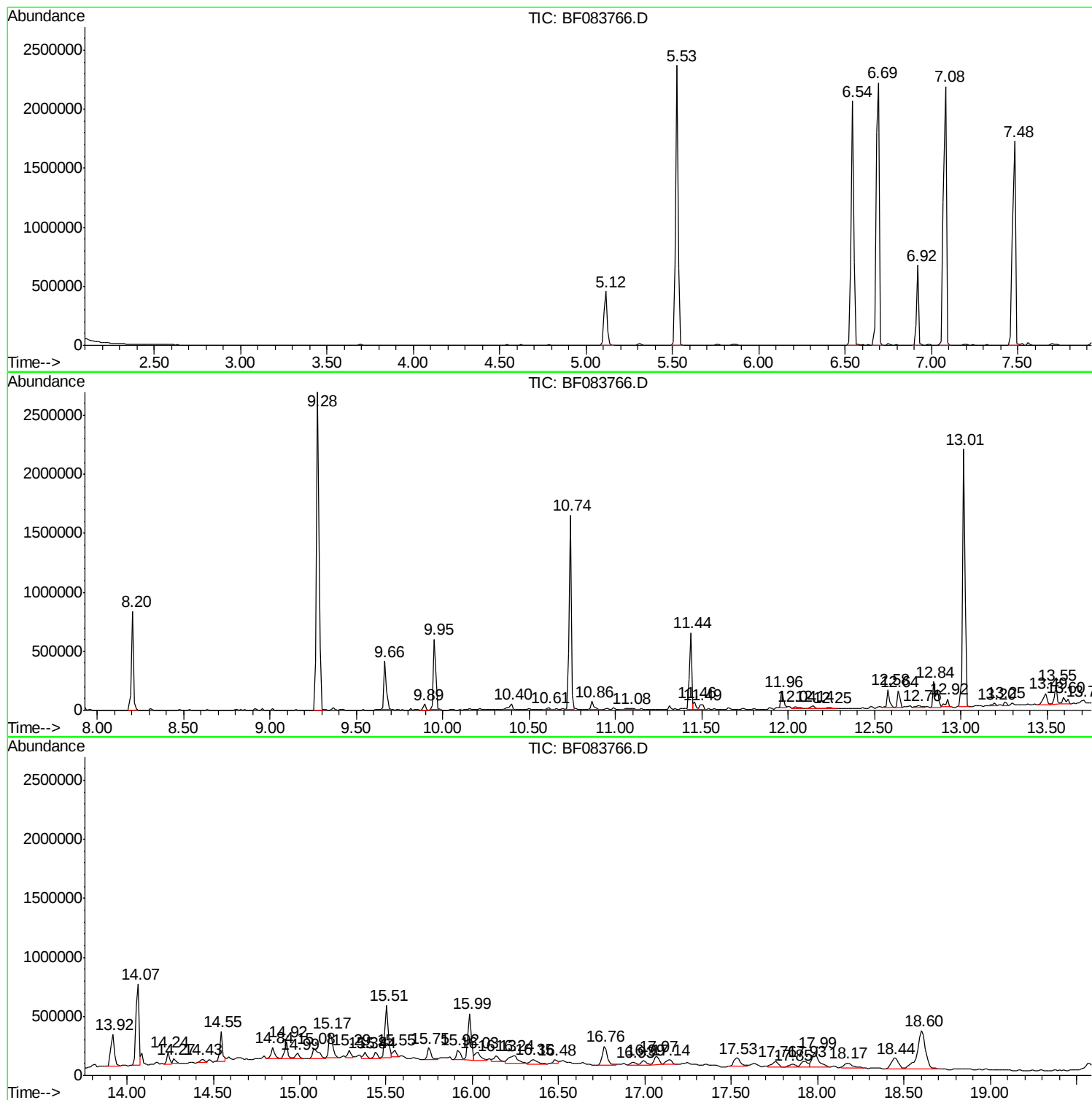
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Instrument :
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SB-07(0-0.16)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121315.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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Instrument :
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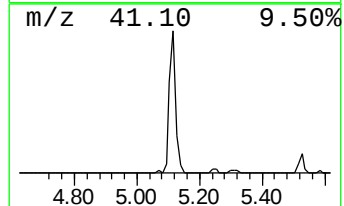
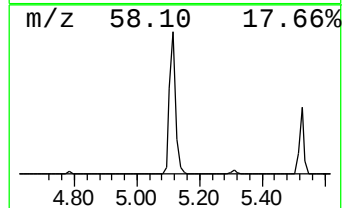
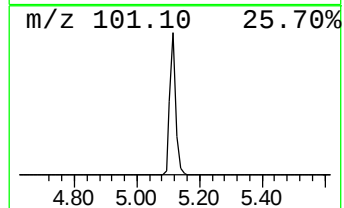
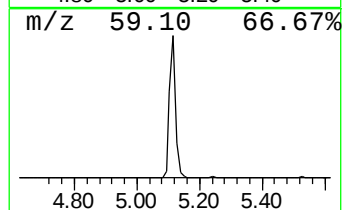
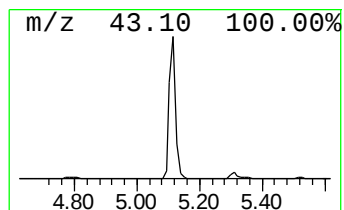
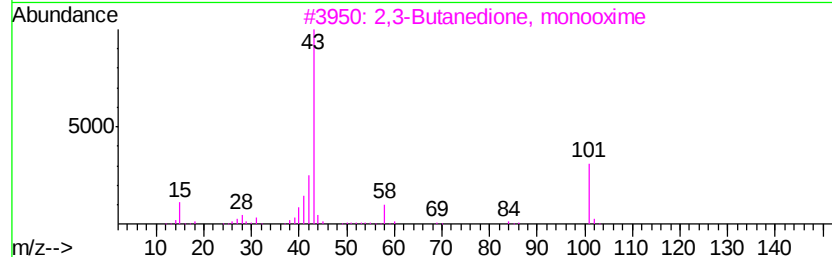
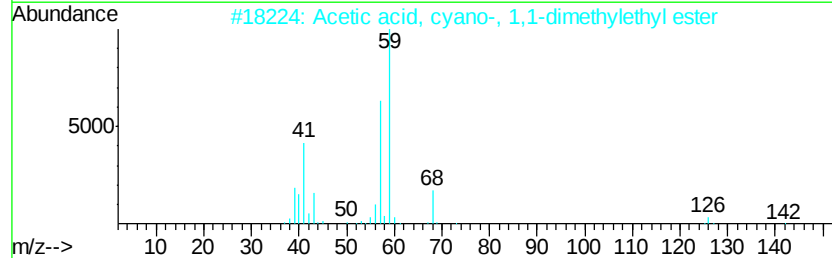
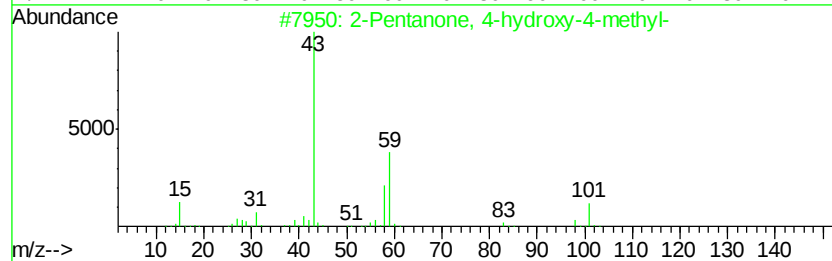
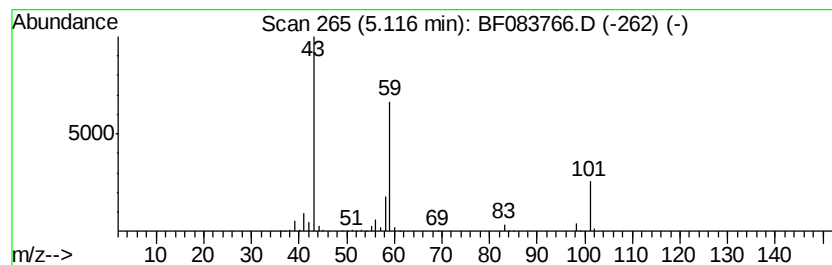
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121315.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	20.86 ng	617442	1,4-Dichlorobenzene-d4	6.92

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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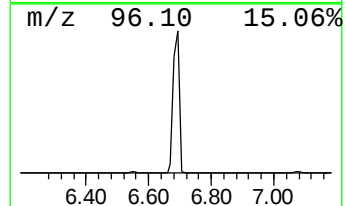
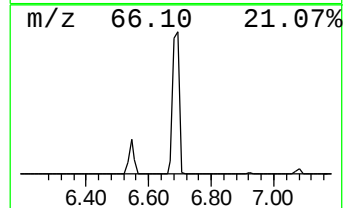
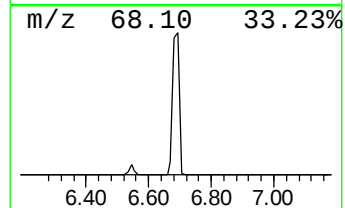
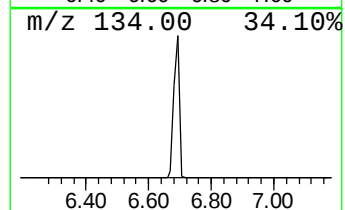
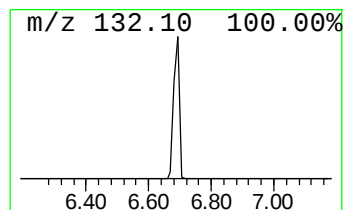
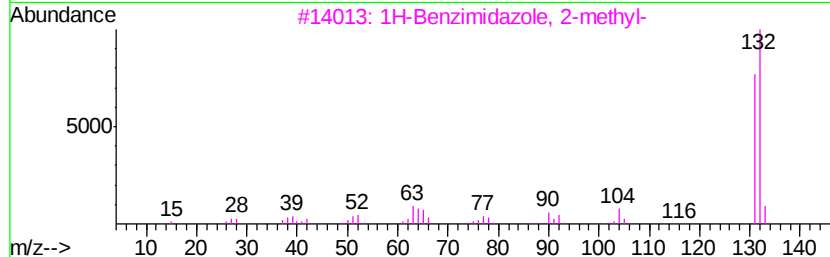
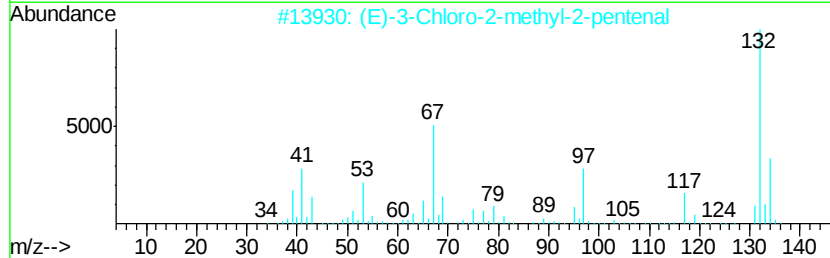
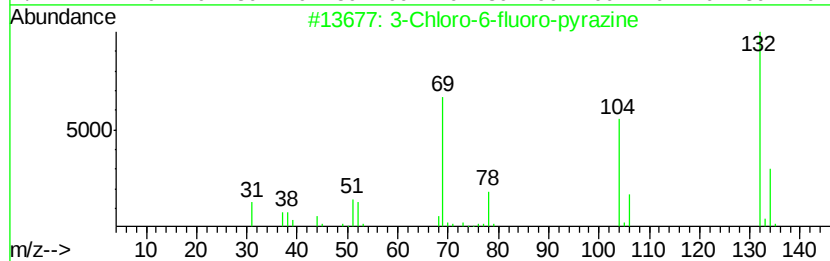
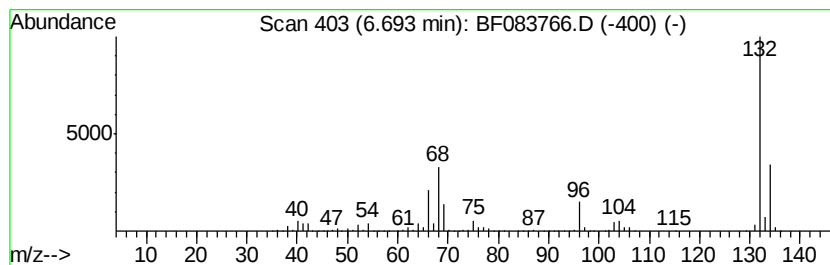
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.69	97.16 ng	2876310	1,4-Dichlorobenzene-d4	6.92

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	38
2	(E)-3-Chloro-2-methyl-2-pentenal			132	C6H9ClO	031357-76-3	25
3	1H-Benzimidazole, 2-methyl-			132	C8H8N2	000615-15-6	12
4	2H-Cyclopenta[d]pyridazine, 2-me...			132	C8H8N2	022291-85-6	9
5	4-Chloro-6-fluoro-pyrimidine			132	C4H2ClFN2	051422-01-6	9



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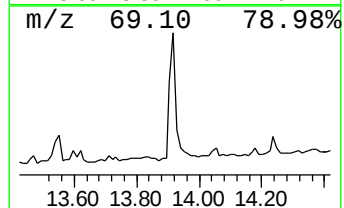
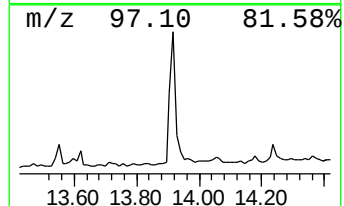
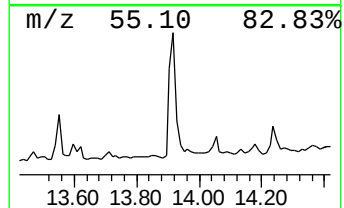
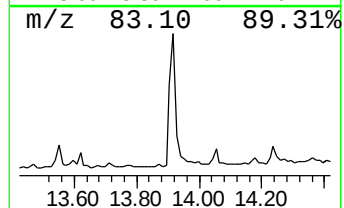
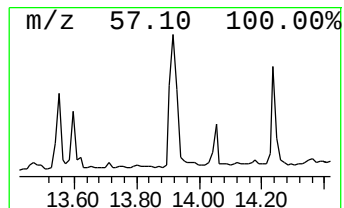
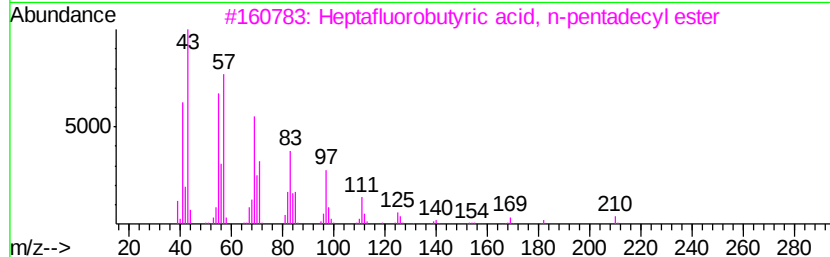
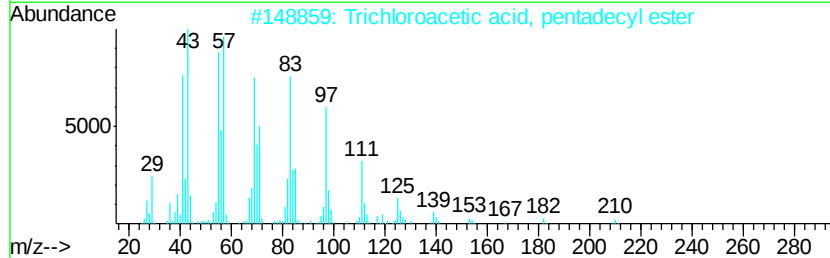
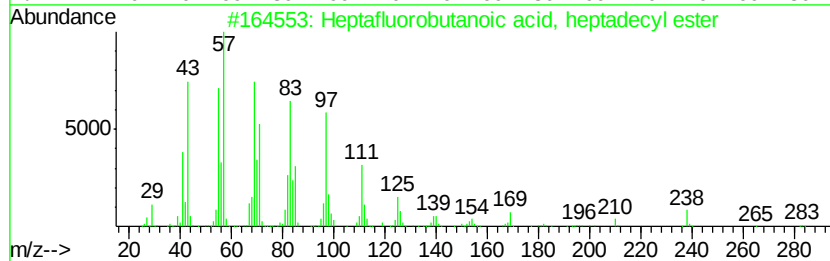
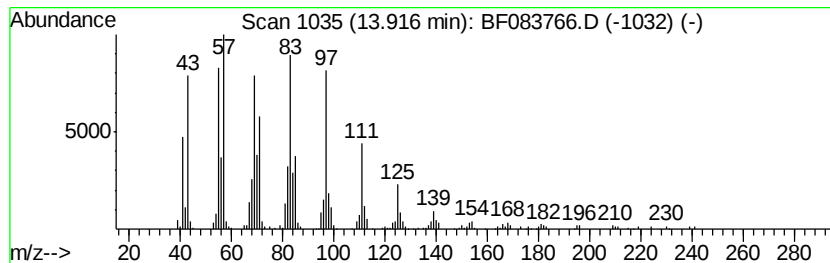
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Heptafluorobutanoic acid, h... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.92	8.75 ng	396920	Chrysene-d12	14.07

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94
2		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	91
3		Heptafluorobutvric acid, n-penta...	424	C19H31F7O2	1000216-79-3	91
4		1-Octadecanol	270	C18H38O	000112-92-5	90
5		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	90



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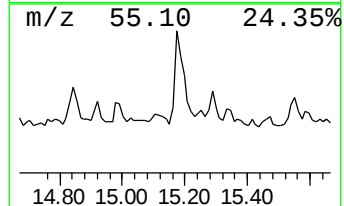
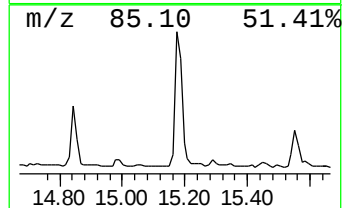
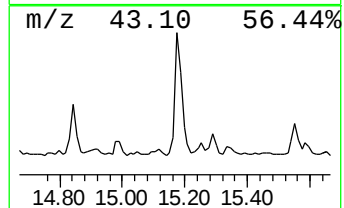
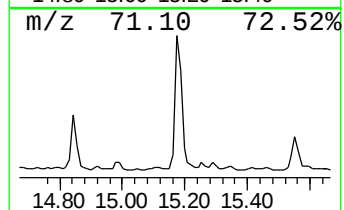
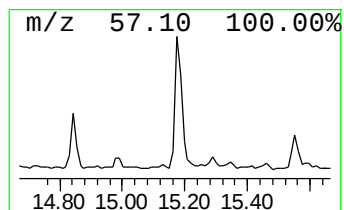
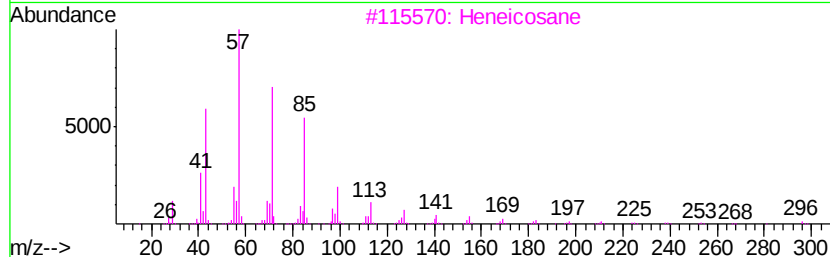
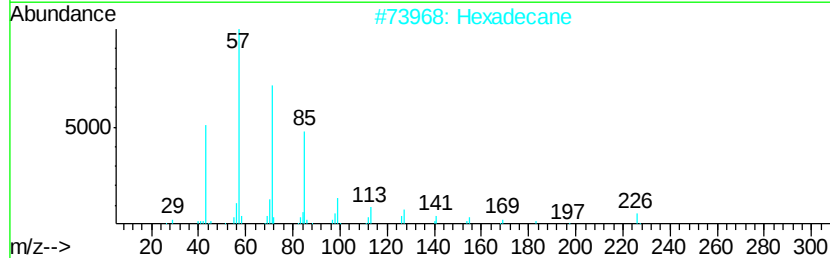
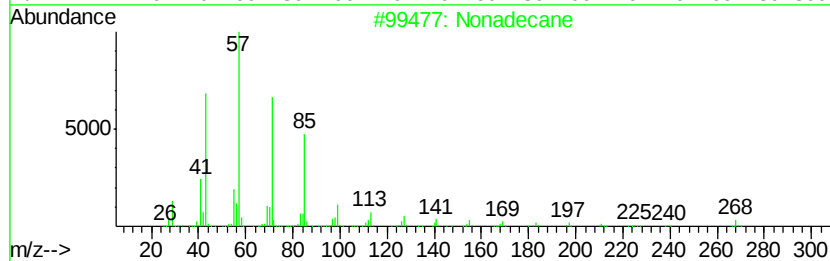
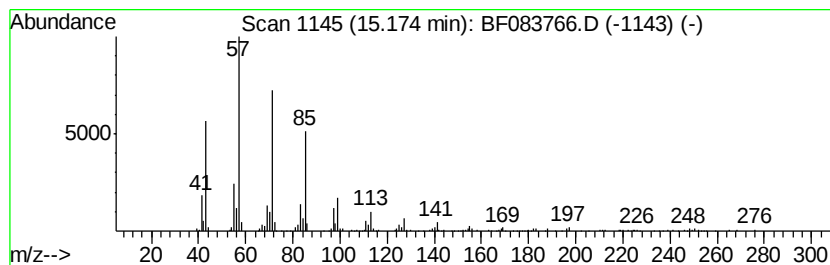
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Nonadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.17	12.67 ng	354179	Perylene-d12	15.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	93
2	Hexadecane		226	C16H34	000544-76-3	91
3	Heneicosane		296	C21H44	000629-94-7	87
4	Pentadecane, 8-heptyl-		310	C22H46	071005-15-7	87
5	Tetratriacontane		479	C34H70	014167-59-0	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121315\
Data File : BF083766.D
Acq On : 14 Dec 2015 8:49
Operator : UM/IZ
Sample : G4648-13
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-07(0-0.16)

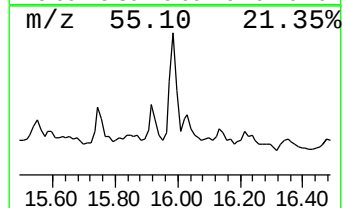
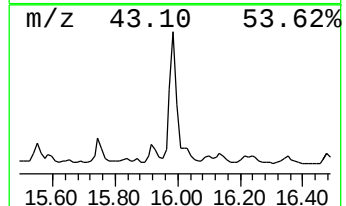
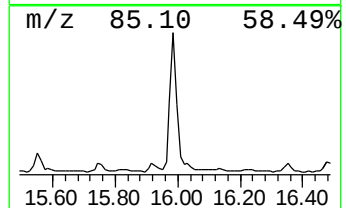
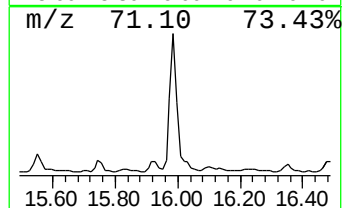
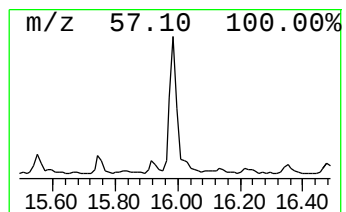
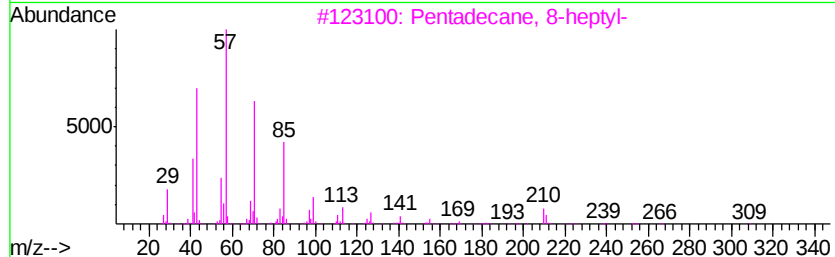
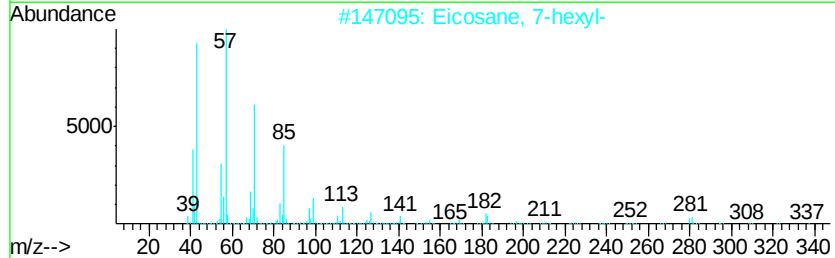
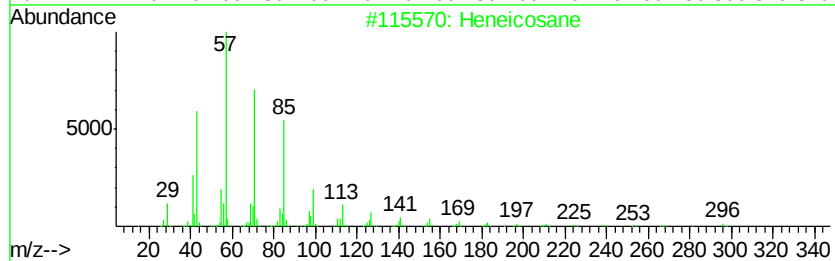
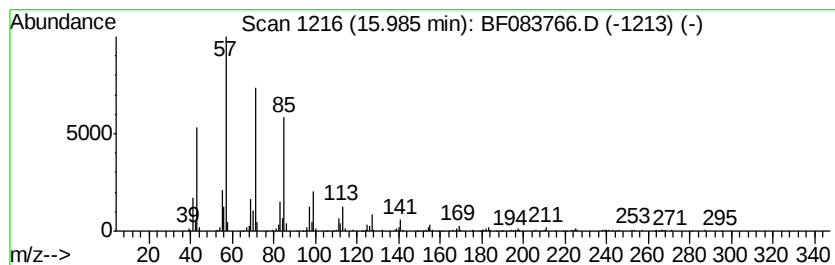
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Heneicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.99	20.95 ng	585585	Perylene-d12	15.51

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	91
3	Pentadecane, 8-heptyl-		310	C22H46	071005-15-7	91
4	Nonadecane		268	C19H40	000629-92-5	90
5	Tetracosane		338	C24H50	000646-31-1	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121315\
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Sample : G4648-13
Misc :
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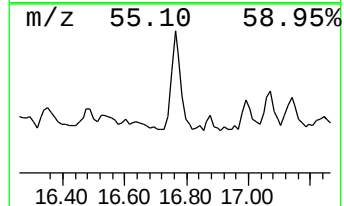
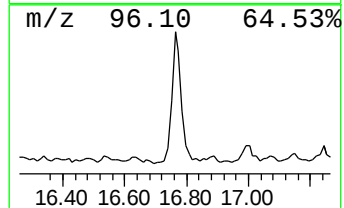
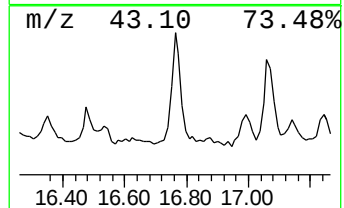
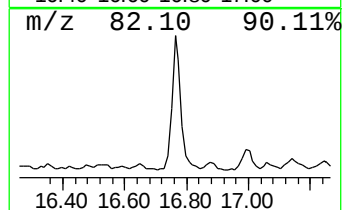
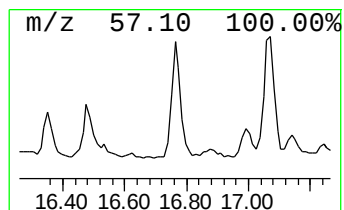
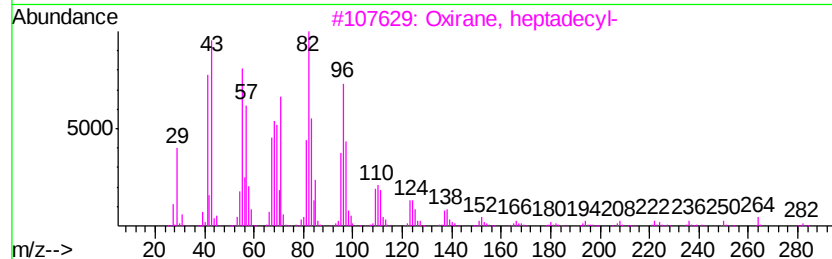
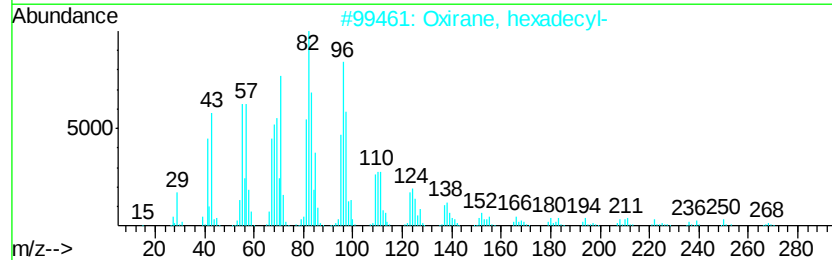
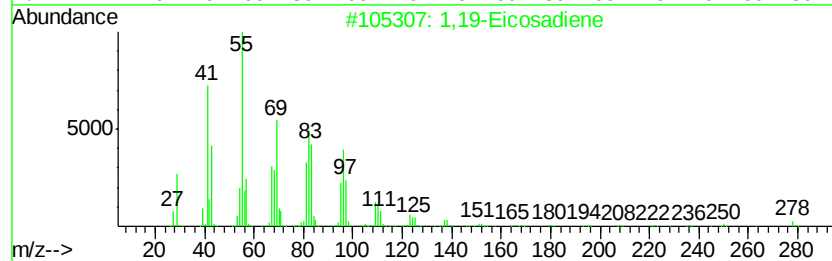
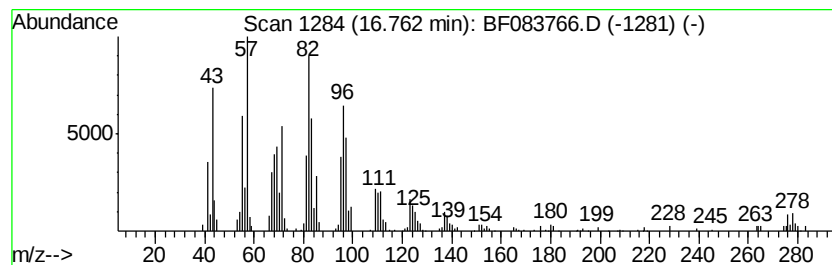
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1,19-Eicosadiene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.76	11.55 ng	322691	Perylene-d12	15.51	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,19-Eicosadiene	278	C20H38	014811-95-1	96
2	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	94
3	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	91
4	Octadecanal	268	C18H36O	000638-66-4	91
5	Bicyclo[10.8.0]eicosane, cis-	278	C20H38	1000155-82-2	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121315\
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Operator : UM/IZ
Sample : G4648-13
Misc :
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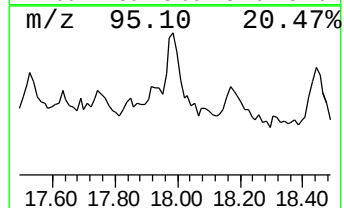
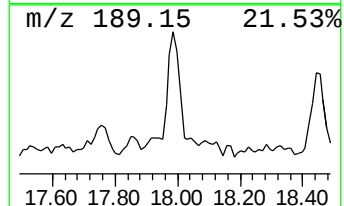
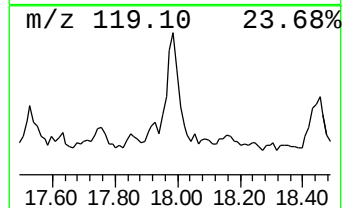
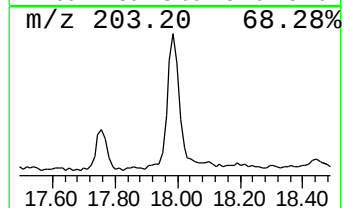
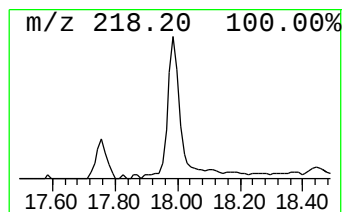
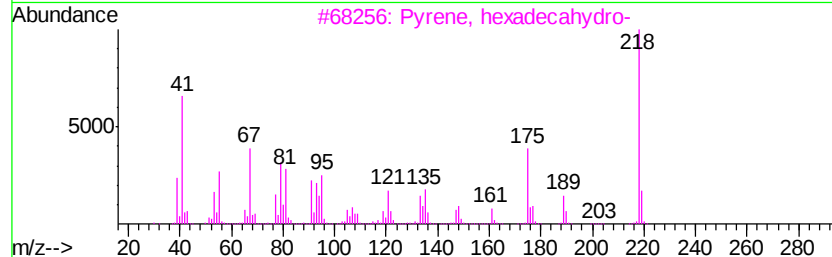
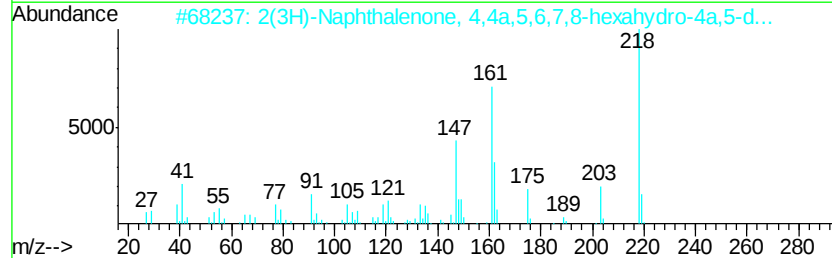
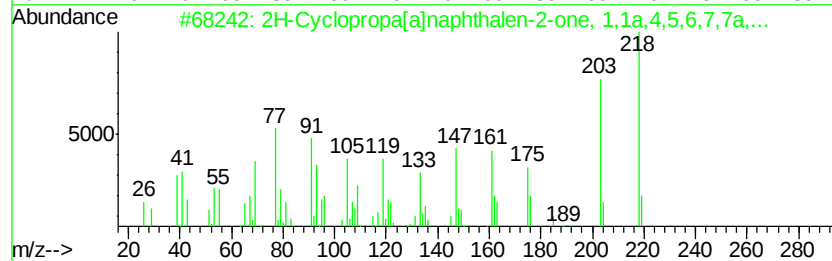
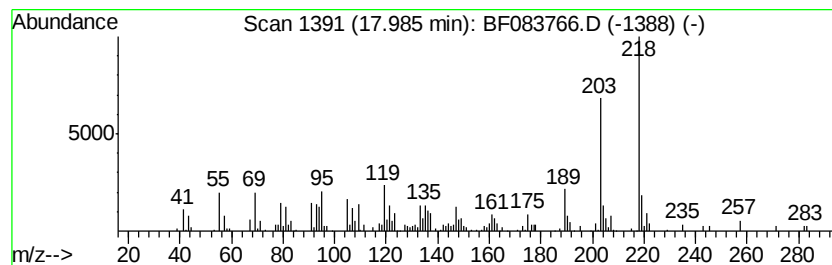
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 2H-Cyclopropa[a]naphthalen-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
17.99	11.77 ng	328917	Perylene-d12	15.51		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2H-Cyclopropa[alnaphthalen-2-one...	218	C15H22O	006831-17-0	50
2		2(3H)-Naphthalenone, 4,4a,5,6,7,...	218	C15H22O	019598-45-9	45
3		Pvrene, hexadecahydro-	218	C16H26	002435-85-0	41
4		Benzo[blnaphtho[1,2-d]furan	218	C16H10O	000239-30-5	38
5		3-Methoxy-6,7,8,9-tetrahydro-dib...	218	C13H14O3	1000186-57-9	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121315\
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Sample : G4648-13
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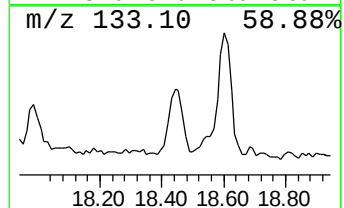
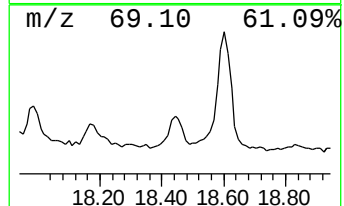
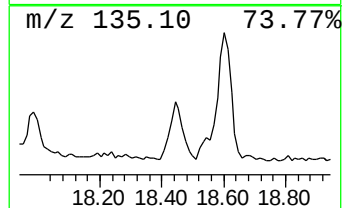
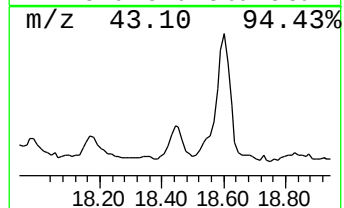
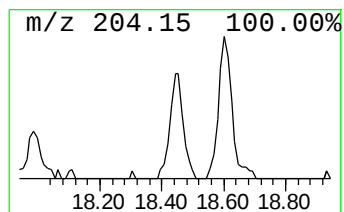
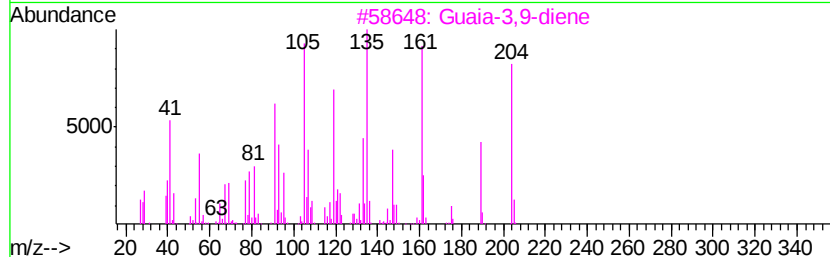
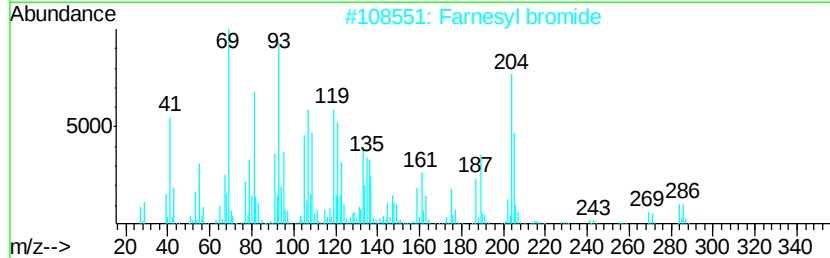
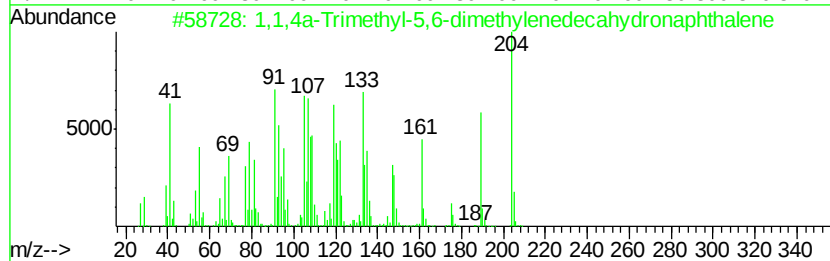
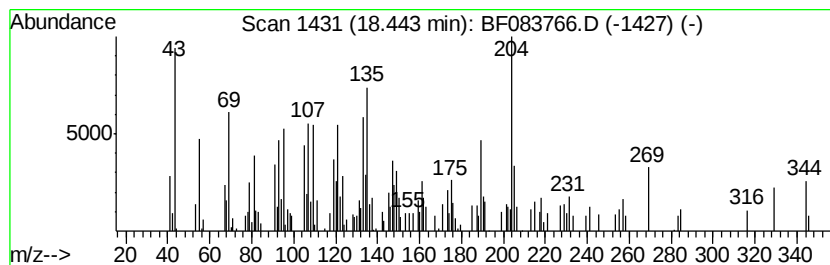
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1,1,4a-Trimethyl-5,6-dimeth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.44	9.20 ng	257080	Perylene-d12	15.51		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1,4a-Trimethyl-5,6-dimethylene...	204	C15H24	1000193-60-8	53	
2	Farnesyl bromide	284	C15H25Br	006874-67-5	53	
3	Guaia-3,9-diene	204	C15H24	000489-83-8	45	
4	1H-Cycloprop[elazulene, 1a,2,3,5...	204	C15H24	021747-46-6	45	
5	Naphthalene, decahydro-4a-methyl...	204	C15H24	017066-67-0	43	



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121315\
Data File : BF083766.D
Acq On : 14 Dec 2015 8:49
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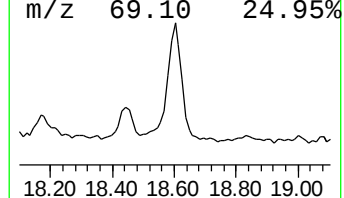
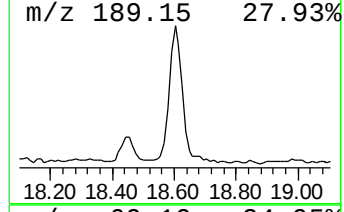
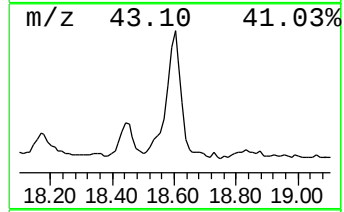
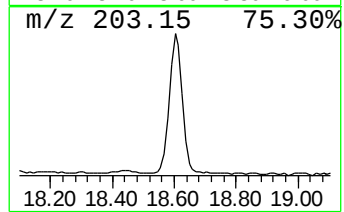
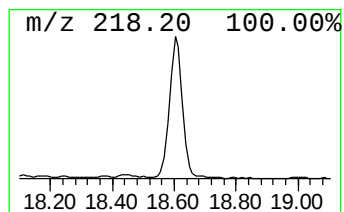
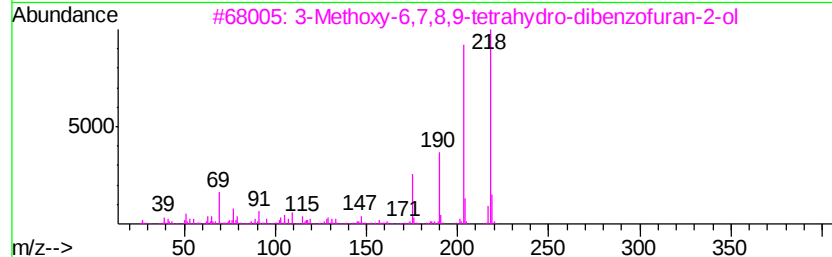
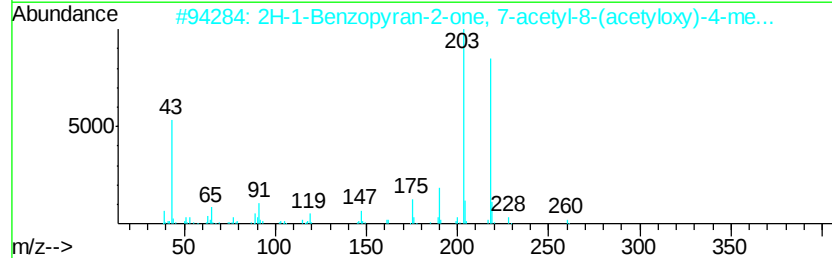
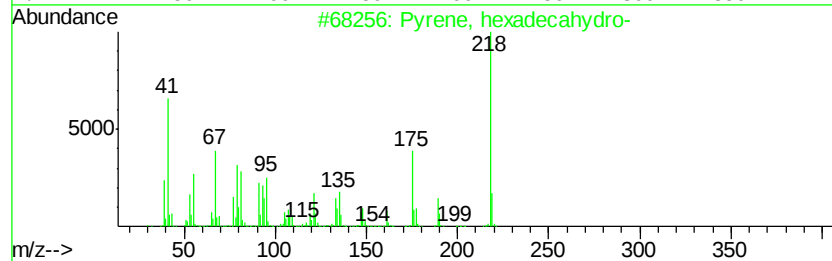
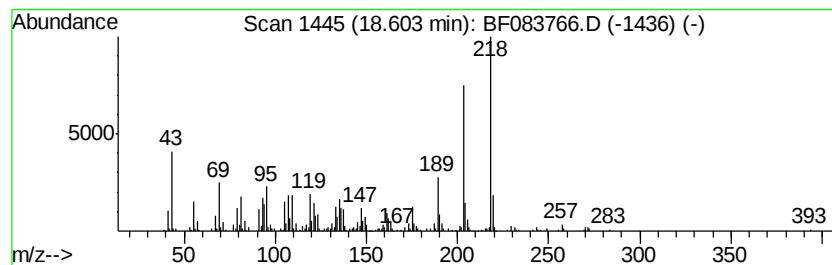
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pyrene, hexadecahydro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.60	39.80 ng	1112230	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, hexadecahydro-	218	C16H26	002435-85-0	62
2		2H-1-Benzopyran-2-one, 7-acetyl-...	260	C14H12O5	129812-32-4	53
3		3-Methoxy-6,7,8,9-tetrahydro-dib...	218	C13H14O3	1000186-57-9	52
4		2,3-Diethoxyquinoxaline	218	C12H14N2O2	057050-66-5	50
5		Benzo[b]naphtho[2,3-d]furan	218	C16H10O	000243-42-5	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121315\
Data File : BF083766.D
Acq On : 14 Dec 2015 8:49
Operator : UM/IZ
Sample : G4648-13
Misc :
ALS Vial : 56 Sample Multiplier: 1

Instrument :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.12	20.9	ng	617442	1	6.92	592054	20.0
unknown6.69	6.69	97.2	ng	2876310	1	6.92	592054	20.0
Heptafluorobutano...	13.92	8.8	ng	396920	5	14.07	907282	20.0
Nonadecane	15.17	12.7	ng	354179	6	15.51	558900	20.0
Heneicosane	15.99	20.9	ng	585585	6	15.51	558900	20.0
1,19-Eicosadiene	16.76	11.6	ng	322691	6	15.51	558900	20.0
2H-Cyclopropa[a]n...	17.99	11.8	ng	328917	6	15.51	558900	20.0
1,1,4a-Trimethyl-...	18.44	9.2	ng	257080	6	15.51	558900	20.0
Pyrene, hexadecah...	18.60	39.8	ng	1112230	6	15.51	558900	20.0