

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110615\
 Data File : BF082733.D
 Acq On : 5 Nov 2015 17:04
 Operator : UM/IZ
 Sample : G4261-12
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DW4

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-BF103015.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	2.133	3	4	18	rVB2	106383	186769	3.33%	0.366%
2	5.059	256	260	269	rVB	1792730	2809577	50.15%	5.500%
3	5.470	293	296	299	rBV	3865643	5190992	92.65%	10.162%
4	6.499	383	386	389	rBV	4347748	4983386	88.95%	9.755%
5	6.636	395	398	400	rBV	5736939	5554595	99.14%	10.873%
6	6.865	416	418	420	rBV	677547	545232	9.73%	1.067%
7	7.025	429	432	434	rBV	4902021	4230739	75.51%	8.282%
8	7.425	464	467	470	rBV	2837491	3429874	61.22%	6.714%
9	8.145	528	530	533	rBV	480364	604174	10.78%	1.183%
10	9.231	622	625	627	rBV	6151799	5602769	100.00%	10.968%
11	9.619	657	659	662	rBV	1299489	1155583	20.63%	2.262%
12	9.905	681	684	686	rBV	861099	731933	13.06%	1.433%
13	10.328	718	721	722	rBV2	50224	68153	1.22%	0.133%
14	10.694	750	753	755	rBV	4041721	3922219	70.01%	7.678%
15	10.819	762	764	768	rVB	44521	60272	1.08%	0.118%
16	11.265	802	803	808	rVB2	42123	81490	1.45%	0.160%
17	11.379	811	813	817	rVV	479954	1028677	18.36%	2.014%
18	11.459	817	820	821	rVB2	82799	80034	1.43%	0.157%
19	11.928	859	861	863	rBV2	217105	279686	4.99%	0.547%
20	11.996	865	867	872	rBV2	121982	164315	2.93%	0.322%
21	12.202	882	885	887	rBV3	49464	93617	1.67%	0.183%
22	12.591	917	919	922	rBV	705392	887571	15.84%	1.737%
23	12.705	927	929	931	rBV2	83645	113385	2.02%	0.222%
24	12.819	936	939	942	rBV2	599297	735139	13.12%	1.439%
25	12.979	950	953	955	rVB	4286308	4594019	82.00%	8.993%
26	13.151	965	968	970	rVB	157231	177051	3.16%	0.347%
27	13.219	972	974	975	rBV	190897	158312	2.83%	0.310%
28	13.768	1020	1022	1023	rBV	105016	117630	2.10%	0.230%
29	13.871	1029	1031	1036	rBV2	113171	199807	3.57%	0.391%
30	14.020	1041	1044	1045	rBV2	1001581	1166850	20.83%	2.284%
31	14.511	1085	1087	1088	rBV2	58231	88112	1.57%	0.172%
32	15.048	1131	1134	1139	rVV	253349	564310	10.07%	1.105%
33	15.140	1140	1142	1145	rVB2	75033	94157	1.68%	0.184%
34	15.334	1157	1159	1162	rVB	180199	214918	3.84%	0.421%

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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-BF103015.M
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35	15.391	1162	1164	1167	rVB	190606	233522	4.17%	0.457%
36	15.448	1167	1169	1171	rBV	300910	443813	7.92%	0.869%
37	16.843	1289	1291	1296	rVB2	118044	267217	4.77%	0.523%
38	17.266	1325	1328	1334	rVB	86966	224334	4.00%	0.439%

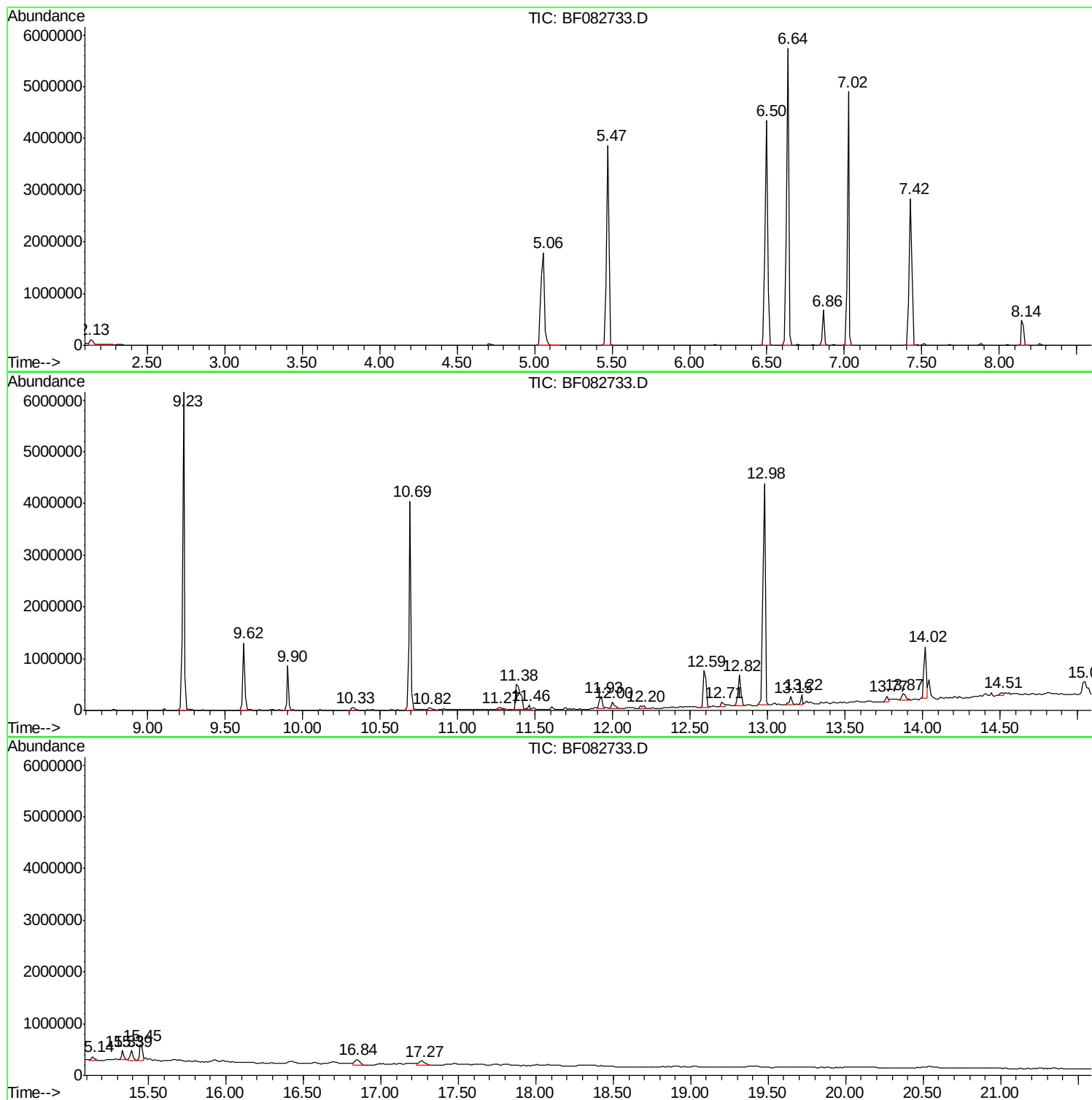
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF103015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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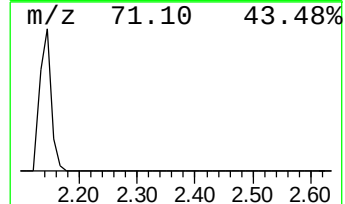
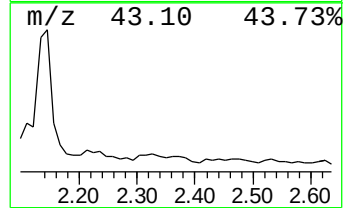
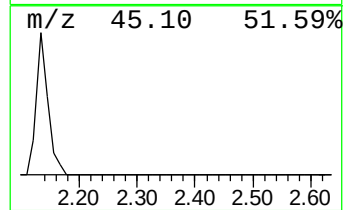
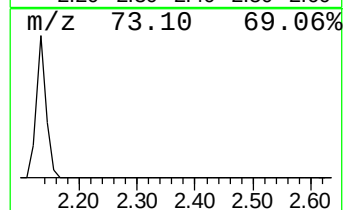
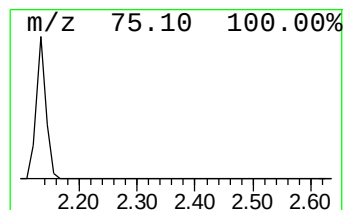
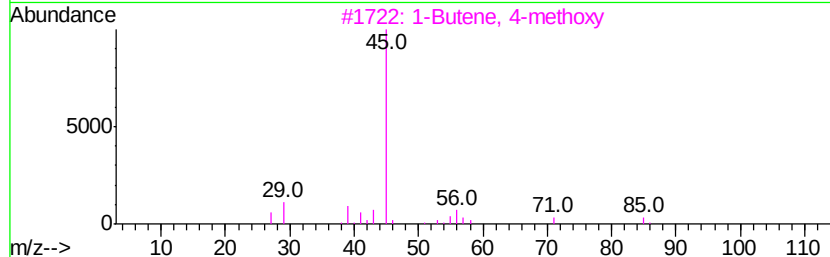
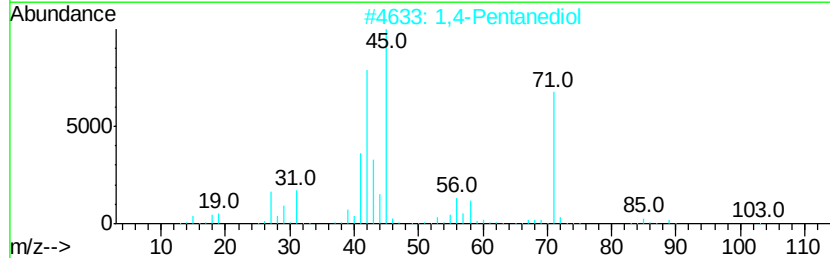
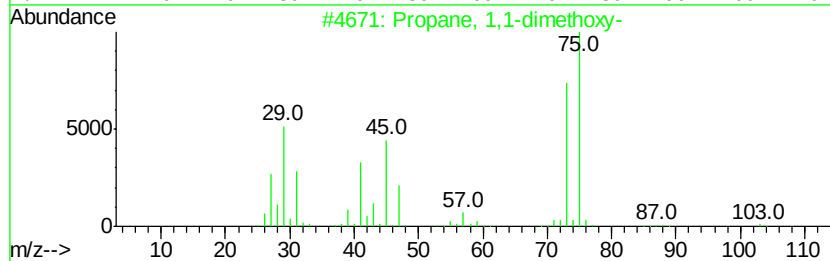
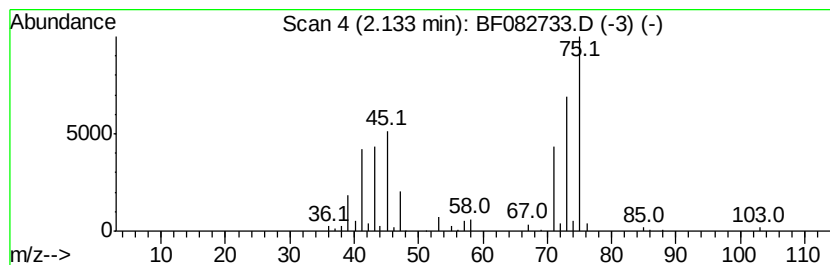
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Propane, 1,1-dimethoxy- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.13	6.85 ng	186769	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	59
2		1,4-Pentanediol	104	C5H12O2	000626-95-9	12
3		1-Butene, 4-methoxy	86	C5H10O	004696-30-4	9
4		Azetidine, 1-methyl-	71	C4H9N	004923-79-9	9
5		Propanal, 3-methoxy-	88	C4H8O2	002806-84-0	9



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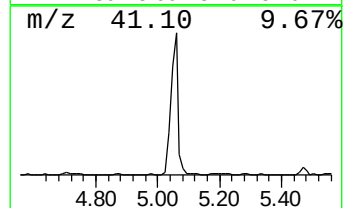
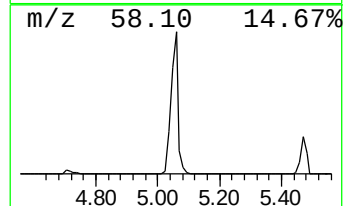
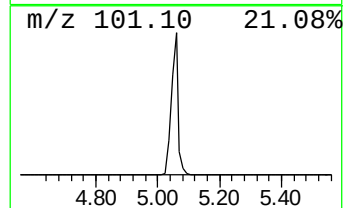
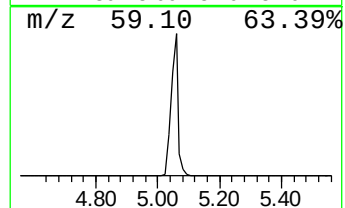
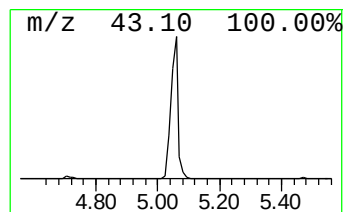
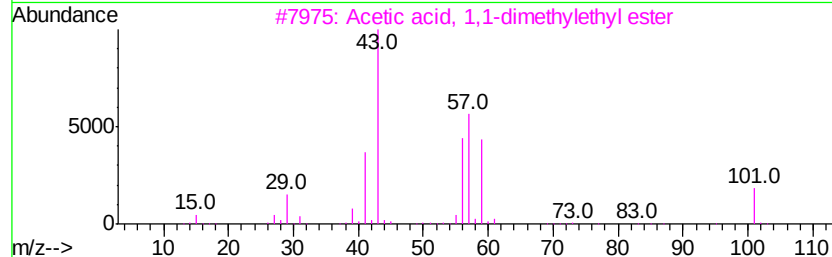
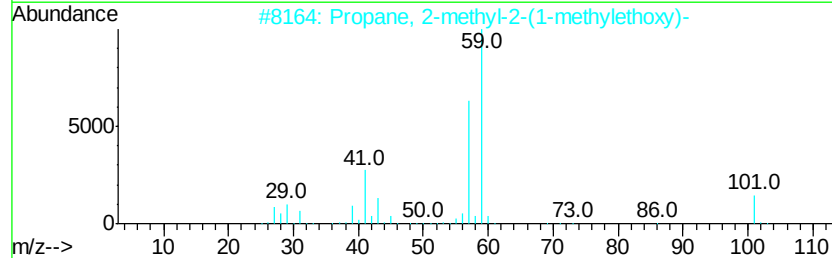
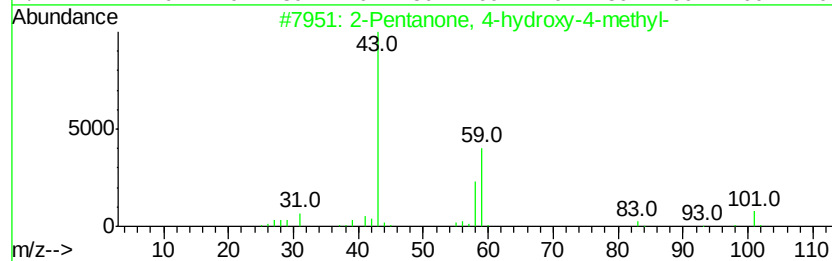
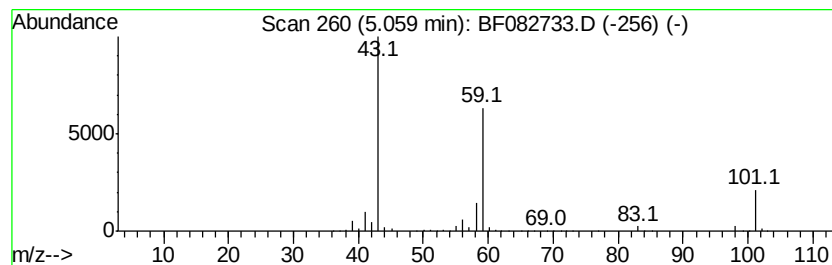
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.06	103.06 ng	2809580	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	53
3		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
5		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17



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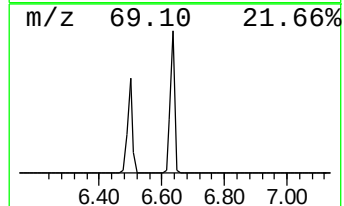
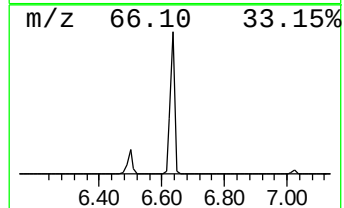
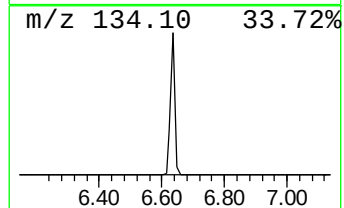
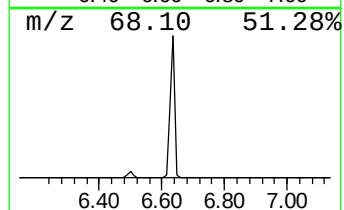
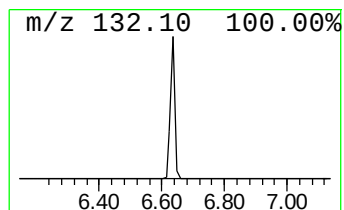
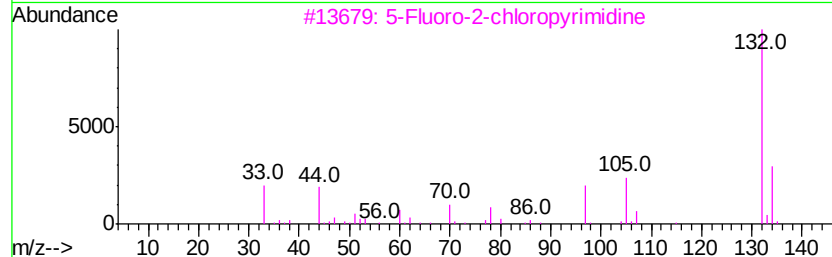
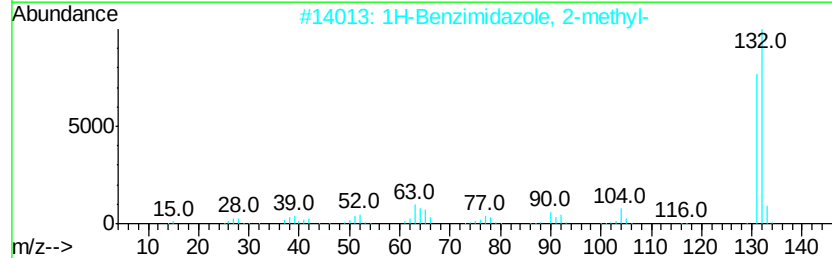
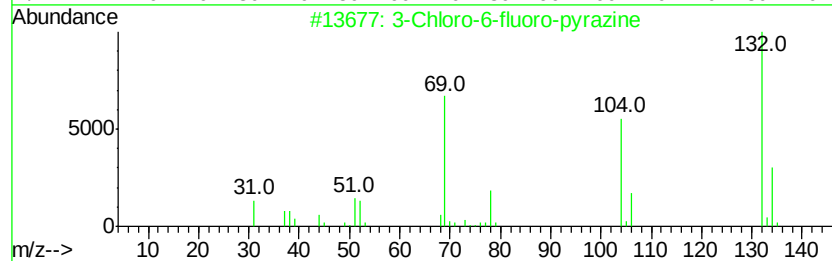
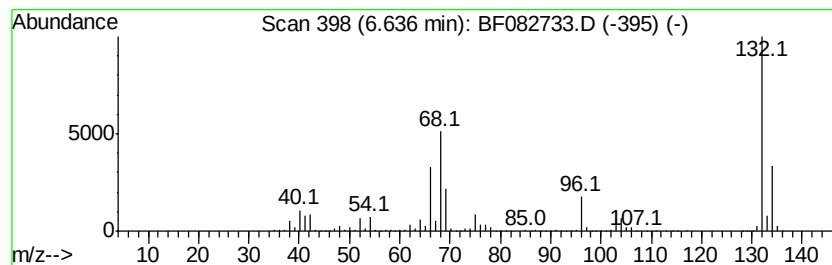
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown6.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	203.75 ng	5554600	1,4-Dichlorobenzene-d4	6.86

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	10



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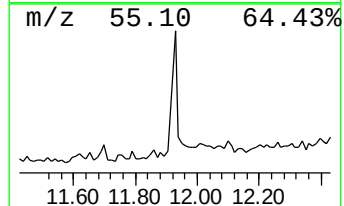
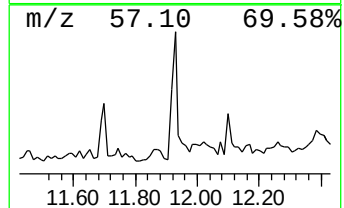
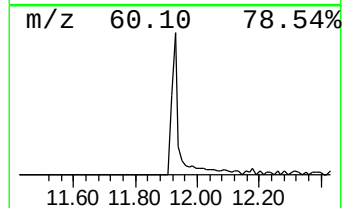
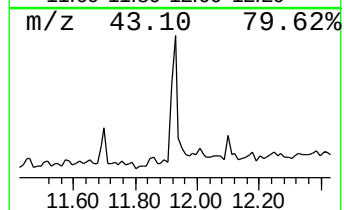
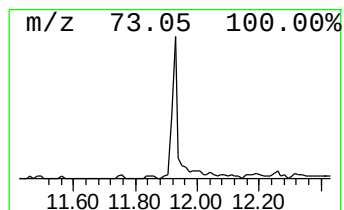
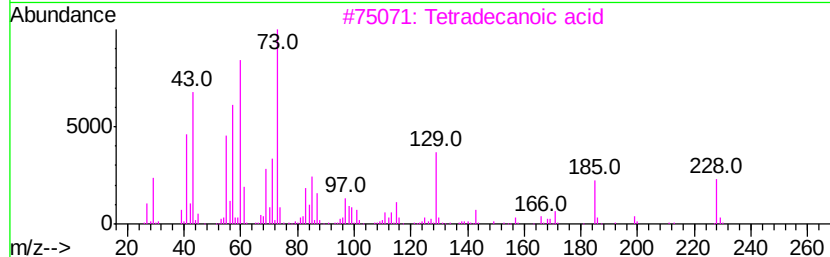
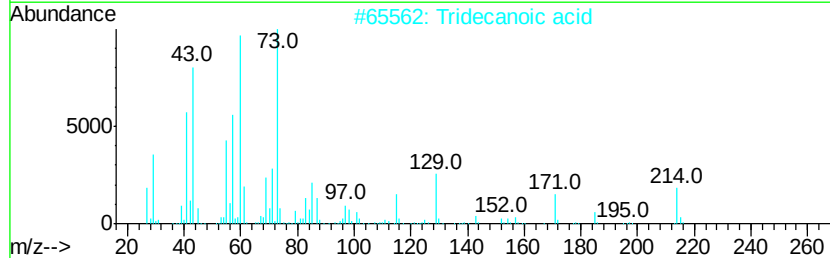
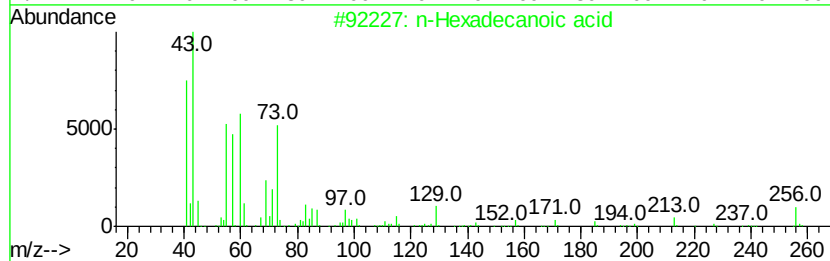
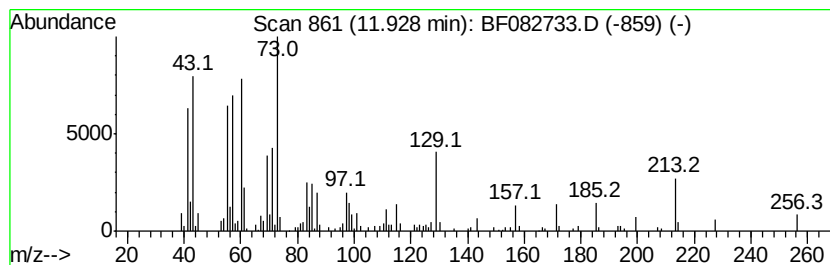
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	5.44 ng	279686	Phenanthrene-d10	11.39

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
2		Tridecanoic acid	214	C13H26O2	000638-53-9	89
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
4		n-Decanoic acid	172	C10H20O2	000334-48-5	76
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	72



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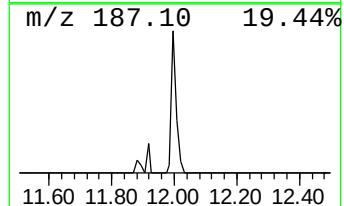
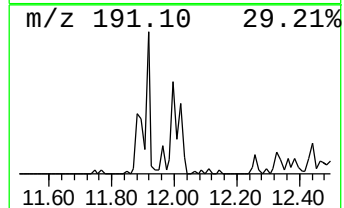
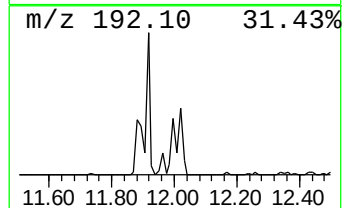
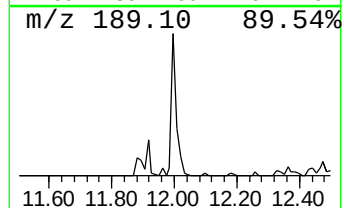
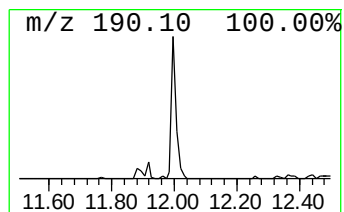
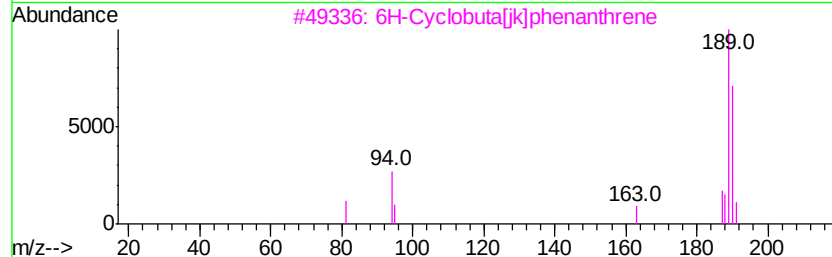
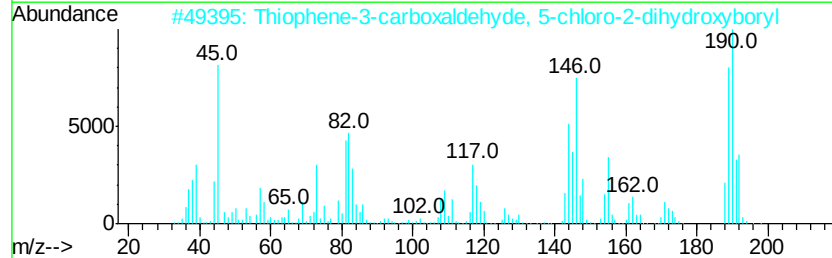
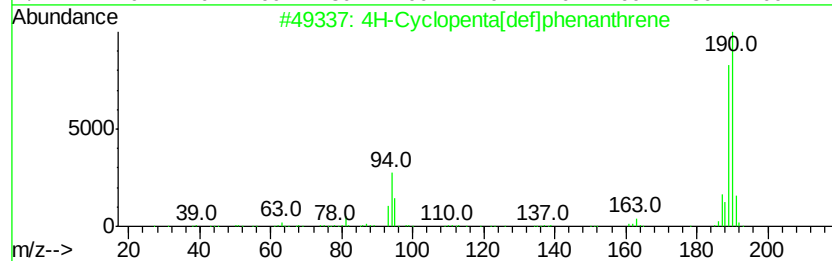
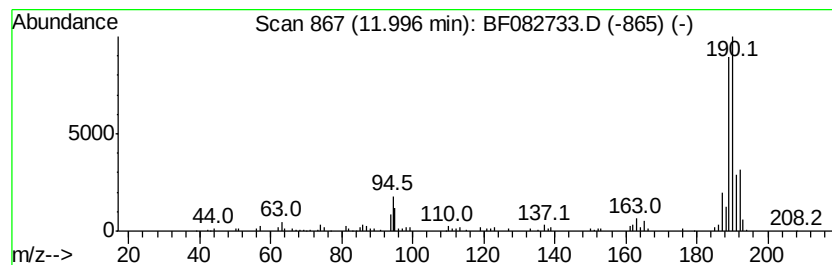
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	3.19 ng	164315	Phenanthrene-d10	11.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	53
3		6H-Cyclobuta[flk]phenanthrene	190	C15H10	083469-43-6	33
4		Pyrimidine, 2-chloro-5-phenyl-	190	C10H7ClN2	022536-62-5	25
5		1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	25



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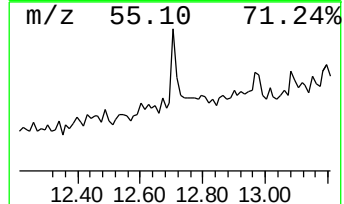
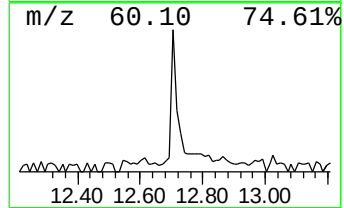
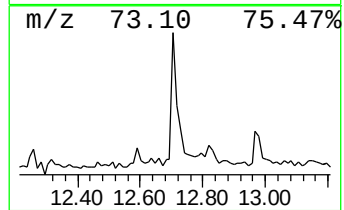
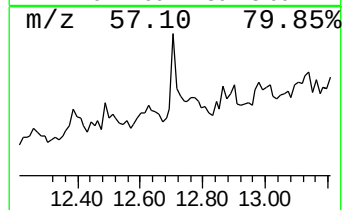
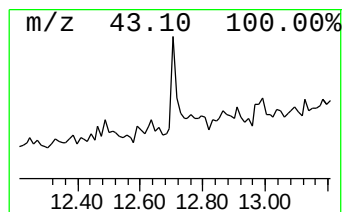
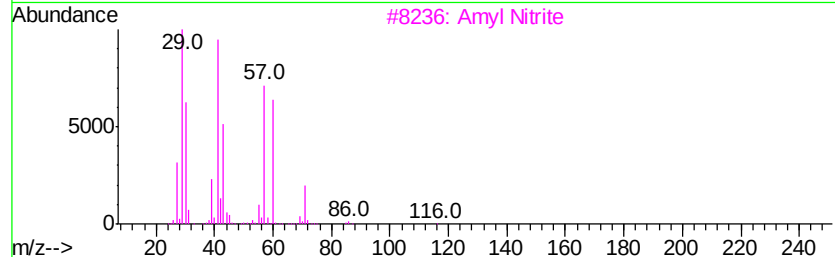
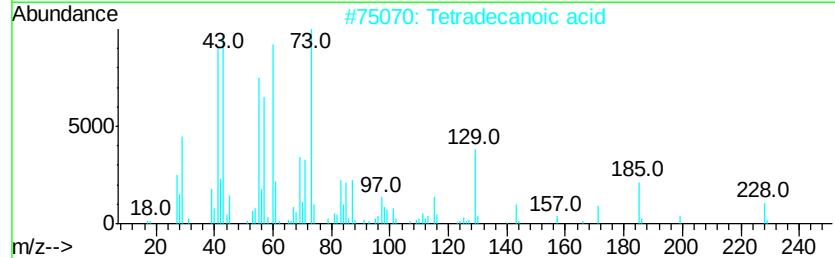
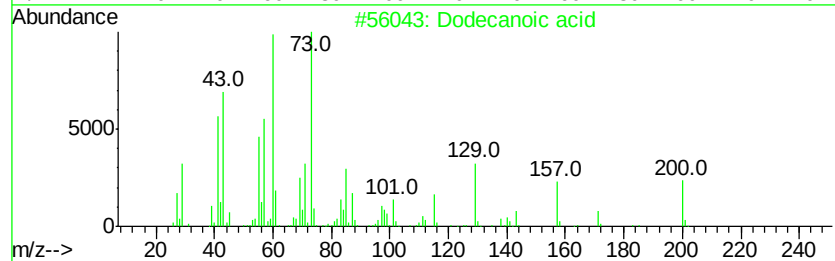
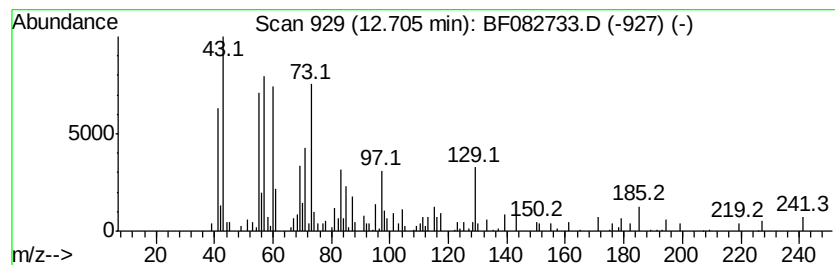
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Dodecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.71	2.20 ng	113385	Phenanthrene-d10	11.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecanoic acid	200	C12H24O2	000143-07-7	53
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	50
3	Amvl Nitrite	117	C5H11NO2	000110-46-3	27
4	Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	11
5	Decane, 2,3,5,8-tetramethyl-	198	C14H30	192823-15-7	11



Data Path : Z:\HPCHEM1\BNA F\DATA\BF110615\
Data File : BF082733.D
Acq On : 5 Nov 2015 17:04
Operator : UM/IZ
Sample : G4261-12
Misc :
ALS Vial : 6 Sample Multiplier: 1

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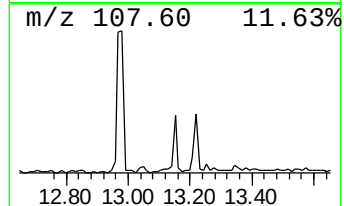
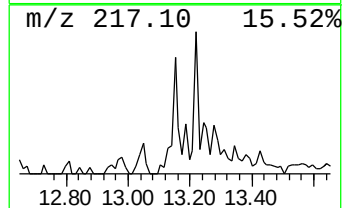
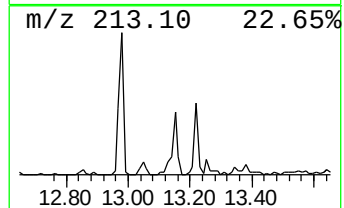
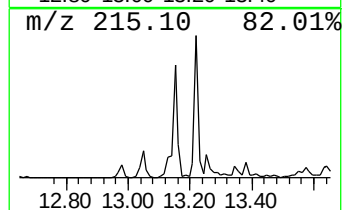
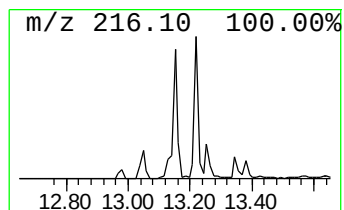
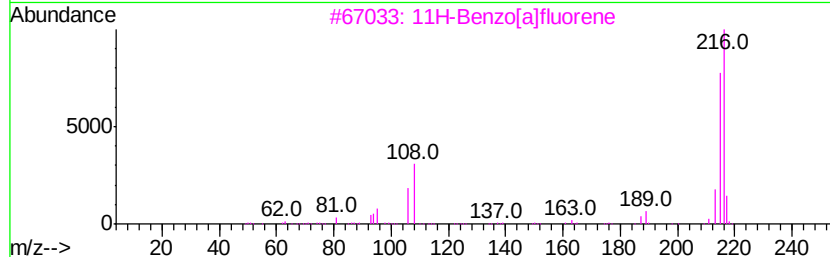
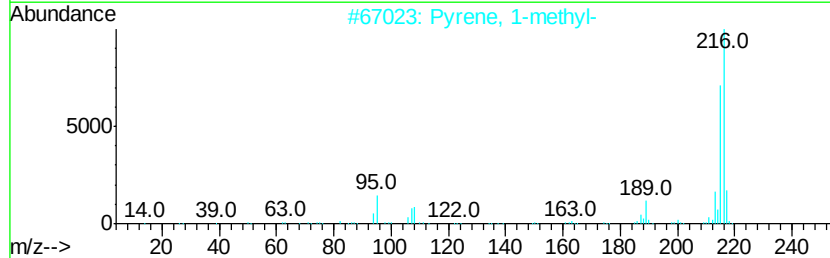
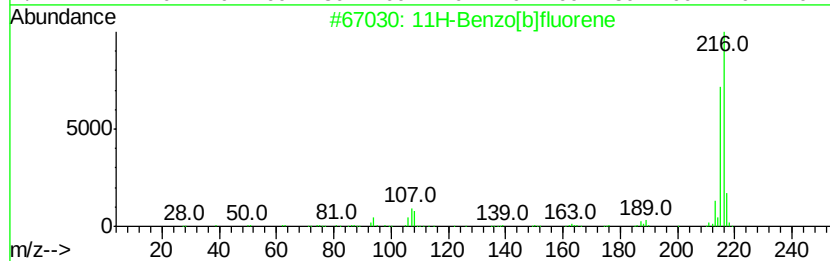
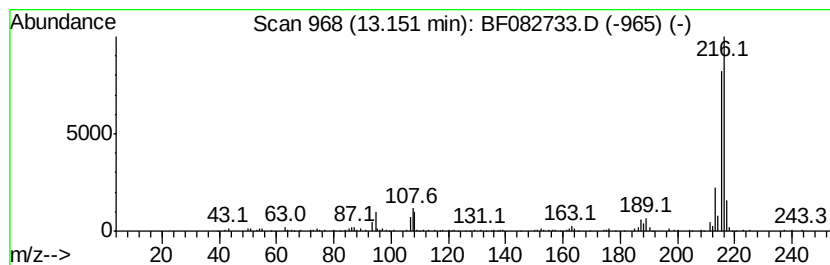
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 11H-Benzo[b]fluorene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.15	3.03 ng	177051	Chrysene-d12	14.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
4			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72



Data Path : Z:\HPCHEM1\BNA F\DATA\BF110615\
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Misc :
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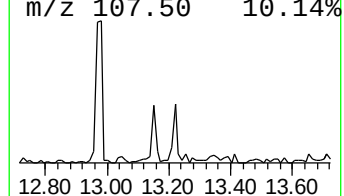
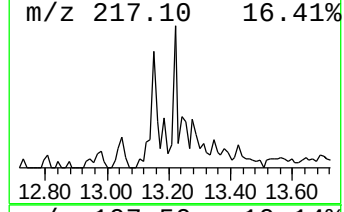
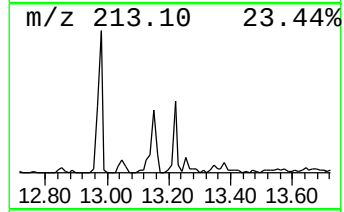
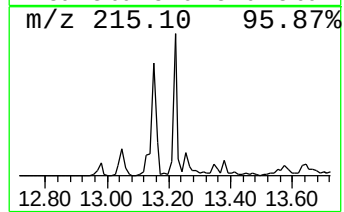
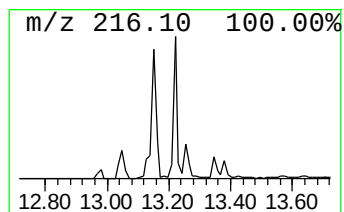
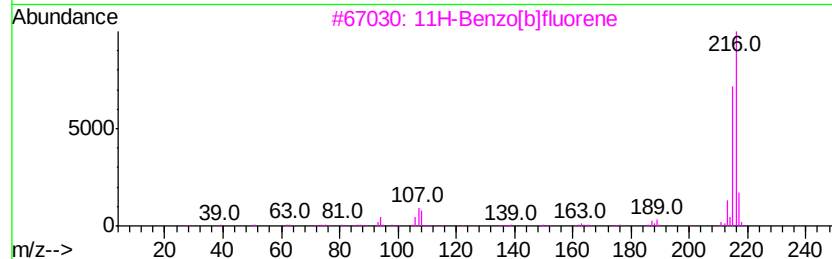
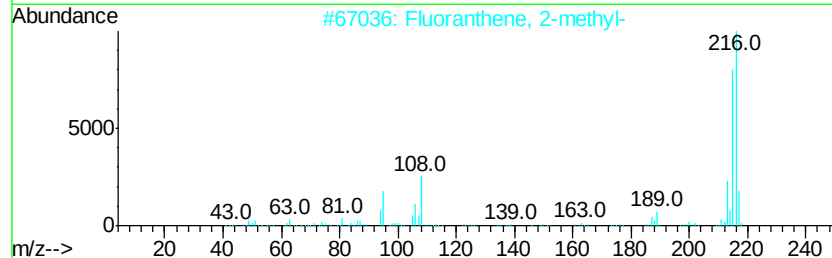
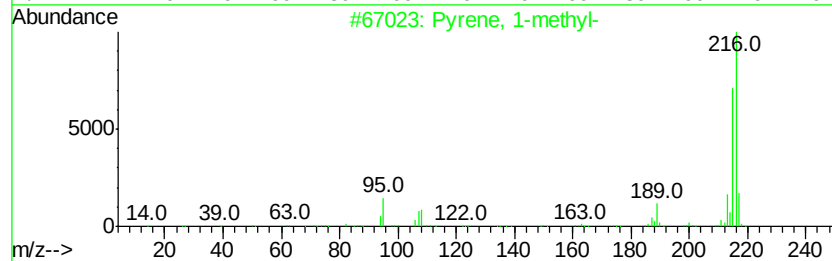
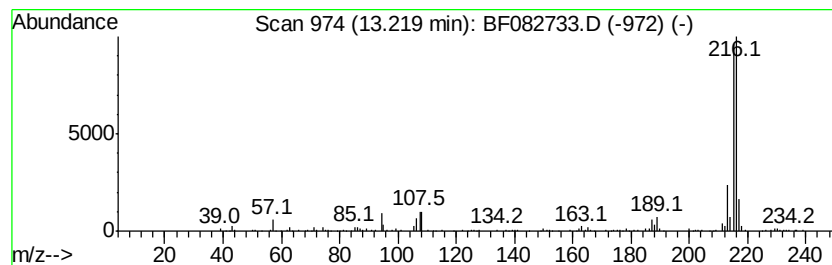
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Pyrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	2.71 ng	158312	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
2		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
3		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
4		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	83
5		7H-Benzo[c]fluorene	216	C17H12	000205-12-9	74



Data Path : Z:\HPCHEM1\BNA F\DATA\BF110615\
Data File : BF082733.D
Acq On : 5 Nov 2015 17:04
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Sample : G4261-12
Misc :
ALS Vial : 6 Sample Multiplier: 1

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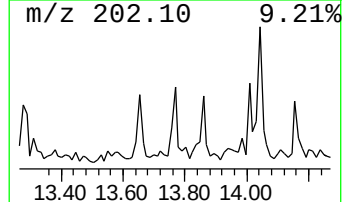
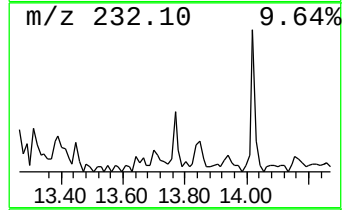
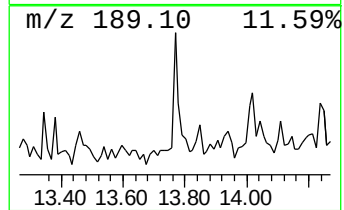
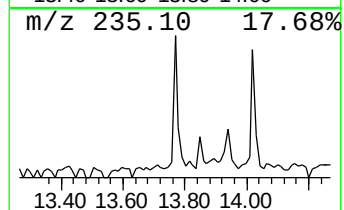
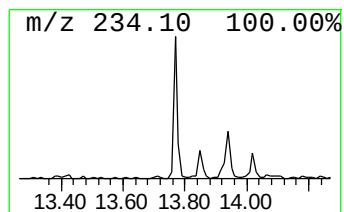
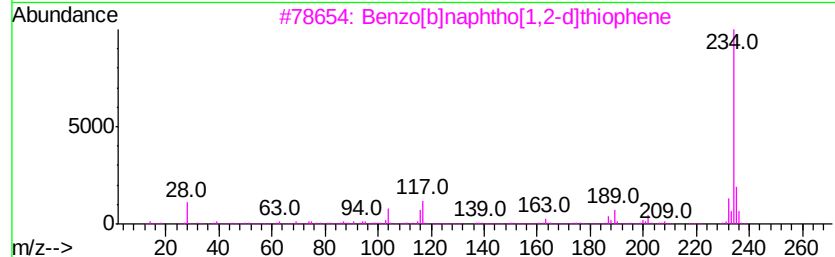
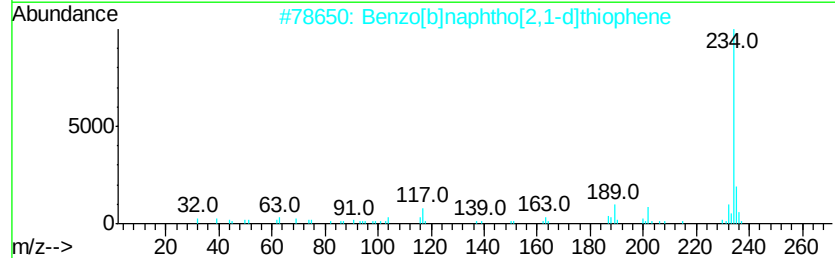
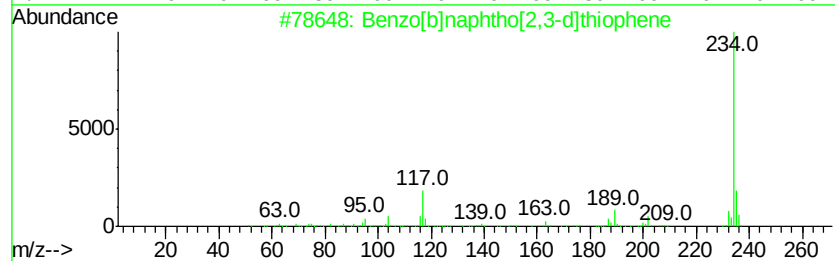
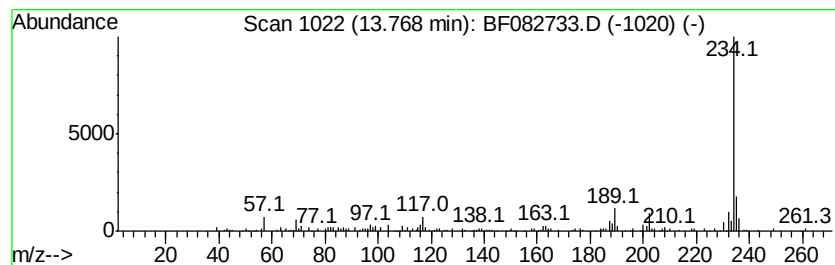
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzo[b]naphtho[2,3-d]thiop... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.77	2.02 ng	117630	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	96
2		Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	95
3		Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	93
4		2-Amino-6-t-butyl-3-cyano-4,5,6,...	234	C13H18N2S	042159-76-2	50
5		4H-1-Benzopyran-2-carboxylic aci...	234	C12H10O5	023866-72-0	42



Data Path : Z:\HPCHEM1\BNA F\DATA\BF110615\
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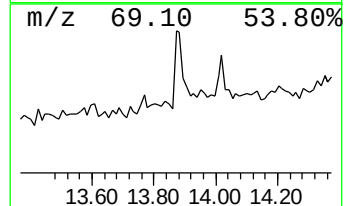
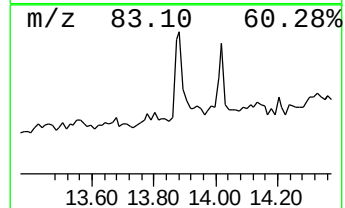
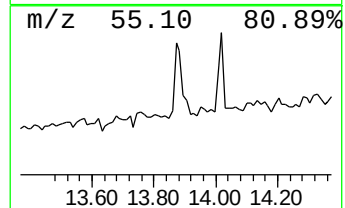
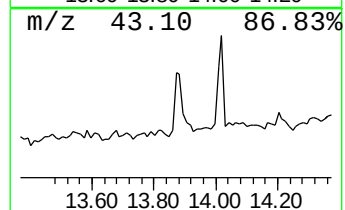
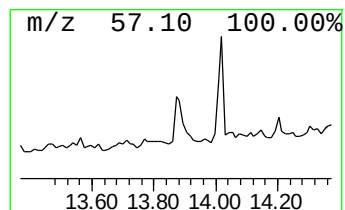
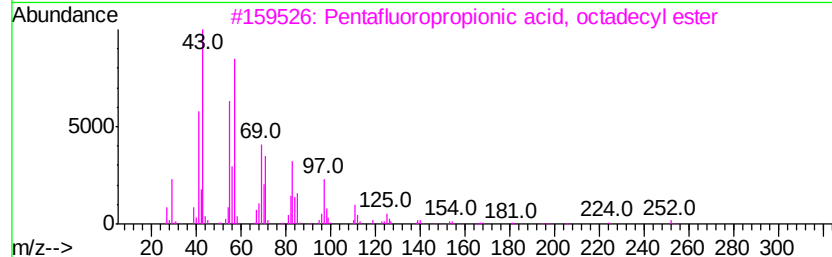
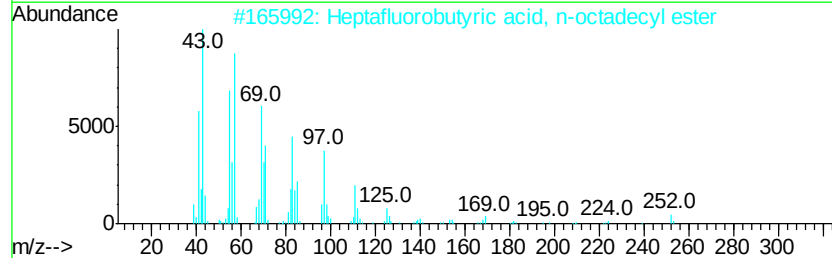
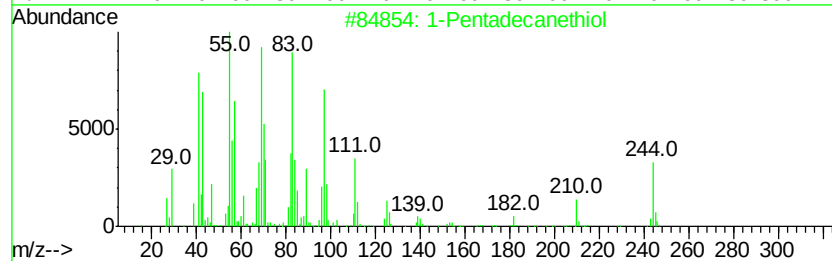
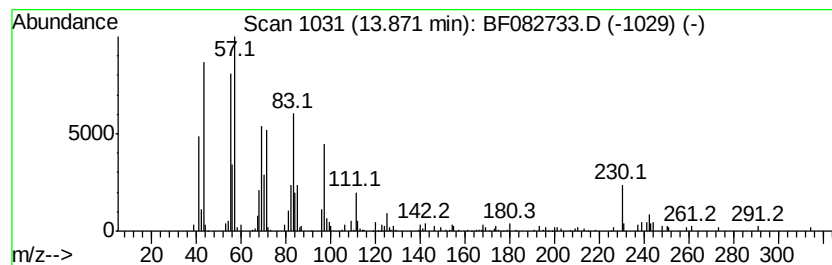
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Pentadecanethiol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	3.42 ng	199807	Chrysene-d12	14.02
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Pentadecanethiol	244 C15H32S	025276-70-4	86
2	Heptafluorobutyric acid, n-octad...	466 C22H37F7O2	000400-57-7	83
3	Pentafluoropropionic acid, octad...	416 C21H37F5O2	1000280-07-7	83
4	17-Pentatriacontene	491 C35H70	006971-40-0	80
5	1-Tetracosanol	354 C24H50O	000506-51-4	74



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110615\
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Sample : G4261-12
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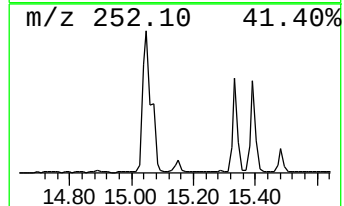
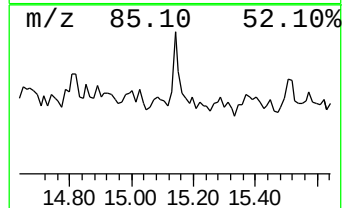
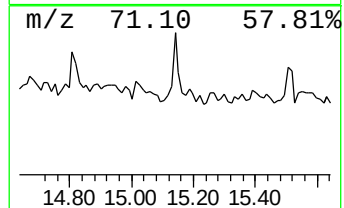
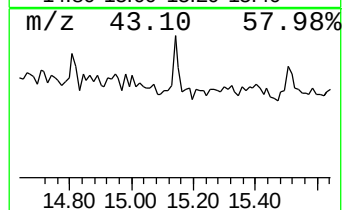
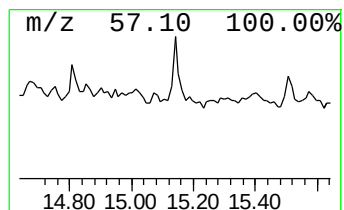
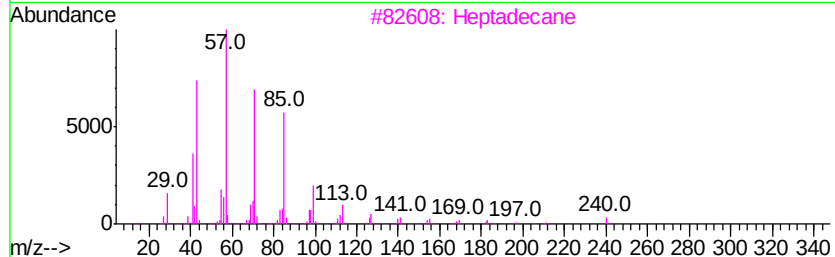
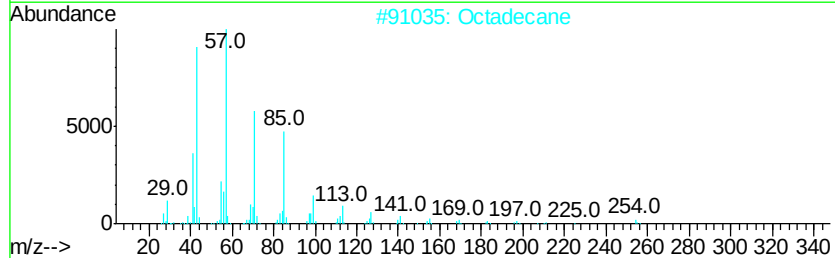
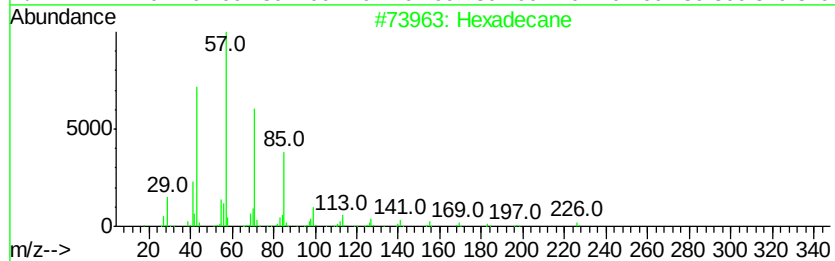
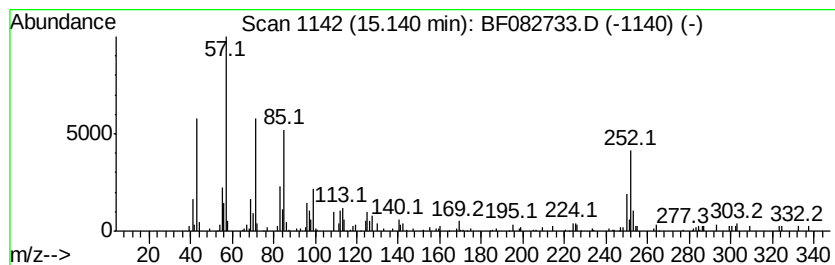
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown15.14 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.14	4.24 ng	94157	Perylene-d12	15.45

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane	226	C16H34	000544-76-3	46
2		Octadecane	254	C18H38	000593-45-3	45
3		Heptadecane	240	C17H36	000629-78-7	45
4		Heneicosane	296	C21H44	000629-94-7	43
5		Docosane	310	C22H46	000629-97-0	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110615\
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1,1-dime...	2.13	6.8	ng	186769	1	6.86	545232	20.0
2-Pentanone, 4-hy...	5.06	103.1	ng	2809580	1	6.86	545232	20.0
unknown6.64	6.64	203.8	ng	5554600	1	6.86	545232	20.0
n-Hexadecanoic acid	11.93	5.4	ng	279686	4	11.39	1028680	20.0
4H-Cyclopenta[def...	12.00	3.2	ng	164315	4	11.39	1028680	20.0
Dodecanoic acid	12.71	2.2	ng	113385	4	11.39	1028680	20.0
11H-Benzo[b]fluorene	13.15	3.0	ng	177051	5	14.02	1166850	20.0
Pyrene, 1-methyl-	13.22	2.7	ng	158312	5	14.02	1166850	20.0
Benzo[b]naphtho[2...	13.77	2.0	ng	117630	5	14.02	1166850	20.0
1-Pentadecanethiol	13.87	3.4	ng	199807	5	14.02	1166850	20.0
unknown15.14	15.14	4.2	ng	94157	6	15.45	443813	20.0