

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092815\
 Data File : BF081864.D
 Acq On : 29 Sep 2015 4:01
 Operator : UM/IZ
 Sample : G3717-03
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-5-9-16-15

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-BF092815.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	5.109	258	260	262	rBV	233977	304622	10.56%	1.313%
2	5.520	293	296	311	rBV	323062	574757	19.92%	2.477%
3	6.549	383	386	396	rBV	166277	376522	13.05%	1.623%
4	6.697	396	399	407	rBV	741326	1227733	42.54%	5.292%
5	6.926	417	419	431	rVV	336632	580483	20.11%	2.502%
6	7.086	431	433	446	rVB	1920242	1999401	69.28%	8.618%
7	7.497	467	469	483	rBV	1307261	1723577	59.72%	7.429%
8	7.817	493	497	502	rBV	49508	162879	5.64%	0.702%
9	7.966	507	510	512	rVB	263308	340611	11.80%	1.468%
10	8.023	512	515	520	rVB4	27213	74509	2.58%	0.321%
11	8.126	522	524	525	rBV	23258	31845	1.10%	0.137%
12	8.217	530	532	542	rVV	675380	869547	30.13%	3.748%
13	8.343	542	543	544	rVV	68178	61708	2.14%	0.266%
14	8.366	544	545	550	rVB	158298	164300	5.69%	0.708%
15	8.583	561	564	568	rBV5	13364	32510	1.13%	0.140%
16	8.937	592	595	599	rBV3	17922	50686	1.76%	0.218%
17	9.029	601	603	607	rVV2	19726	41589	1.44%	0.179%
18	9.292	624	626	629	rBV	2172892	2886000	100.00%	12.439%
19	9.406	633	636	641	rVB2	107297	215813	7.48%	0.930%
20	9.635	653	656	659	rBV3	29684	54291	1.88%	0.234%
21	9.692	659	661	664	rBV	345202	341125	11.82%	1.470%
22	9.978	683	686	688	rBV	860013	892607	30.93%	3.847%
23	10.126	697	699	702	rBV2	14631	31277	1.08%	0.135%
24	10.183	702	704	706	rBV2	41082	53516	1.85%	0.231%
25	10.400	721	723	725	rVB2	51064	58502	2.03%	0.252%
26	10.446	725	727	730	rBV4	15945	35834	1.24%	0.154%
27	10.515	732	733	735	rBV2	22771	30329	1.05%	0.131%
28	10.766	752	755	760	rBV	1050151	1258289	43.60%	5.423%
29	10.869	763	764	769	rVB2	68634	116136	4.02%	0.501%
30	10.972	771	773	775	rVB2	35980	39257	1.36%	0.169%
31	11.063	779	781	783	rBV2	29599	48650	1.69%	0.210%
32	11.132	785	787	791	rBV4	22902	53433	1.85%	0.230%
33	11.326	802	804	809	rVB3	44182	104149	3.61%	0.449%
34	11.463	813	816	820	rBV	877223	1000142	34.65%	4.311%

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Stop Thrs : 0 Peak Location: TOP

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35	11.703	835	837	838 rBV	34940	44073	1.53%	0.190%
36	11.749	839	841	844 rVB	47585	46565	1.61%	0.201%
37	12.001	860	863	868 rBV	996790	1205115	41.76%	5.194%
38	12.435	899	901	902 rBV	43242	44898	1.56%	0.194%
39	12.469	902	904	907 rVV	63906	89525	3.10%	0.386%
40	12.526	907	909	912 rVB3	38647	52597	1.82%	0.227%
41	12.686	921	923	927 rBV4	53959	146632	5.08%	0.632%
42	12.789	929	932	938 rVV	342409	643319	22.29%	2.773%
43	12.949	944	946	949 rVB	95027	148394	5.14%	0.640%
44	13.052	953	955	958 rBV	2676460	2778321	96.27%	11.975%
45	13.132	961	962	965 rBV	77715	88240	3.06%	0.380%
46	13.612	1002	1004	1006 rBV	41863	49355	1.71%	0.213%
47	13.955	1031	1034	1040 rBV	173002	291691	10.11%	1.257%
48	14.104	1045	1047	1050 rVB	718417	844731	29.27%	3.641%
49	14.950	1120	1121	1126 rVB	44497	63376	2.20%	0.273%
50	15.578	1173	1176	1184 rVB	517142	728419	25.24%	3.140%
51	16.104	1219	1222	1225 rBV3	22112	41601	1.44%	0.179%
52	17.635	1352	1356	1361 rBV5	21793	57543	1.99%	0.248%

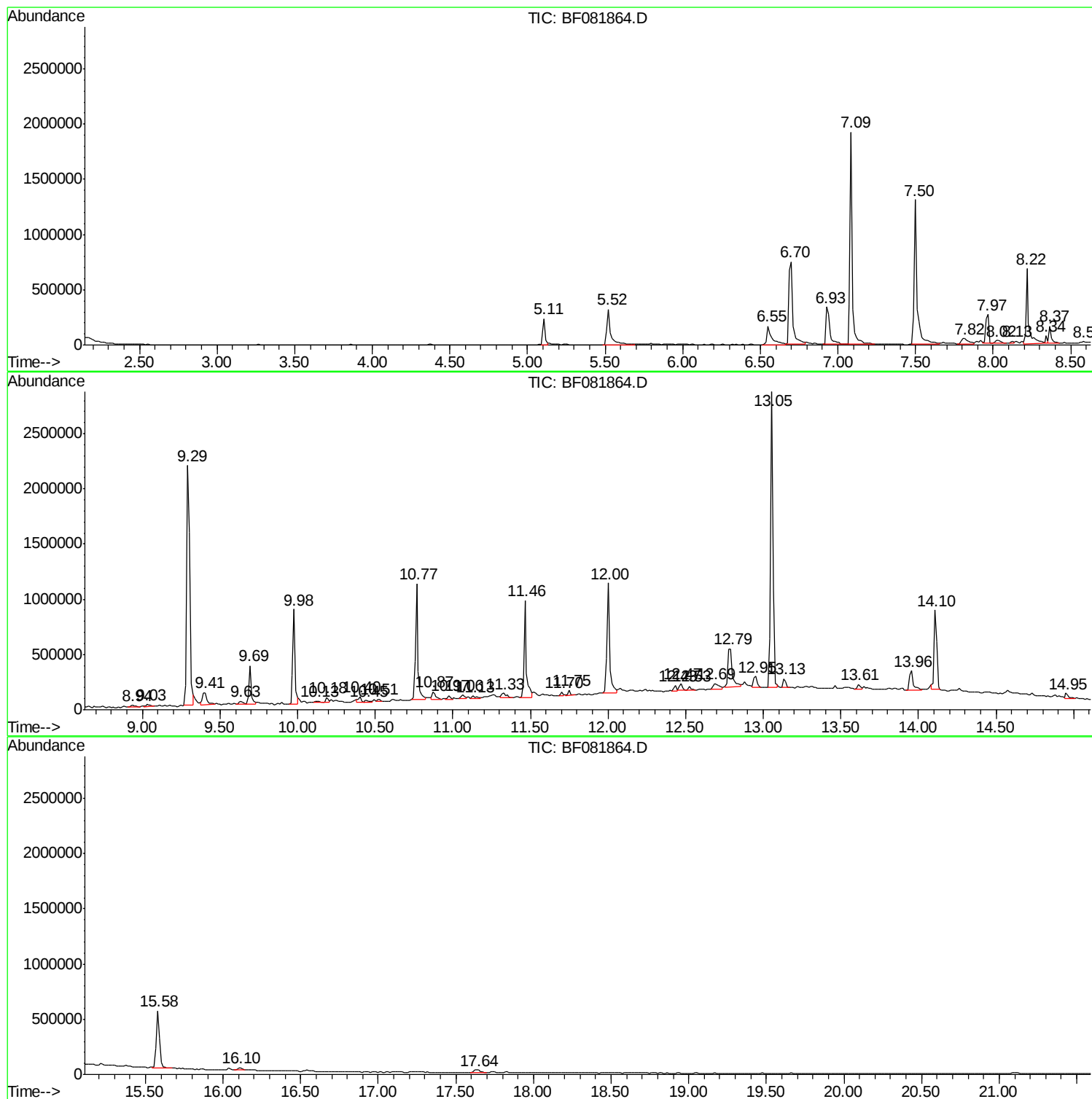
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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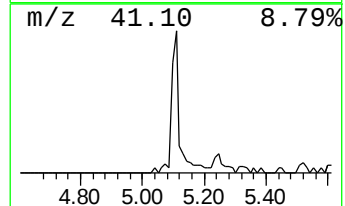
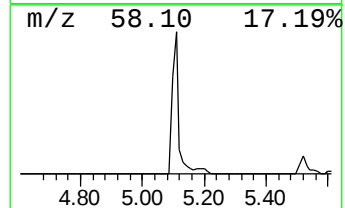
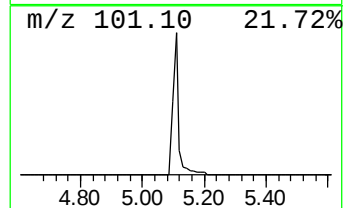
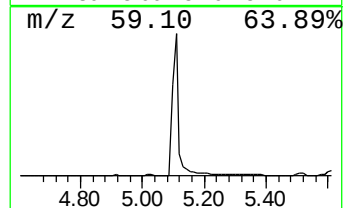
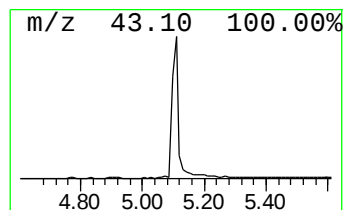
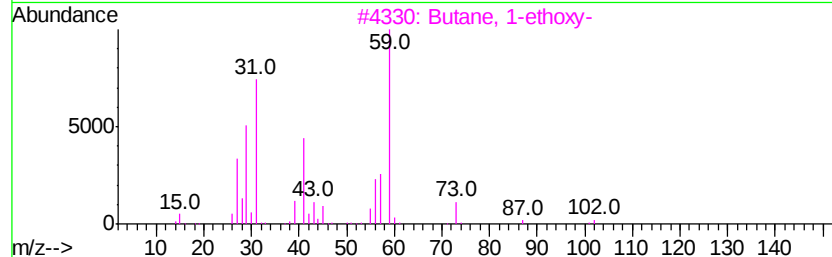
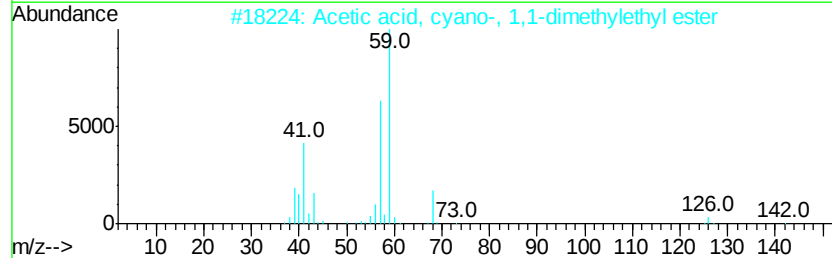
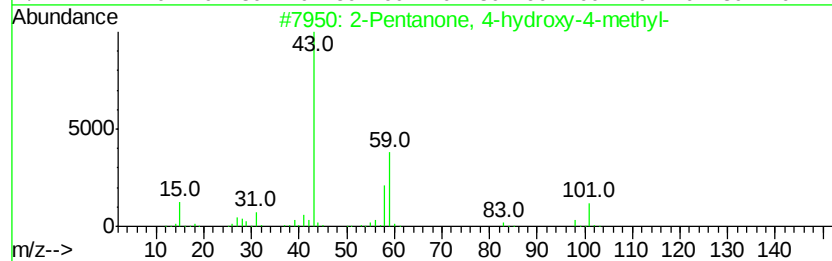
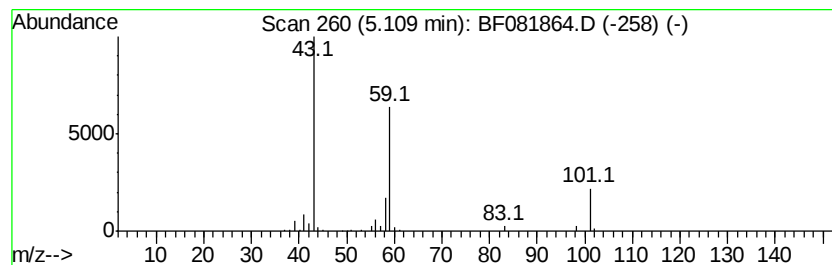
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	10.50 ng	304622	1,4-Dichlorobenzene-d4	6.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Guanidine	59	CH5N3	000113-00-8	9



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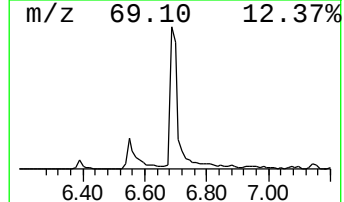
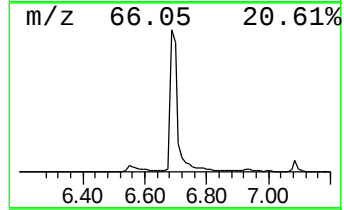
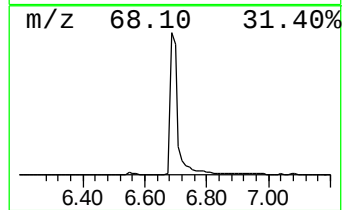
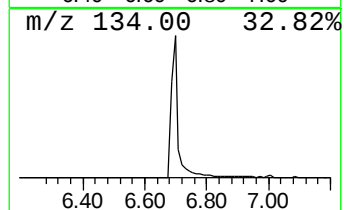
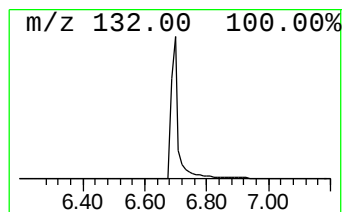
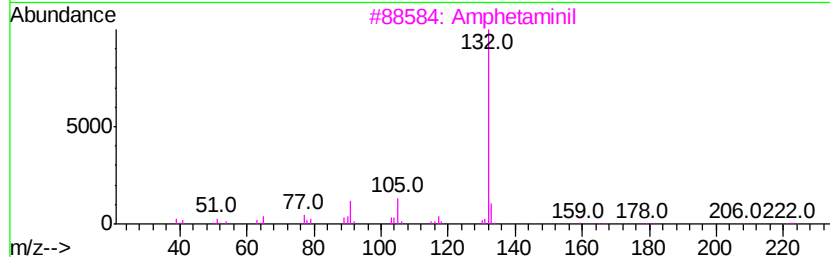
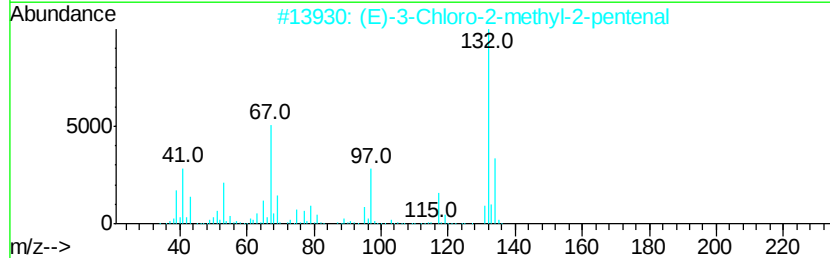
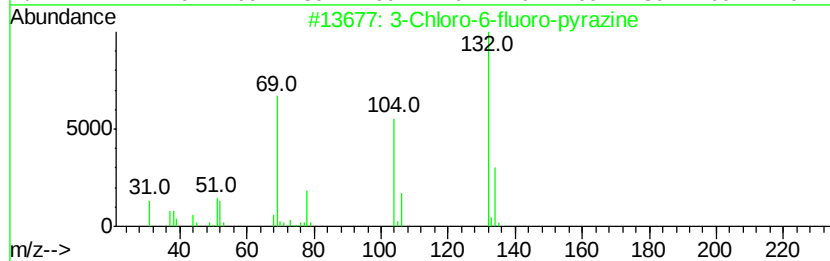
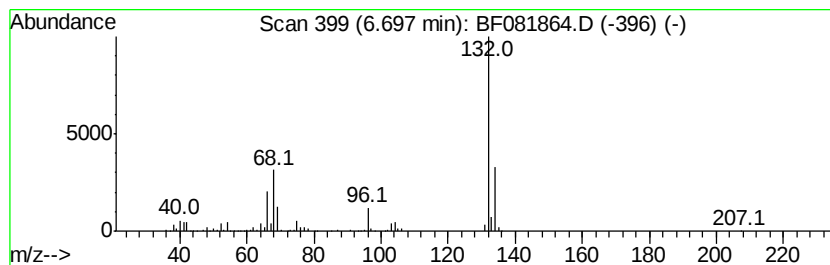
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.70 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	42.30 ng	1227730	1,4-Dichlorobenzene-d4	6.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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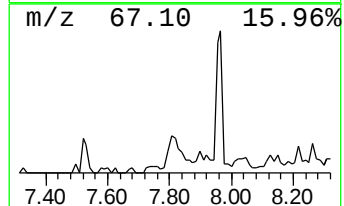
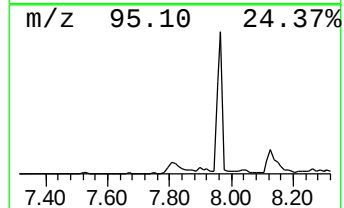
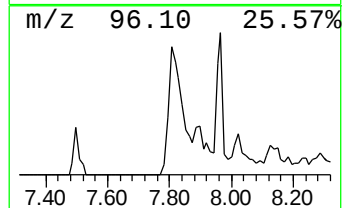
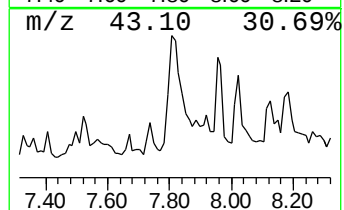
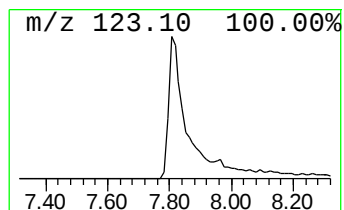
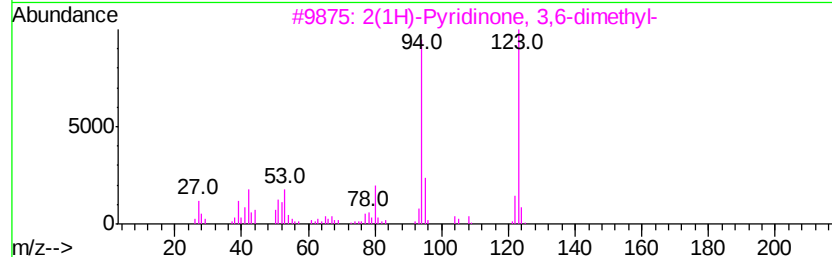
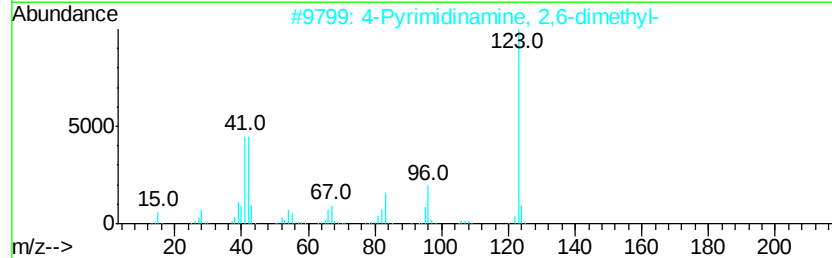
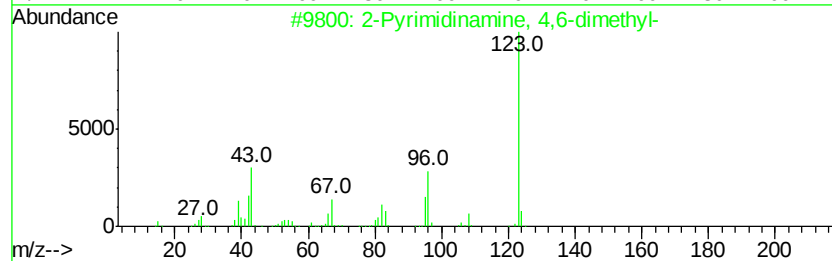
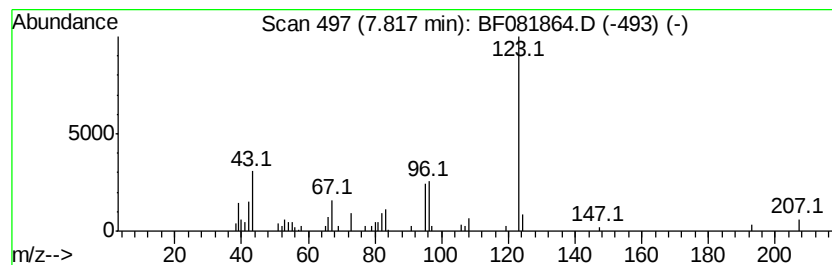
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 2-Pyrimidinamine, 4,6-dimet... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.82	3.75 ng	162879	Napthalene-d8	8.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pyrimidinamine, 4,6-dimethyl-	123	C6H9N3	000767-15-7	93
2		4-Pyrimidinamine, 2,6-dimethyl-	123	C6H9N3	000461-98-3	68
3		2(1H)-Pyridinone, 3,6-dimethyl-	123	C7H9N0	053428-02-7	50
4		3,4-dimethyl-1H-pyrrole-2-carbox...	123	C7H9N0	019713-89-4	46
5		Pyridine, 1,2,3,6-tetrahydro-1-(...	153	C9H15N0	057150-46-6	45



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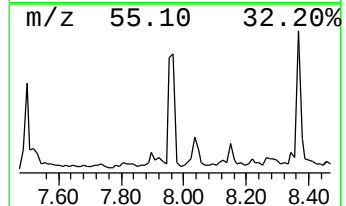
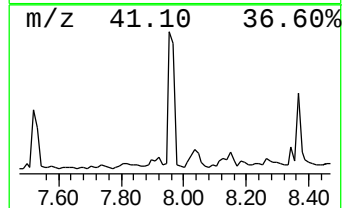
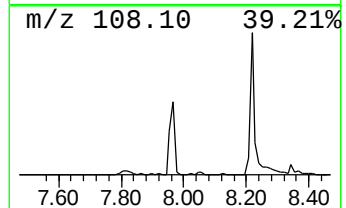
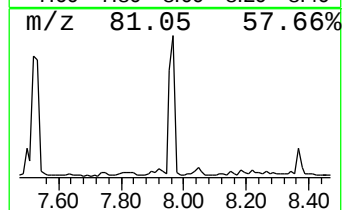
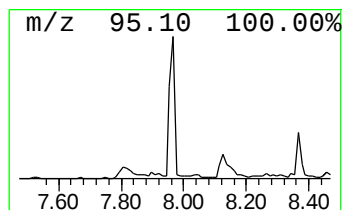
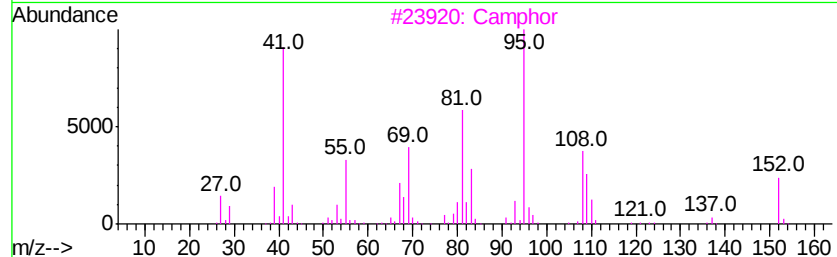
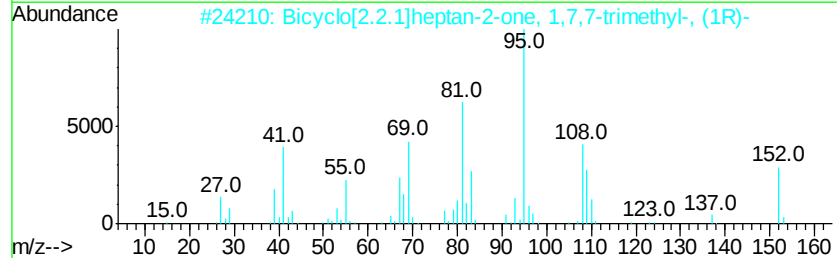
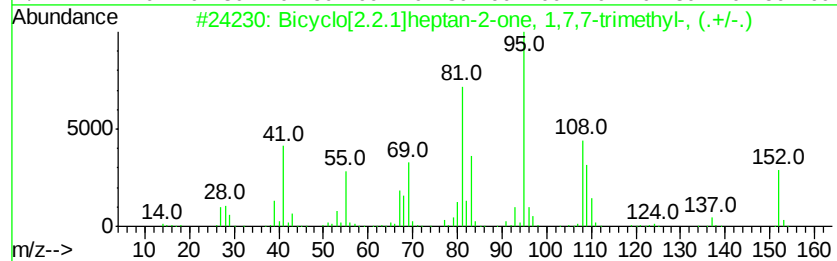
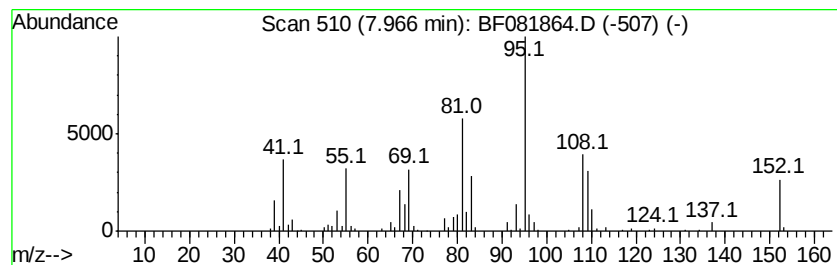
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.97	7.83 ng	340611	Naphthalene-d8	8.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	021368-68-3	97
2		Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-49-3	97
3		Camphor	152	C10H16O	000076-22-2	96
4		Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	93
5		2-Butanone, 4-(5-methyl-2-furanyl)-	152	C9H12O2	013679-56-6	43



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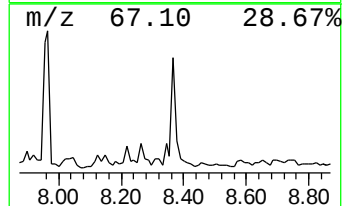
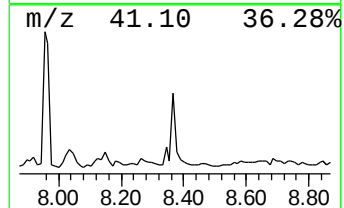
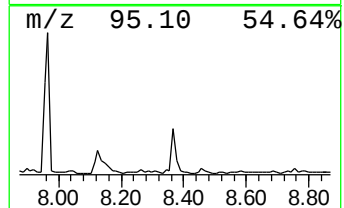
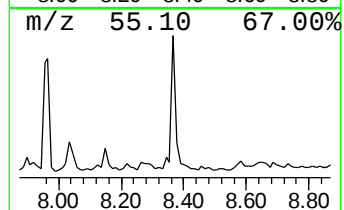
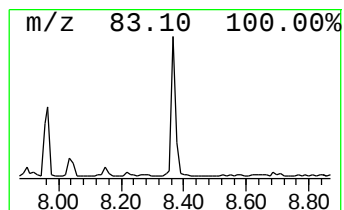
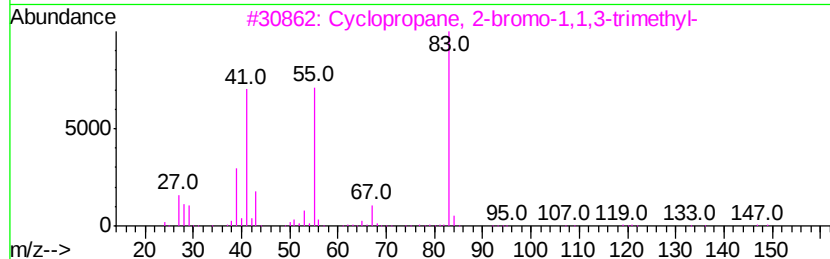
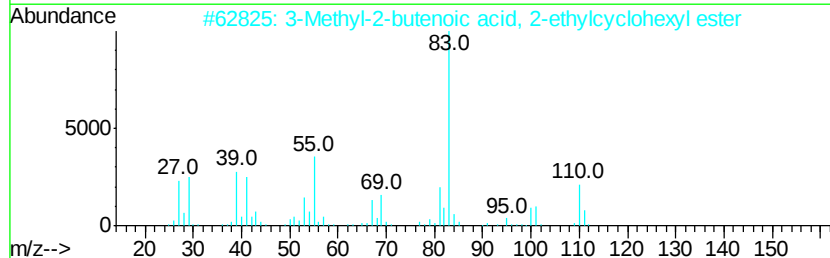
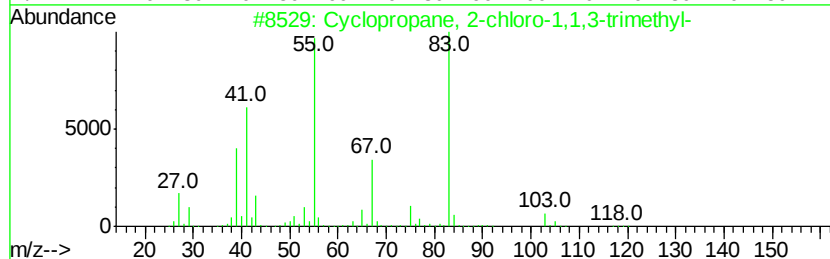
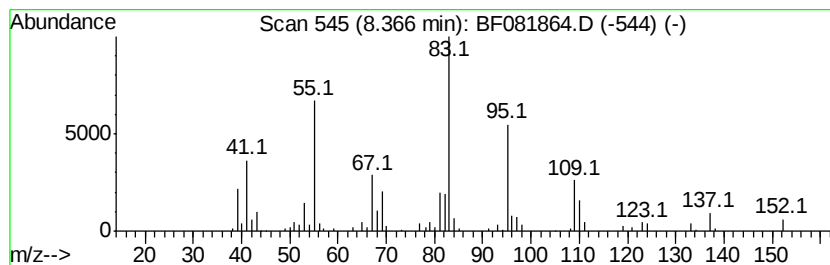
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown8.37 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.37	3.78 ng	164300	Napthalene-d8	8.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropane, 2-chloro-1,1,3-tri...	118	C6H11Cl	098485-99-5	38
2		3-Methyl-2-butenic acid, 2-ethy...	210	C13H22O2	1000279-23-5	38
3		Cyclopropane, 2-bromo-1,1,3-trim...	162	C6H11Br	036617-00-2	35
4		Cyclopentane, 1,2-dimethyl-3-met...	110	C8H14	1000150-99-3	30
5		Bicyclo[2.2.1]heptane, 2-butyl-	152	C11H20	061177-16-0	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092815\
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Acq On : 29 Sep 2015 4:01
Operator : UM/IZ
Sample : G3717-03
Misc :
ALS Vial : 24 Sample Multiplier: 1

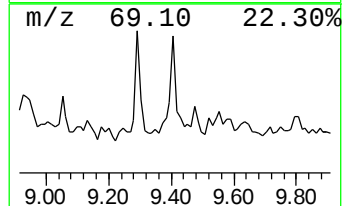
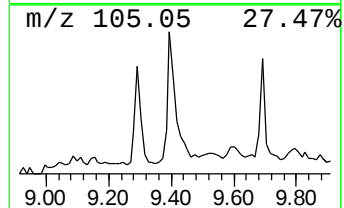
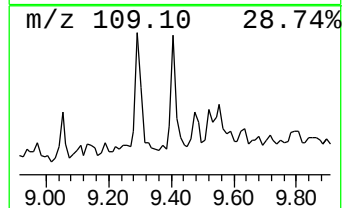
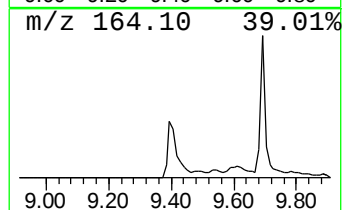
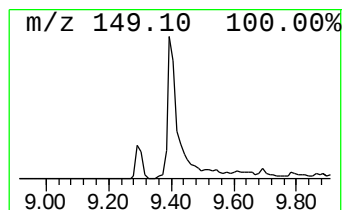
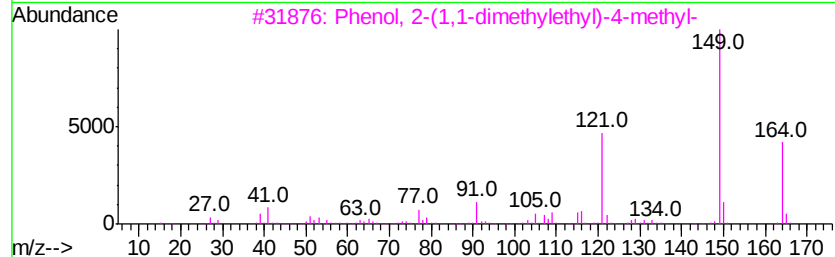
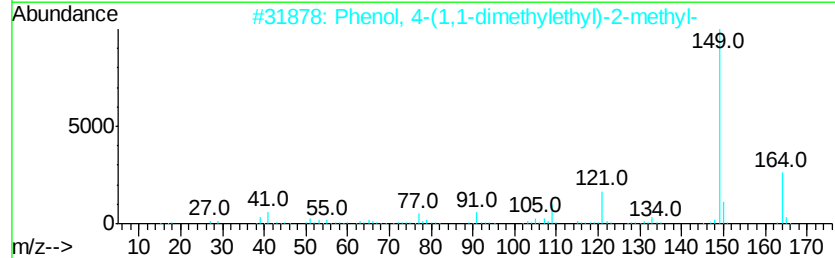
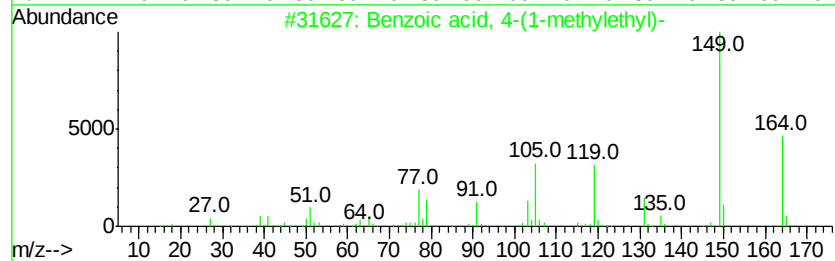
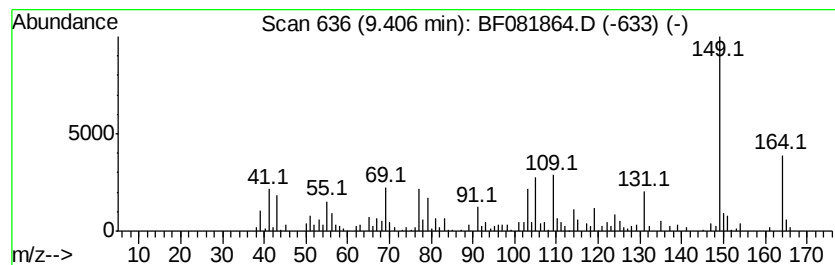
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Benzoic acid, 4-(1-methylethyl) Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.41	4.84 ng	215813	Acenaphthene-d10	9.98
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzoic acid, 4-(1-methylethyl)-	164 C10H12O2	000536-66-3	76
2	Phenol, 4-(1,1-dimethylethyl)-2-...	164 C11H16O	000098-27-1	52
3	Phenol, 2-(1,1-dimethylethyl)-4-...	164 C11H16O	002409-55-4	49
4	Phenol, 2-(1,1-dimethylethyl)-5-...	164 C11H16O	000088-60-8	47
5	Phenol, 2-(1,1-dimethylethyl)-6-...	164 C11H16O	002219-82-1	46



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092815\
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Instrument :
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ClientSampleId :
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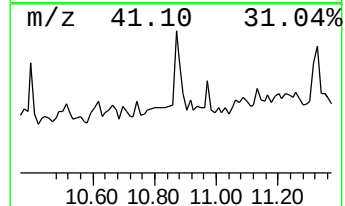
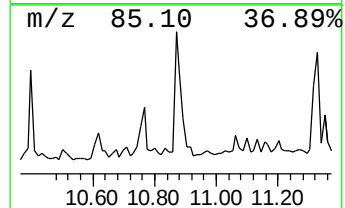
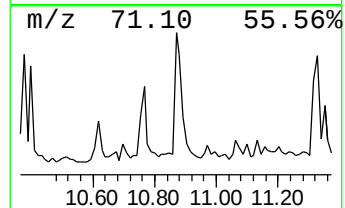
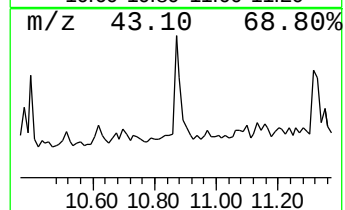
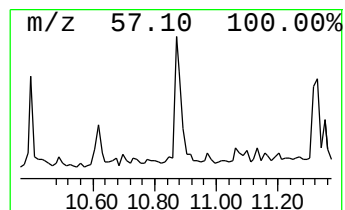
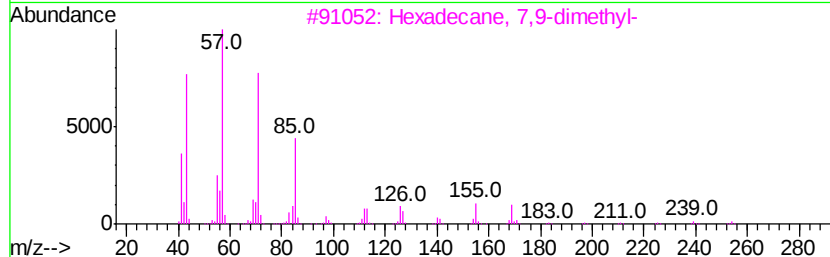
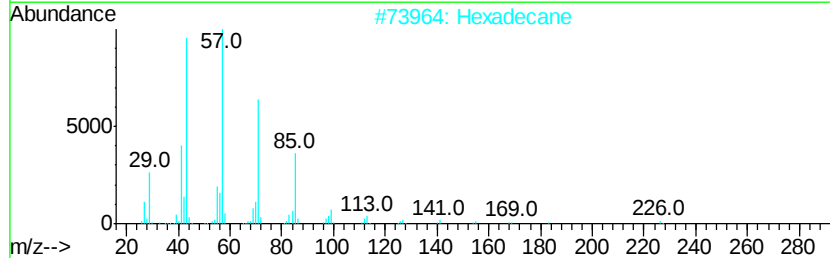
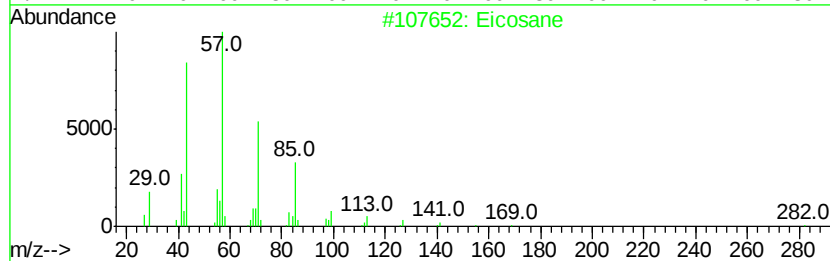
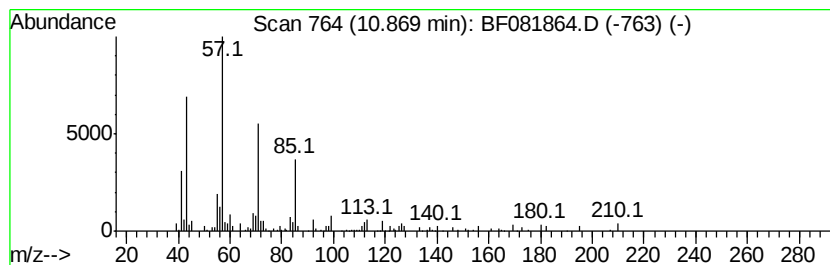
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Eicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.87	2.32 ng	116136	Phenanthrene-d10	11.46

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	64
2	Hexadecane		226	C16H34	000544-76-3	64
3	Hexadecane, 7,9-dimethyl-		254	C18H38	021164-95-4	64
4	Tetradecane, 1-iodo-		324	C14H29I	019218-94-1	59
5	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092815\
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ClientSampleId :
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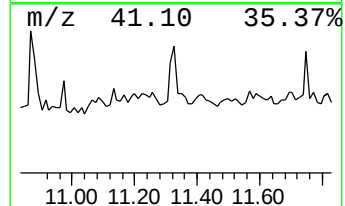
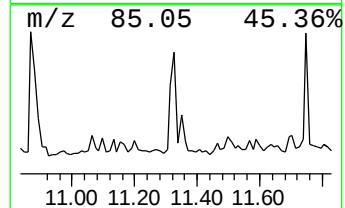
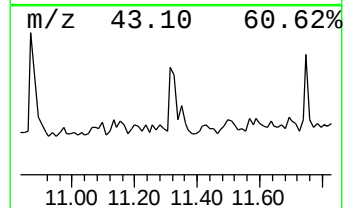
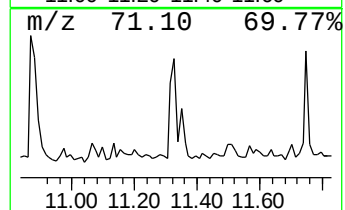
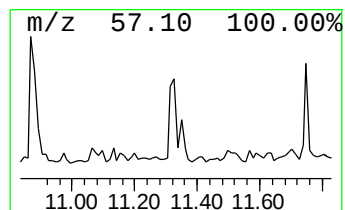
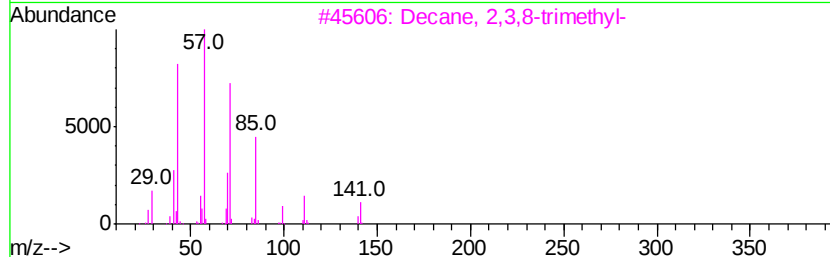
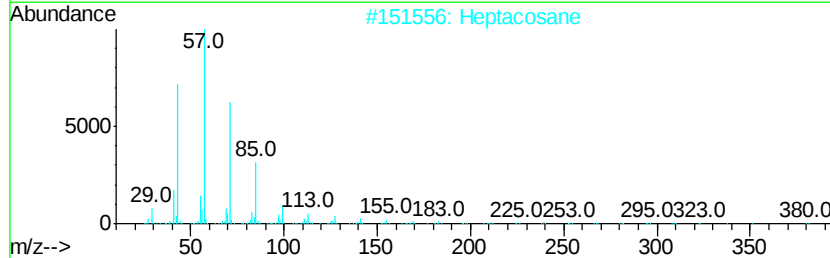
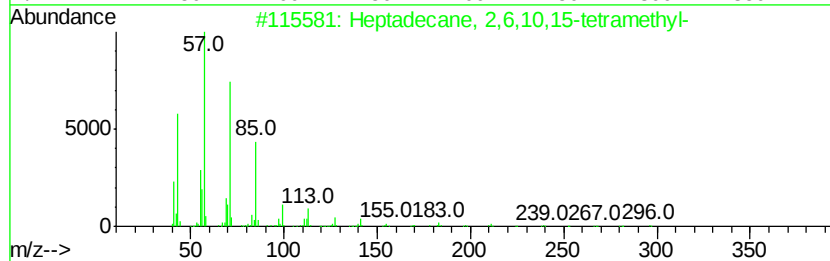
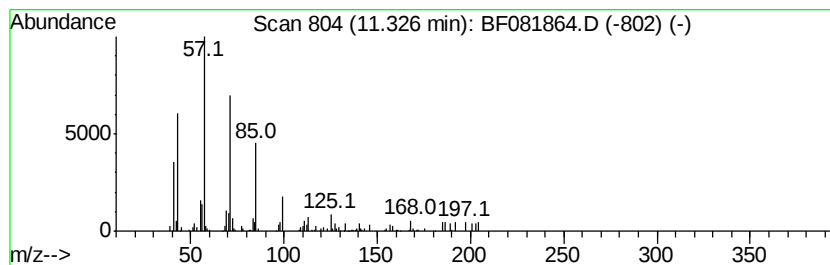
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptadecane, 2,6,10,15-tetr... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.33	2.08 ng	104149	Phenanthrene-d10	11.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	68
2		Heptacosane	380	C27H56	000593-49-7	59
3		Decane, 2,3,8-trimethyl-	184	C13H28	062238-14-6	59
4		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	59
5		Tetradecane	198	C14H30	000629-59-4	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092815\
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Acq On : 29 Sep 2015 4:01
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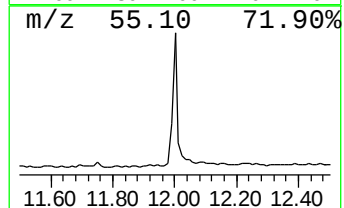
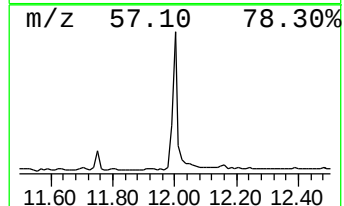
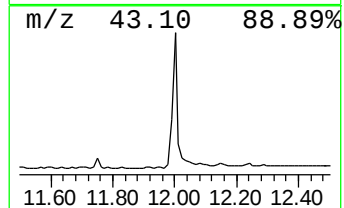
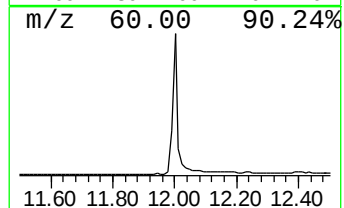
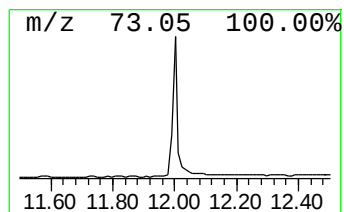
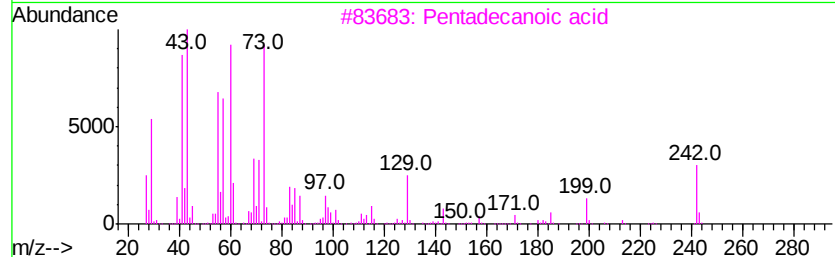
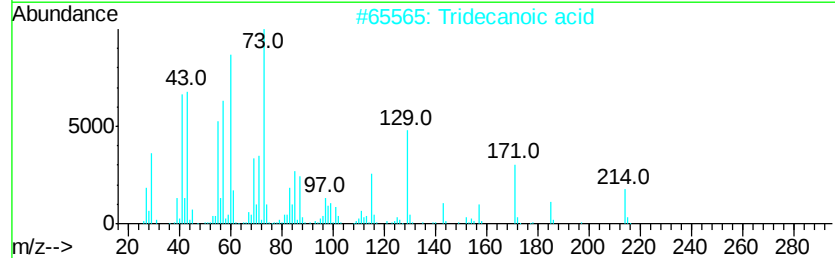
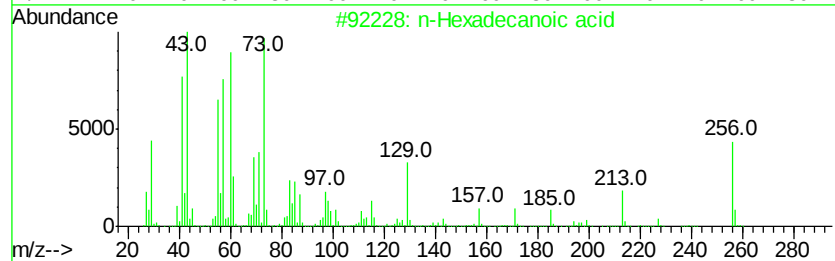
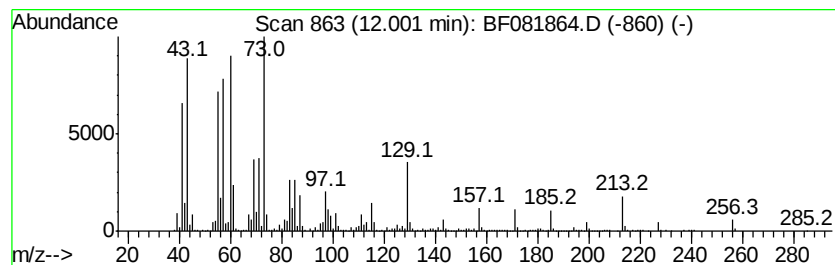
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	24.10 ng	1205120	Phenanthrene-d10	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tridecanoic acid	214	C13H26O2	000638-53-9	96
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	86
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
5	Dodecanoic acid	200	C12H24O2	000143-07-7	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092815\
Data File : BF081864.D
Acq On : 29 Sep 2015 4:01
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Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-5-9-16-15

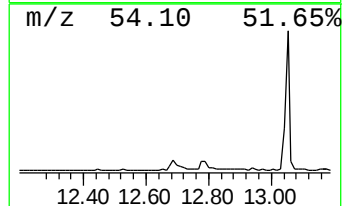
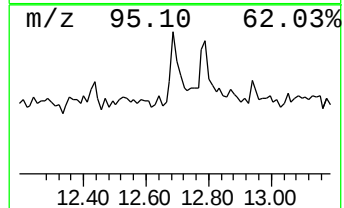
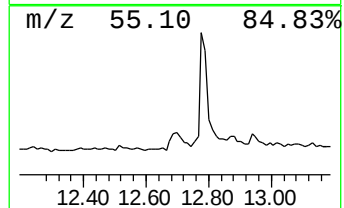
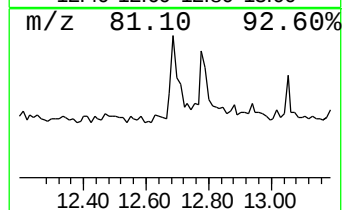
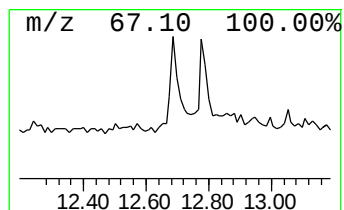
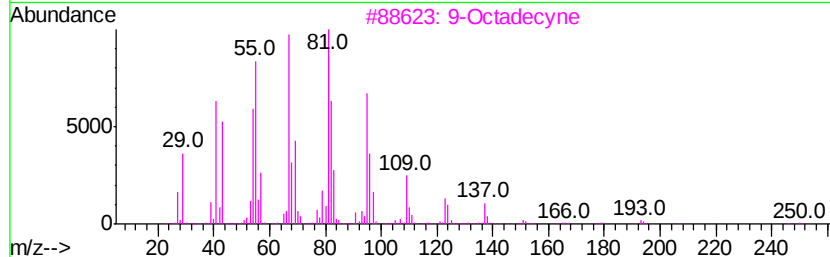
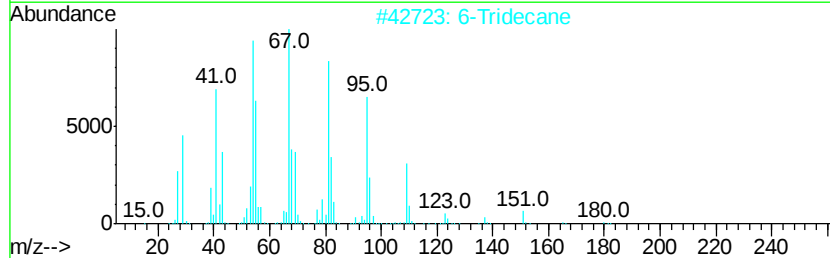
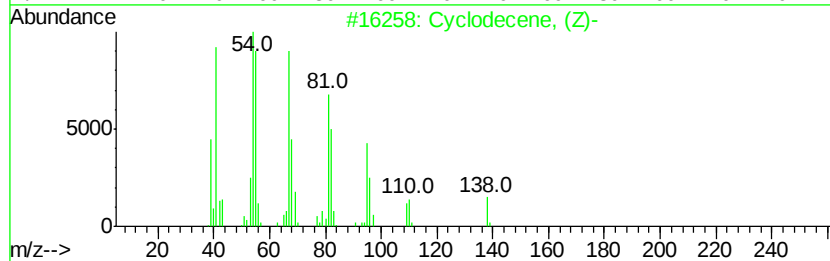
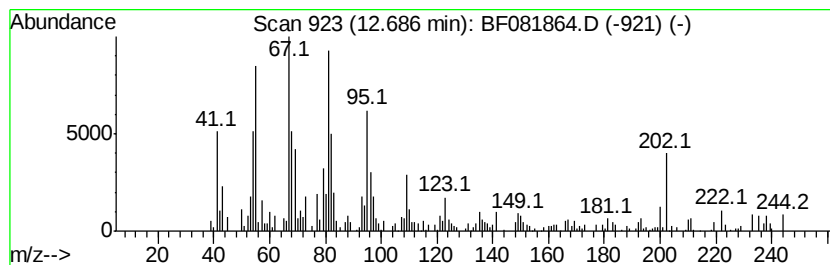
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Cyclodecene, (Z)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.69	2.93 ng	146632	Phenanthrene-d10	11.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclodecene, (Z)-	138	C10H18	000935-31-9	81
2	6-Tridecane	180	C13H24	042371-66-4	68
3	9-Octadecyne	250	C18H34	035365-59-4	64
4	7-Hexadecyne	222	C16H30	074685-28-2	64
5	7-Heptadecyne, 1-chloro-	270	C17H31Cl	056554-74-6	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092815\
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ClientSampleId :
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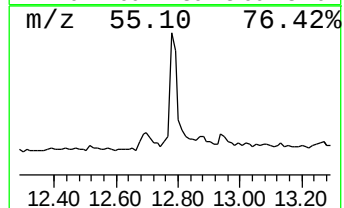
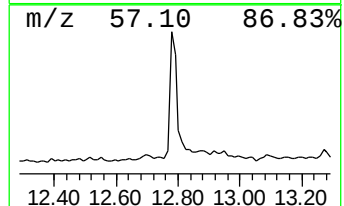
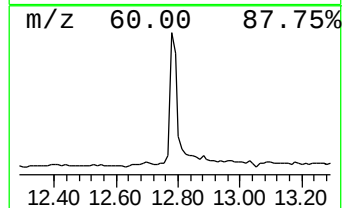
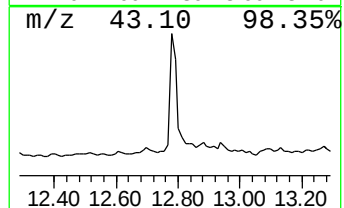
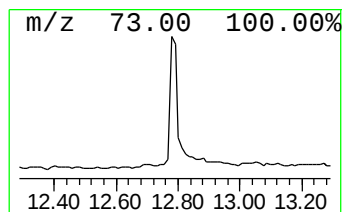
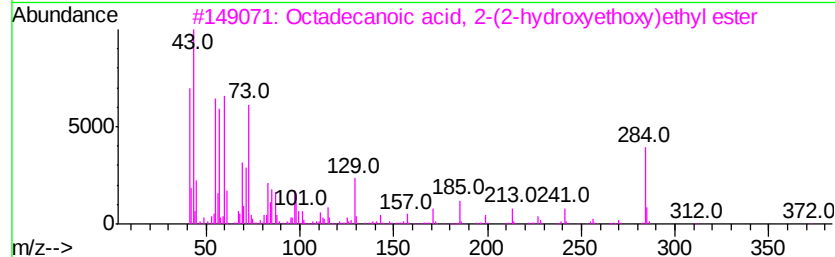
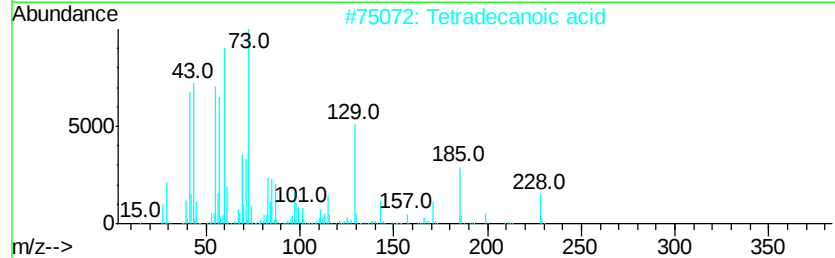
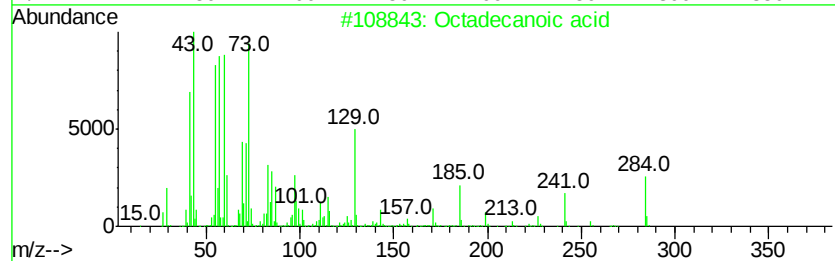
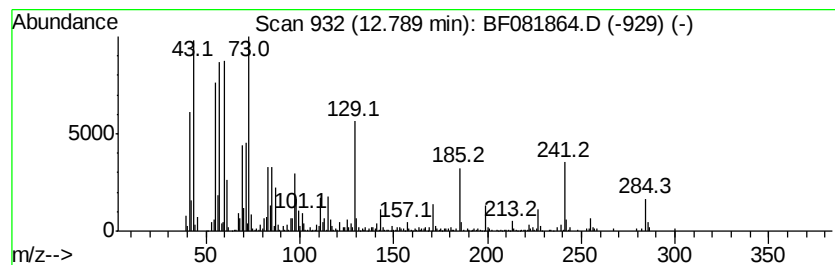
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.79	15.23 ng	643319	Chrysene-d12	14.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	74
3		Octadecanoic acid, 2-(2-hydroxyve...	372	C22H44O4	000106-11-6	72
4		Tridecanoic acid	214	C13H26O2	000638-53-9	72
5		Undecanoic acid	186	C11H22O2	000112-37-8	49



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092815\
Data File : BF081864.D
Acq On : 29 Sep 2015 4:01
Operator : UM/IZ
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Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-5-9-16-15

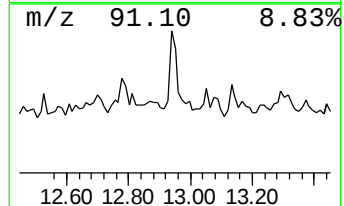
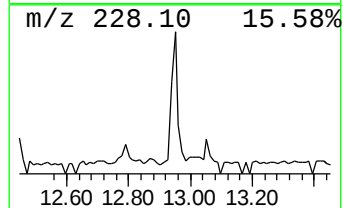
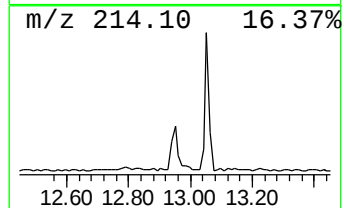
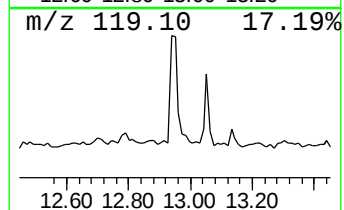
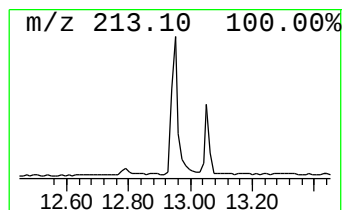
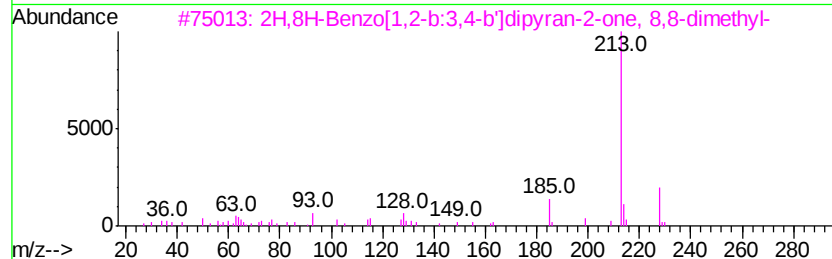
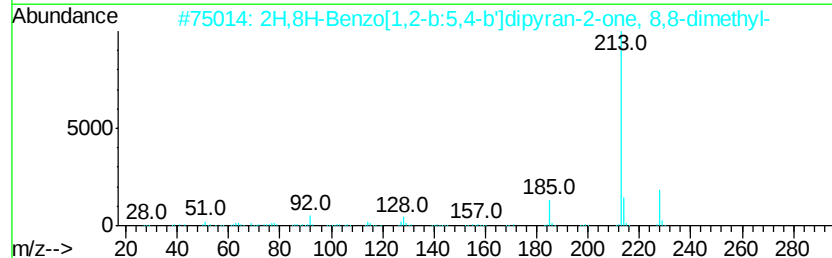
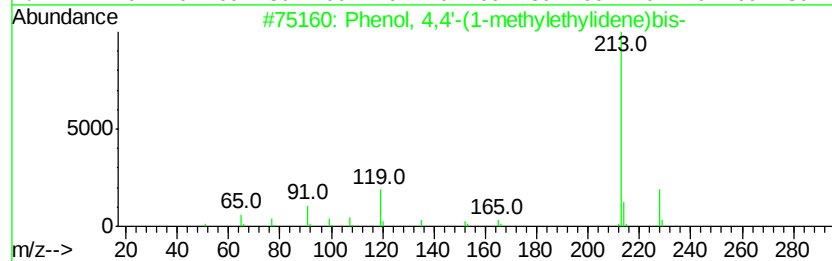
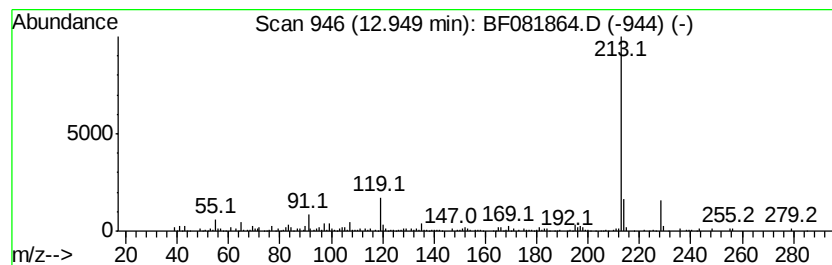
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Phenol, 4,4'-(1-methylethyl... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.95	3.51 ng	148394	Chrysene-d12	14.10

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	93
2			2H,8H-Benzo[1,2-b:5,4-b']dipyran...	228	C14H12O3	000553-19-5	64
3			2H,8H-Benzo[1,2-b:3,4-b']dipyran...	228	C14H12O3	000523-59-1	50
4			2,2-Bis-(4-hydroxyphenyl)-butane	242	C16H18O2	000077-40-7	50
5			9-Propanoyl-1,2,3,4,5,6,7,8-octa...	242	C17H22O	1000255-01-4	42



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092815\
Data File : BF081864.D
Acq On : 29 Sep 2015 4:01
Operator : UM/IZ
Sample : G3717-03
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-5-9-16-15

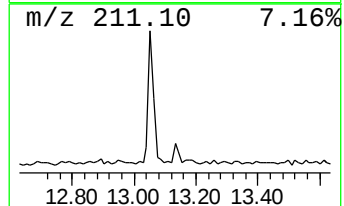
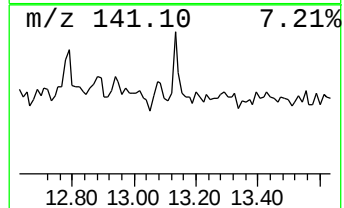
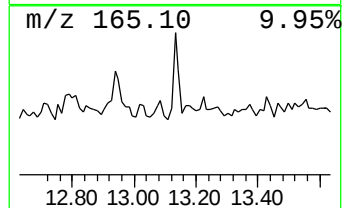
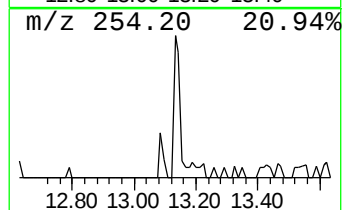
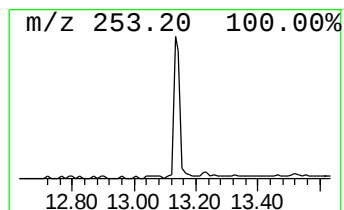
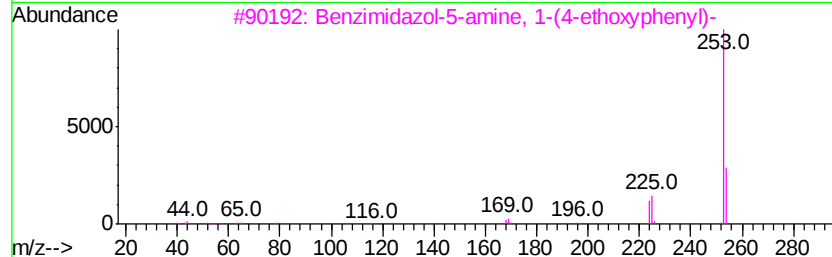
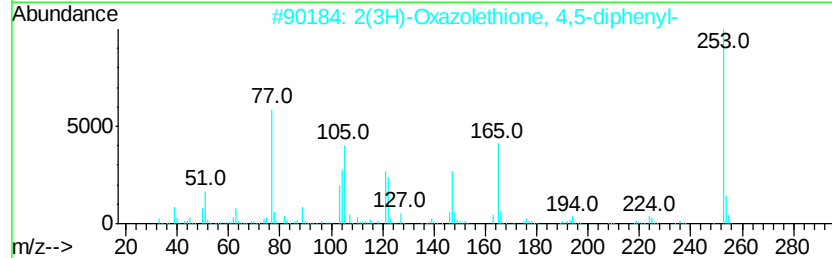
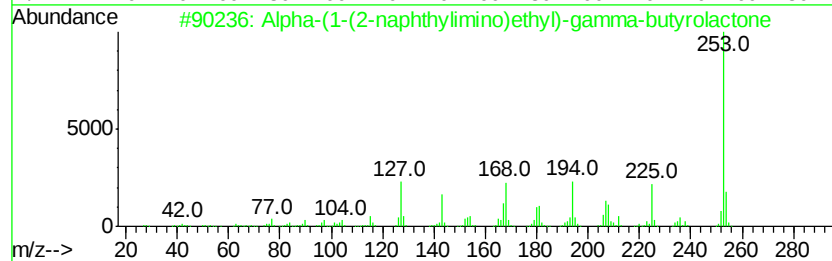
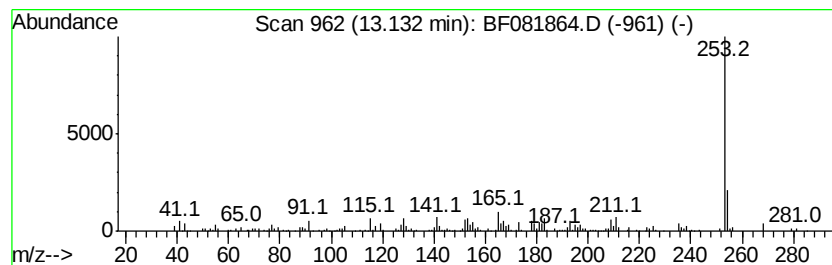
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Alpha-(1-(2-naphthylimino)e... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.13	2.09 ng	88240	Chrysene-d12	14.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Alpha-(1-(2-naphthylimino)ethyl)...	253	C16H15NO2	1000240-01-3	50
2		2(3H)-Oxazolethione, 4,5-diphenyl-	253	C15H11NOS	006670-13-9	50
3		Benzimidazol-5-amine, 1-(4-ethox...	253	C15H15N3O	007104-62-3	45
4		3,6,6-Trimethyl-1-O-tolyl-1,5,6,...	268	C17H20N2O	1000295-04-0	45
5		1,5-Diphenyl-2H-1,2,4-triazoline...	253	C14H11N3S	005055-74-3	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF092815\
Data File : BF081864.D
Acq On : 29 Sep 2015 4:01
Operator : UM/IZ
Sample : G3717-03
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-5-9-16-15

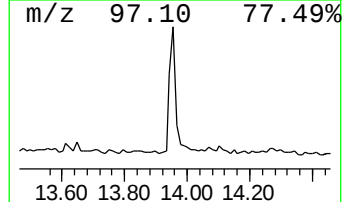
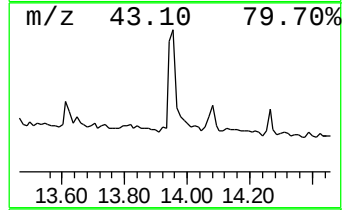
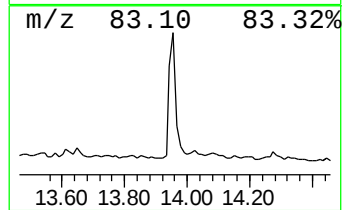
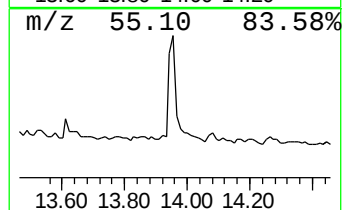
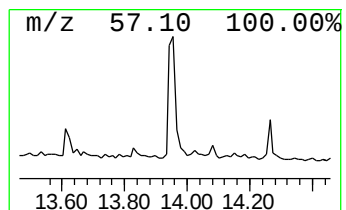
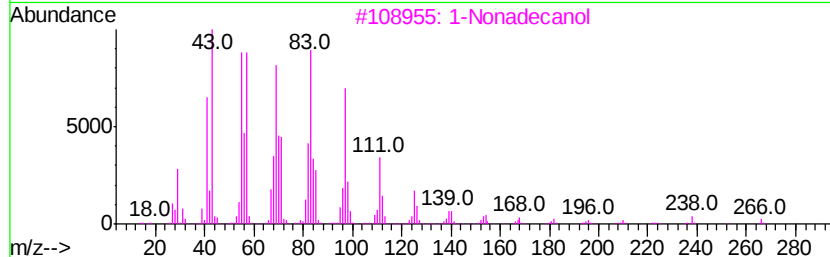
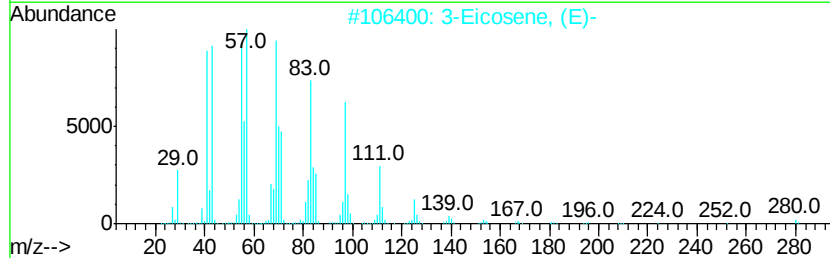
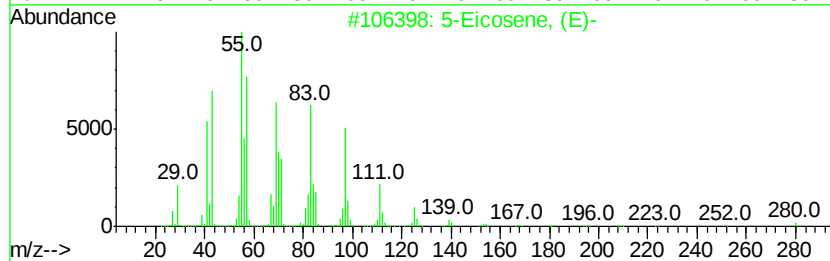
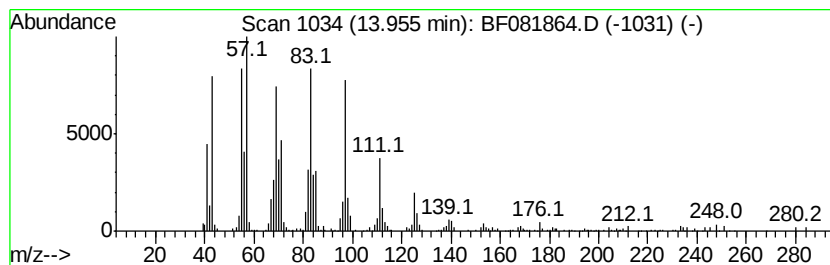
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 5-Eicosene, (E)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.96	6.91 ng	291691	Chrysene-d12	14.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Eicosene, (E)-	280	C20H40	074685-30-6	98
2		3-Eicosene, (E)-	280	C20H40	074685-33-9	95
3		1-Nonadecanol	284	C19H40O	001454-84-8	95
4		9-Eicosene, (E)-	280	C20H40	074685-29-3	94
5		Cyclopentadecane	210	C15H30	000295-48-7	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092815\
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 Sample : G3717-03
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-5-9-16-15

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.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...	5.11	10.5	ng	304622	1	6.93	580483	20.0
unknown6.70	6.70	42.3	ng	1227730	1	6.93	580483	20.0
2-Pyrimidinamine, ...	7.82	3.8	ng	162879	2	8.22	869547	20.0
Bicyclo[2.2.1]hep...	7.97	7.8	ng	340611	2	8.22	869547	20.0
unknown8.37	8.37	3.8	ng	164300	2	8.22	869547	20.0
Benzoic acid, 4-(...	9.41	4.8	ng	215813	3	9.98	892607	20.0
Eicosane	10.87	2.3	ng	116136	4	11.46	1000140	20.0
Heptadecane, 2,6,...	11.33	2.1	ng	104149	4	11.46	1000140	20.0
n-Hexadecanoic acid	12.00	24.1	ng	1205120	4	11.46	1000140	20.0
Cyclodecene, (Z)-	12.69	2.9	ng	146632	4	11.46	1000140	20.0
Octadecanoic acid	12.79	15.2	ng	643319	5	14.10	844731	20.0
Phenol, 4,4'-(1-m...	12.95	3.5	ng	148394	5	14.10	844731	20.0
Alpha-(1-(2-napht...	13.13	2.1	ng	88240	5	14.10	844731	20.0
5-Eicosene, (E)-	13.96	6.9	ng	291691	5	14.10	844731	20.0