

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092815\
 Data File : BF081868.D
 Acq On : 29 Sep 2015 6:03
 Operator : UM/IZ
 Sample : G3717-07
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PZ-8-9-16-15

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	267	rBV	357302	416138	11.53%	1.614%
2	5.520	294	296	305	rBV	919956	1126409	31.21%	4.369%
3	6.549	384	386	396	rBV	445117	692424	19.19%	2.686%
4	6.697	396	399	401	rBV	2134390	2263789	62.72%	8.781%
5	6.937	417	420	428	rBV	336101	586388	16.25%	2.275%
6	7.086	431	433	442	rBV	2340429	2413280	66.87%	9.361%
7	7.280	447	450	453	rVB	34062	38247	1.06%	0.148%
8	7.452	461	465	467	rBV2	41882	74354	2.06%	0.288%
9	7.497	467	469	472	rBV	1495738	1728364	47.89%	6.704%
10	7.634	479	481	487	rVB2	35132	75749	2.10%	0.294%
11	7.749	489	491	493	rBV	41745	40010	1.11%	0.155%
12	7.794	493	495	498	rBV	57973	62339	1.73%	0.242%
13	7.966	506	510	513	rBV5	17371	46679	1.29%	0.181%
14	8.126	521	524	525	rVV2	46248	65674	1.82%	0.255%
15	8.149	525	526	528	rVB	39921	48370	1.34%	0.188%
16	8.217	530	532	539	rBV	658342	788283	21.84%	3.058%
17	8.583	562	564	568	rVB3	29031	42071	1.17%	0.163%
18	8.652	568	570	573	rBV3	32268	53672	1.49%	0.208%
19	8.697	573	574	577	rBV2	48709	40682	1.13%	0.158%
20	8.755	577	579	580	rBV	108047	89996	2.49%	0.349%
21	8.857	584	588	589	rBV2	25429	63253	1.75%	0.245%
22	8.949	592	596	600	rBV5	46259	152758	4.23%	0.593%
23	9.063	604	606	609	rVB3	27353	46588	1.29%	0.181%
24	9.166	611	615	616	rBV3	25504	42441	1.18%	0.165%
25	9.257	621	623	624	rVV2	88933	105324	2.92%	0.409%
26	9.303	624	627	629	rVV	2605086	3609123	100.00%	13.999%
27	9.383	633	634	639	rVV5	56142	111084	3.08%	0.431%
28	9.463	639	641	642	rVV	43644	73018	2.02%	0.283%
29	9.486	642	643	645	rVV	91991	77242	2.14%	0.300%
30	9.520	645	646	650	rVB2	30182	41445	1.15%	0.161%
31	9.589	650	652	654	rBV2	36083	62284	1.73%	0.242%
32	9.635	654	656	658	rVV	79196	88556	2.45%	0.343%
33	9.692	658	661	663	rVB2	64235	95172	2.64%	0.369%
34	9.783	667	669	671	rBV2	40176	68654	1.90%	0.266%

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 Stop Thrs : 0 Peak Location: TOP

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35	9.829	671	673	676	rVB3	19069	39848	1.10%	0.155%
36	9.978	683	686	688	rVV	929333	939882	26.04%	3.646%
37	10.023	688	690	693	rVV3	61281	125257	3.47%	0.486%
38	10.115	696	698	699	rBV	31234	37371	1.04%	0.145%
39	10.229	706	708	710	rBV2	47876	100669	2.79%	0.390%
40	10.275	710	712	713	rVV2	31977	53647	1.49%	0.208%
41	10.309	713	715	718	rVB3	43767	94276	2.61%	0.366%
42	10.526	732	734	735	rBV2	36071	46458	1.29%	0.180%
43	10.652	743	745	747	rVB2	34672	45076	1.25%	0.175%
44	10.709	748	750	752	rBV2	55204	93806	2.60%	0.364%
45	10.766	752	755	763	rVB	1827985	2231272	61.82%	8.655%
46	10.972	771	773	775	rBV2	34186	58978	1.63%	0.229%
47	11.338	801	805	807	rBV3	71148	139097	3.85%	0.540%
48	11.463	814	816	819	rBV	824317	876802	24.29%	3.401%
49	11.806	845	846	851	rVB	48672	74453	2.06%	0.289%
50	11.989	860	862	866	rBV2	79463	106566	2.95%	0.413%
51	12.778	929	931	936	rVB4	29846	43208	1.20%	0.168%
52	13.052	952	955	958	rBV	2870554	3470146	96.15%	13.460%
53	13.647	1005	1007	1014	rVB	113398	137068	3.80%	0.532%
54	13.944	1031	1033	1040	rVB2	28999	48028	1.33%	0.186%
55	14.104	1045	1047	1050	rBV	788882	926720	25.68%	3.595%
56	15.578	1173	1176	1186	rVB	535509	762361	21.12%	2.957%

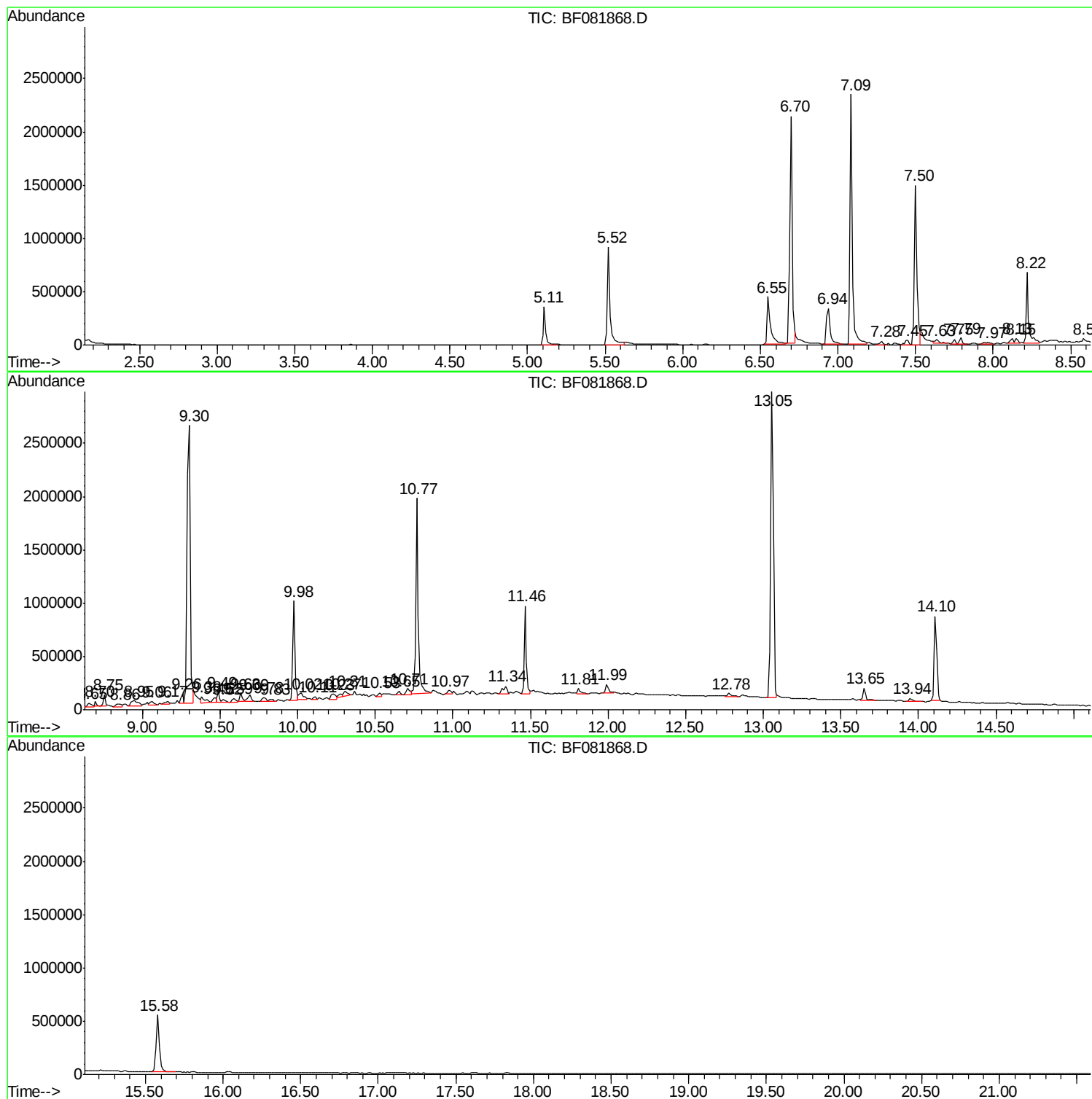
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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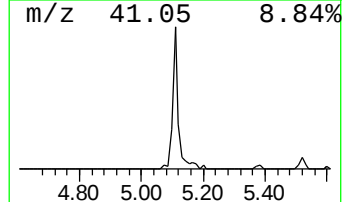
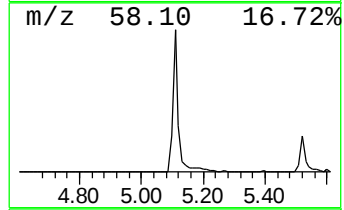
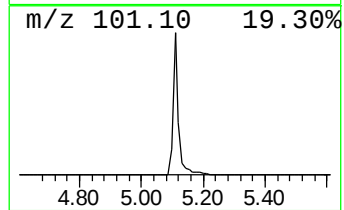
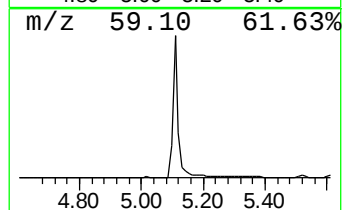
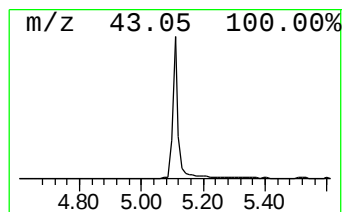
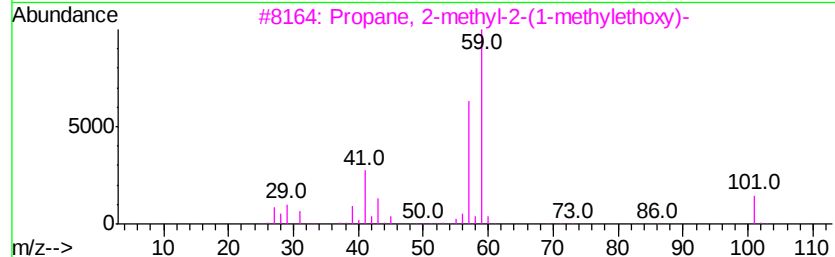
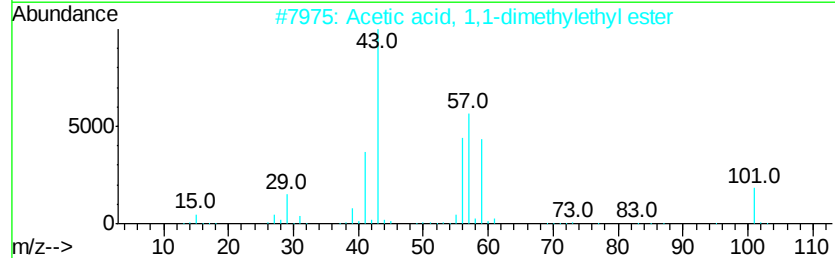
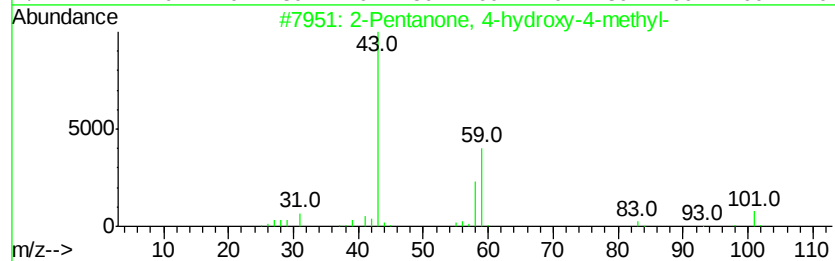
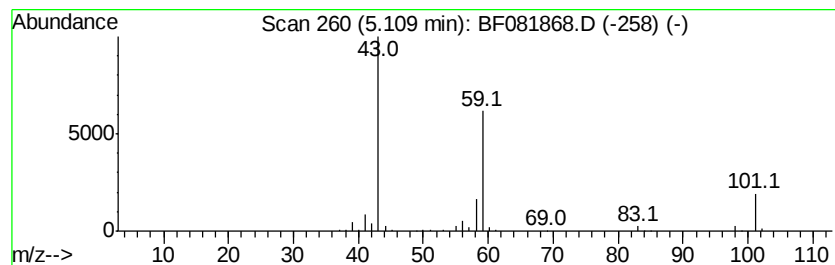
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	14.19 ng	416138	1,4-Dichlorobenzene-d4	6.94

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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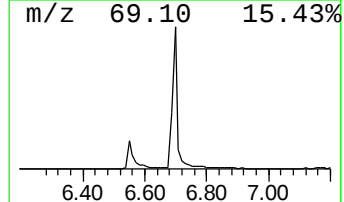
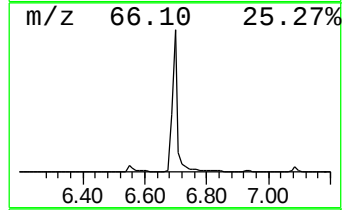
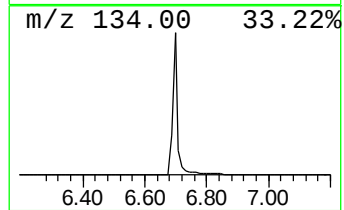
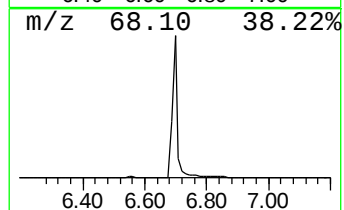
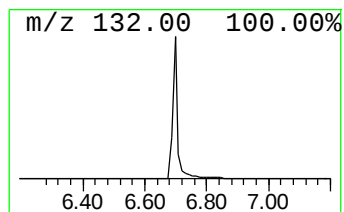
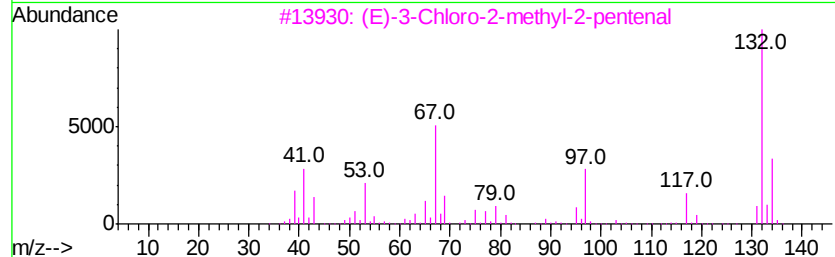
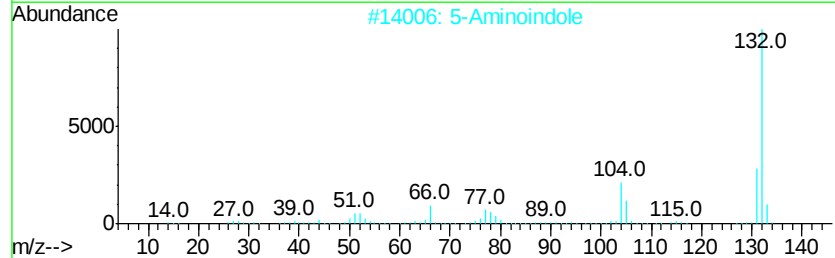
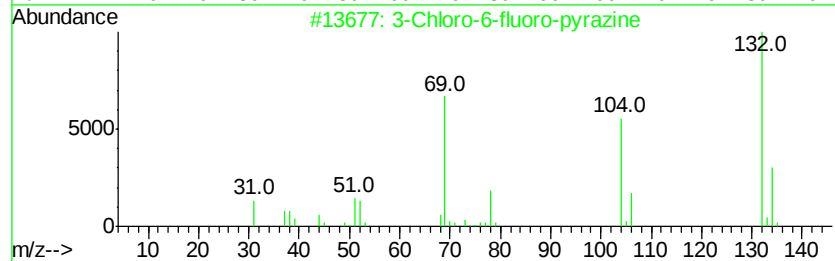
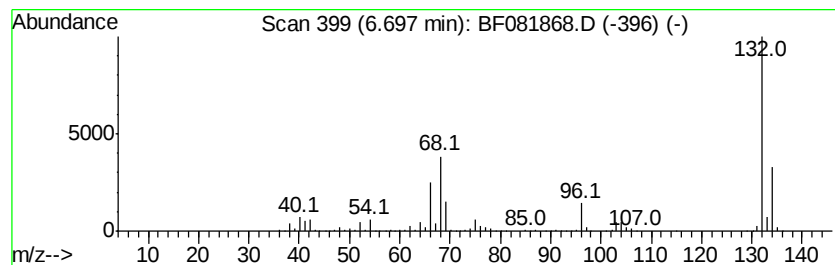
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.70 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.70	77.21 ng	2263790	1,4-Dichlorobenzene-d4	6.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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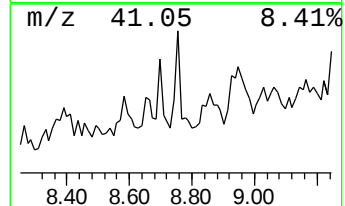
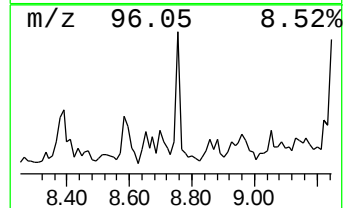
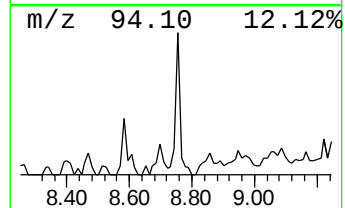
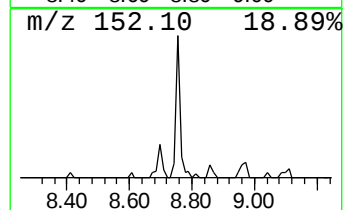
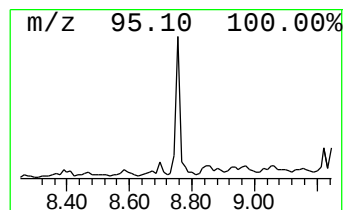
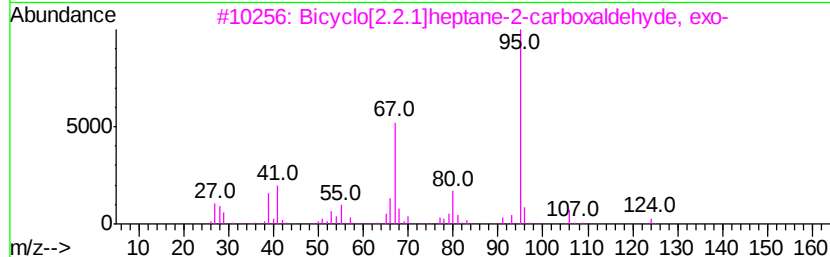
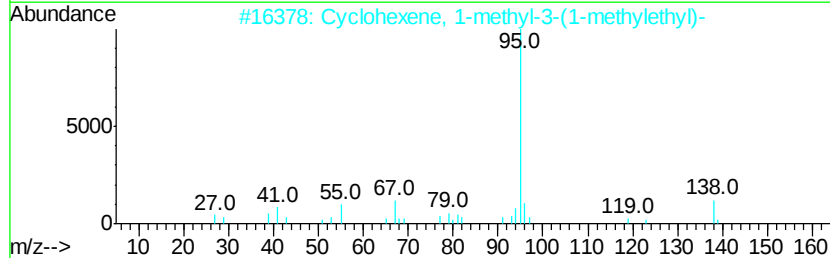
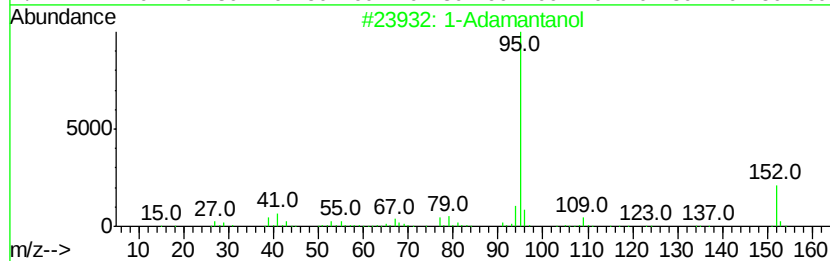
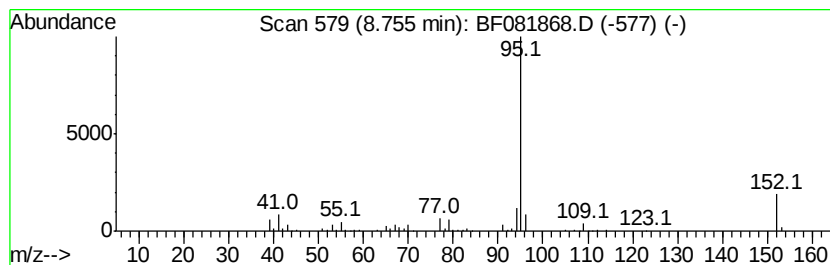
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1-Adamantanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.75	2.28 ng	89996	Naphthalene-d8	8.22

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Adamantanol		152	C10H16O	000768-95-6	91
2	Cyclohexene, 1-methyl-3-(1-methy...		138	C10H18	013828-31-4	39
3	Bicyclo[2.2.1]heptane-2-carboxal...		124	C8H12O	003574-55-8	38
4	1,3-Dimethyl-1-cyclohexene		110	C8H14	002808-76-6	38
5	5,7-Octadien-4-one, 2,6-dimethyl...		152	C10H16O	006752-80-3	38



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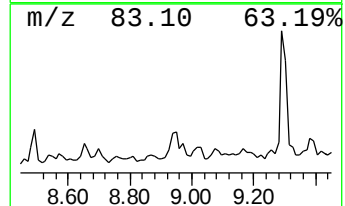
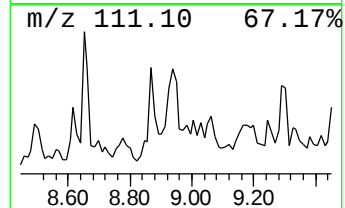
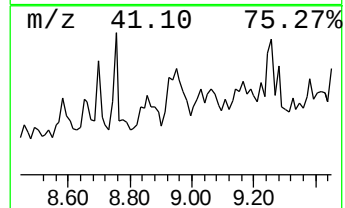
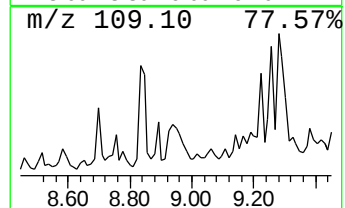
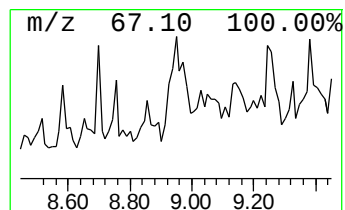
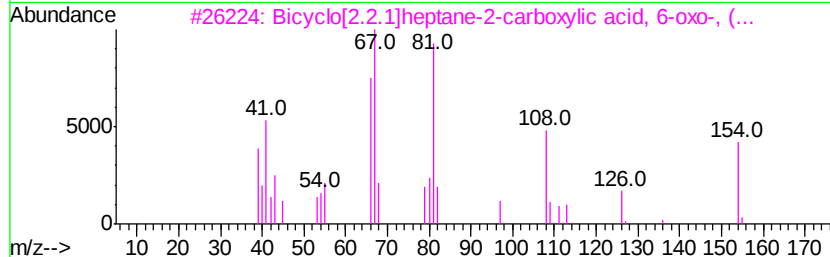
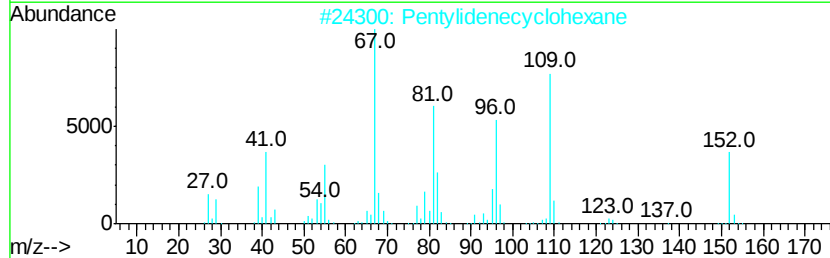
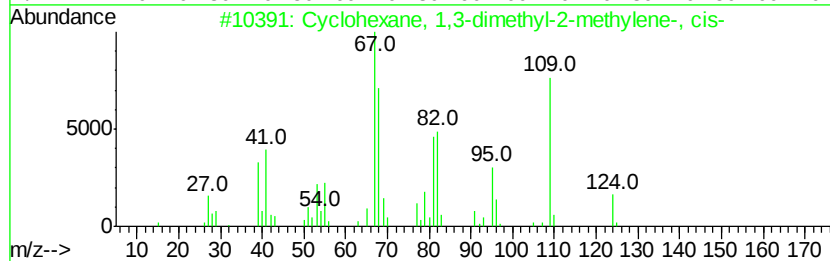
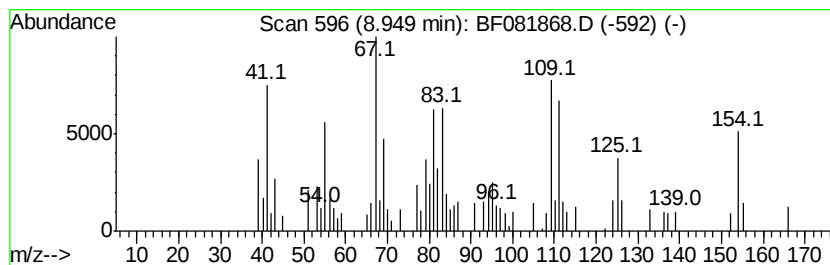
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 unknown8.95 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.95	3.88 ng	152758	Naphthalene-d8	8.22

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane, 1,3-dimethyl-2-meth...	124	C9H16	019781-47-6	38
2		Pentylidenecyclohexane	152	C11H20	039546-79-7	38
3		Bicyclo[2.2.1]heptane-2-carboxyl...	154	C8H10O3	1000141-75-5	30
4		1(2H)-Naphthalenone, octahydro-8...	166	C11H18O	000770-62-7	30
5		Cyclohexane, 1,3-dimethyl-2-meth...	124	C9H16	020348-74-7	25



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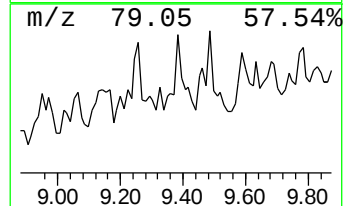
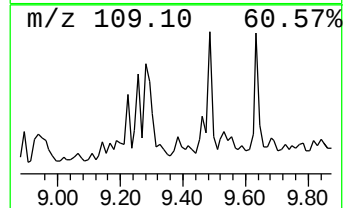
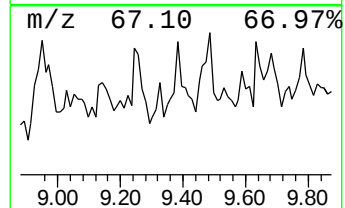
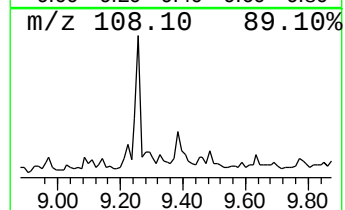
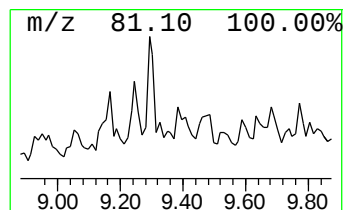
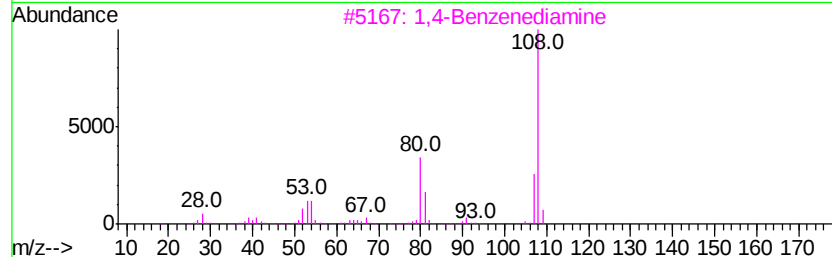
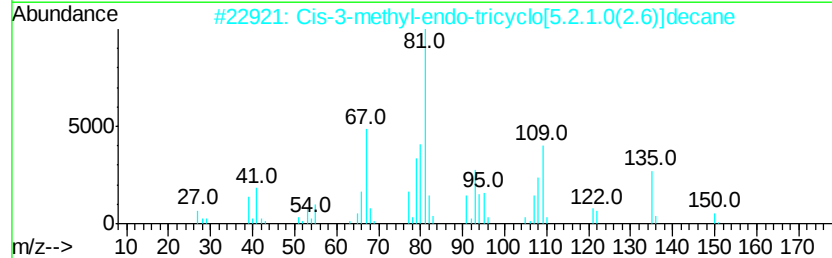
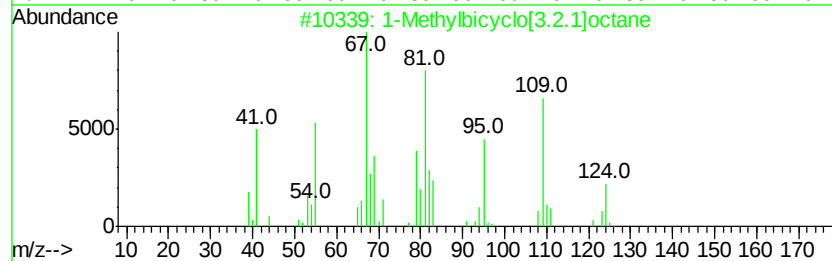
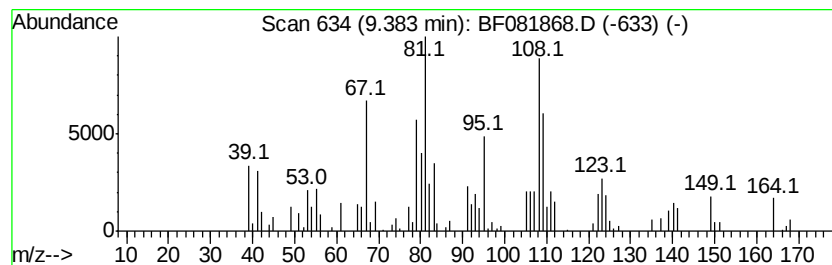
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown9.38 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.38	2.36 ng	111084	Acenaphthene-d10	9.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Methylbicyclo[3.2.1]octane	124	C9H16	119972-41-7	46
2		Cis-3-methyl-endo-tricyclo[5.2.1...]	150	C11H18	1000215-29-0	35
3		1,4-Benzenediamine	108	C6H8N2	000106-50-3	27
4		1H-Pyrrole-2-carboxaldehyde, 5-m...	109	C6H7NO	001192-79-6	27
5		Cyclopentanecarboxylic acid, 2-m...	140	C8H12O2	110550-98-6	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092815\
Data File : BF081868.D
Acq On : 29 Sep 2015 6:03
Operator : UM/IZ
Sample : G3717-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

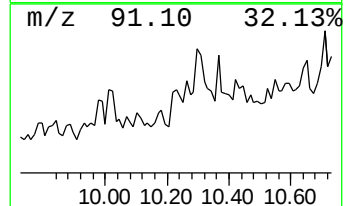
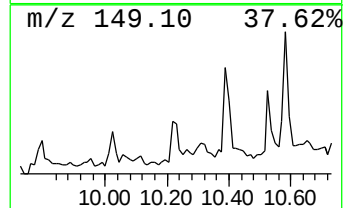
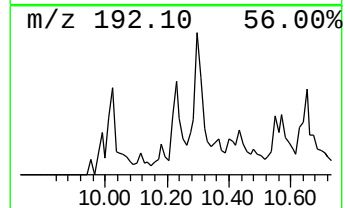
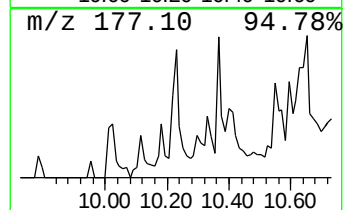
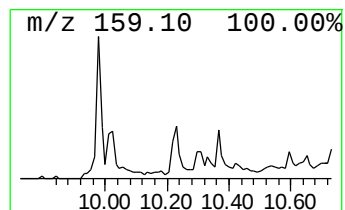
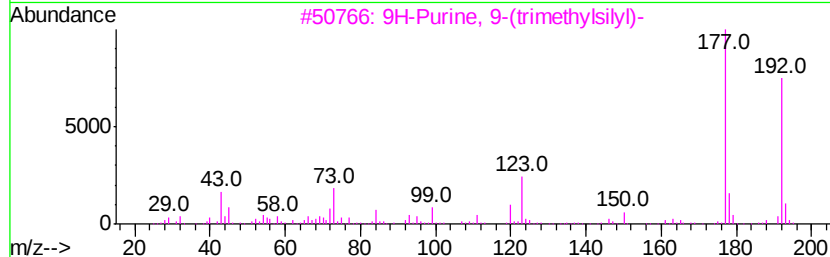
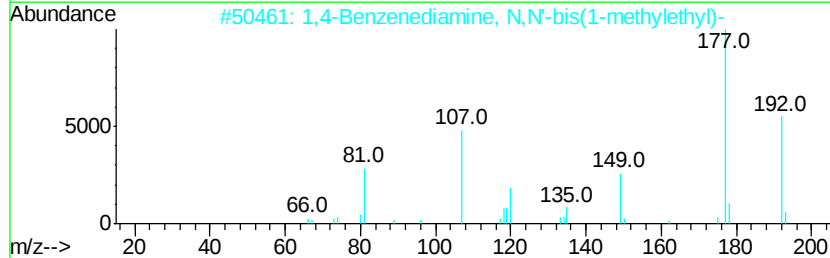
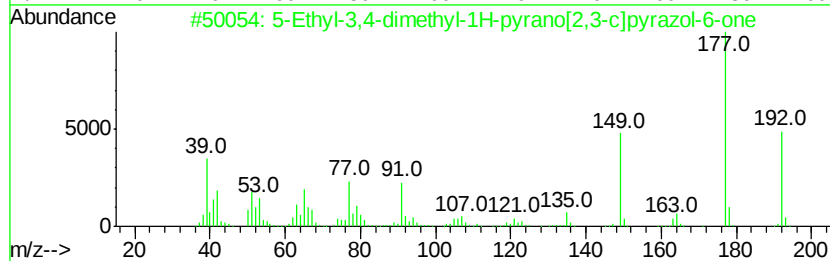
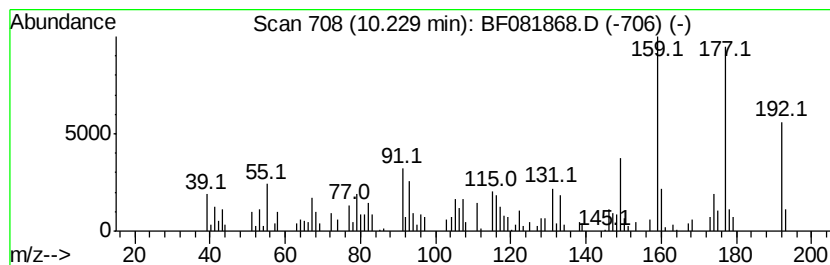
Instrument :
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ClientSampleId :
PZ-8-9-16-15

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

Peak Number 8 unknown10.23 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.23	2.14 ng	100669	Acenaphthene-d10	9.98	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Ethyl-3,4-dimethyl-1H-pyrano[2...	192	C10H12N2O2	1000296-60-7	45
2	1,4-Benzenediamine, N,N'-bis(1-m...	192	C12H20N2	004251-01-8	43
3	9H-Purine, 9-(trimethylsilyl)-	192	C8H12N4Si	032865-85-3	41
4	2,6-Diisopropylanisole	192	C13H20O	002944-52-7	35
5	1,3,2-Dioxarsenane, 2,5,5-trimet...	192	C6H13AsO2	042541-34-4	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092815\
Data File : BF081868.D
Acq On : 29 Sep 2015 6:03
Operator : UM/IZ
Sample : G3717-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

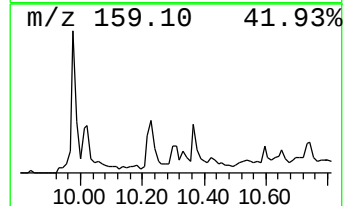
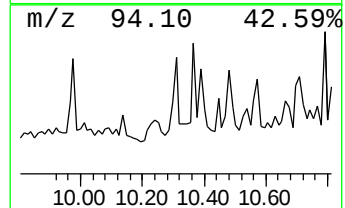
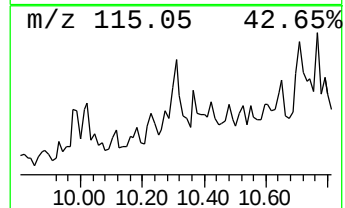
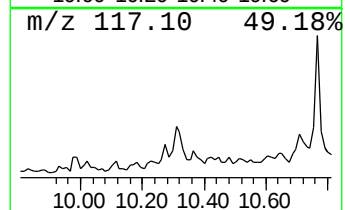
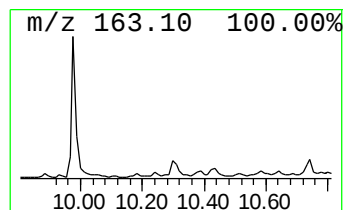
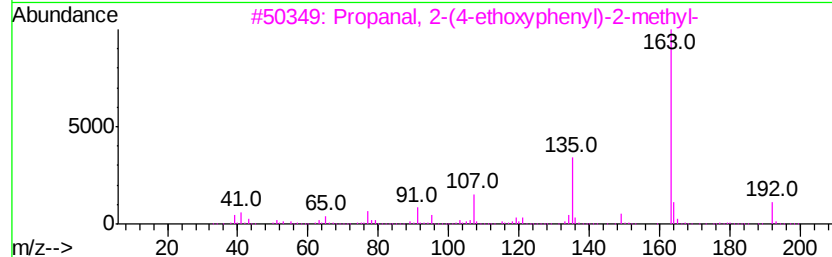
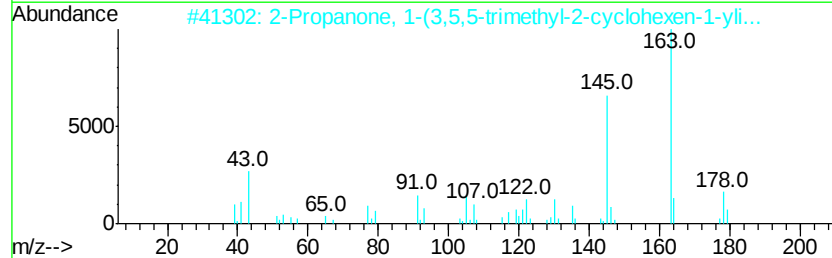
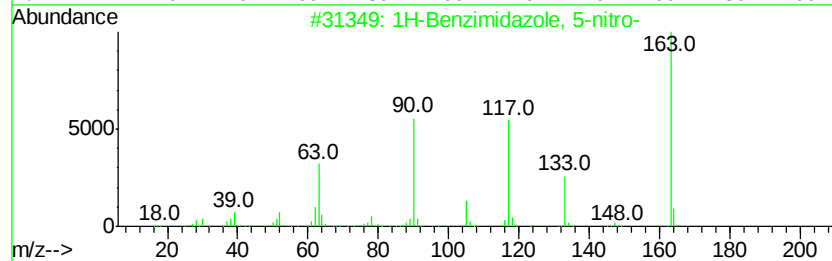
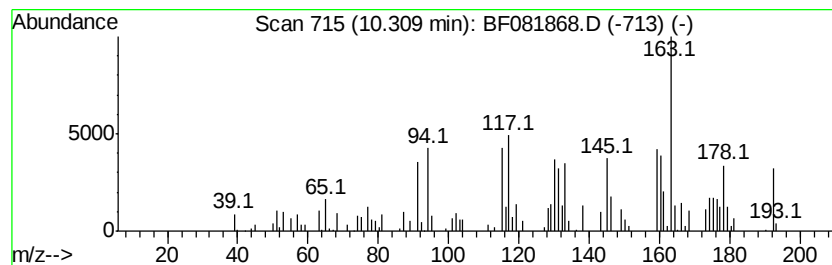
Instrument :
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ClientSampleId :
PZ-8-9-16-15

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

Peak Number 9 unknown10.31 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.31	2.01 ng	94276	Acenaphthene-d10		9.98
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzimidazole, 5-nitro-	163	C7H5N3O2	000094-52-0	27
2	2-Propanone, 1-(3,5,5-trimethyl-...	178	C12H18O	016695-72-0	22
3	Propanal, 2-(4-ethoxyphenyl)-2-m...	192	C12H16O2	093622-71-0	18
4	1,3-Disilaindan, 1,3-dimethyl-	178	C9H14Si2	017864-73-2	16
5	Benzene, 1-(1,1-dimethylethyl)-4...	178	C12H18O	017269-94-2	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092815\
Data File : BF081868.D
Acq On : 29 Sep 2015 6:03
Operator : UM/IZ
Sample : G3717-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PZ-8-9-16-15

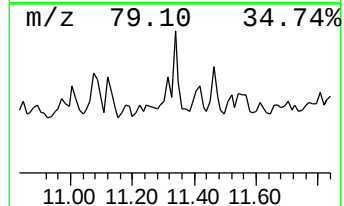
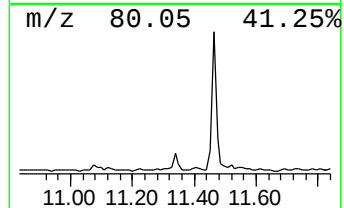
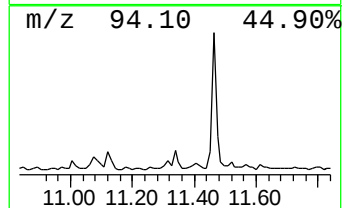
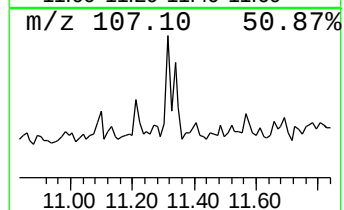
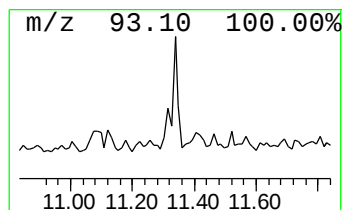
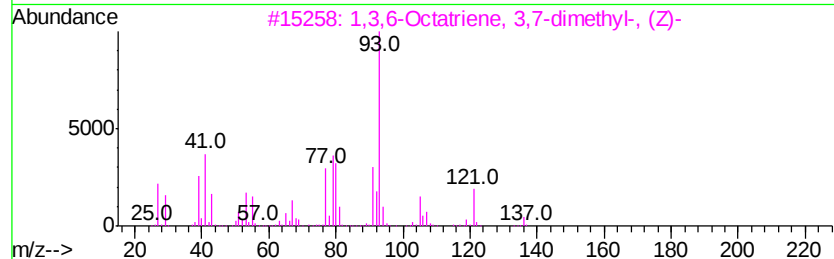
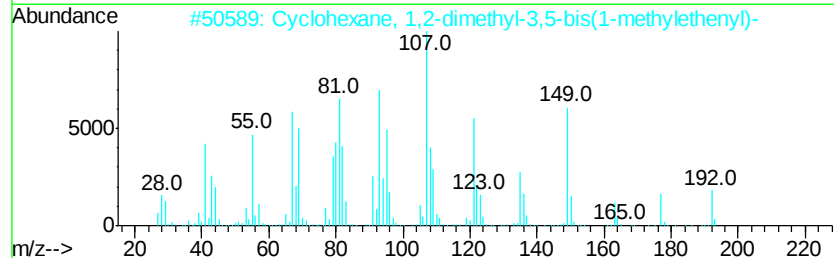
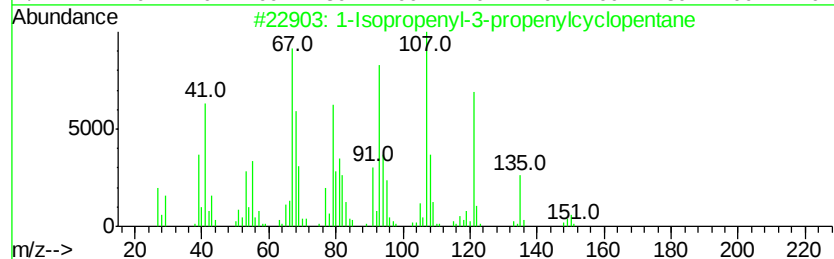
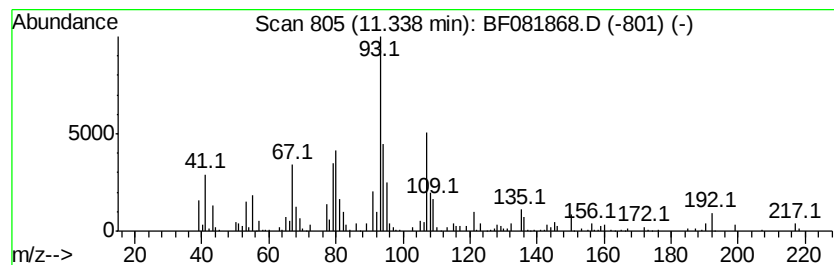
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Isopropenyl-3-propenylcyc... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.34	3.17 ng	139097	Phenanthrene-d10	11.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Isopropenyl-3-propenylcyclopent...	150	C11H18	1000192-81-2	52
2		Cyclohexane, 1,2-dimethyl-3,5-bi...	192	C14H24	062337-99-9	50
3		1,3,6-Octatriene, 3,7-dimethyl-,...	136	C10H16	003338-55-4	43
4		p-Menth-8-ene, 3-methylene-	150	C11H18	1000151-25-7	38
5		Spiro[4.4]non-1-ene	122	C9H14	000873-12-1	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092815\
Data File : BF081868.D
Acq On : 29 Sep 2015 6:03
Operator : UM/IZ
Sample : G3717-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

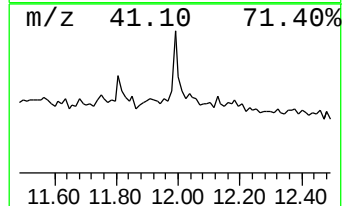
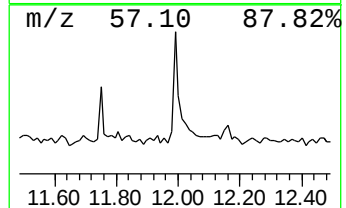
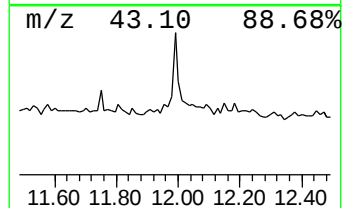
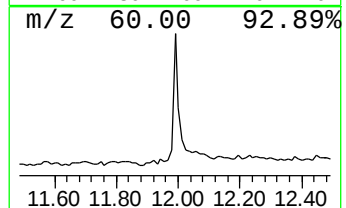
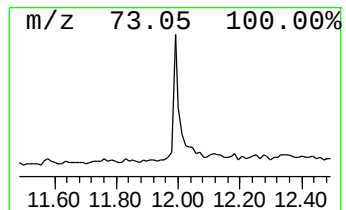
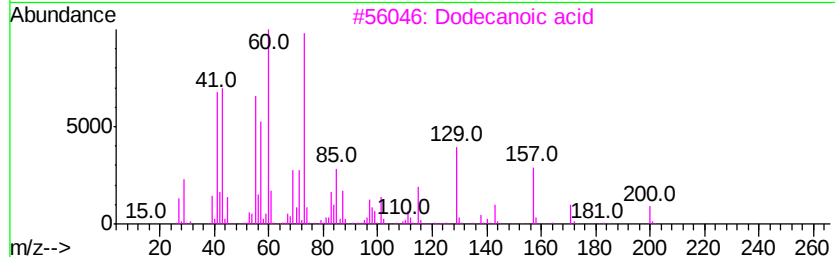
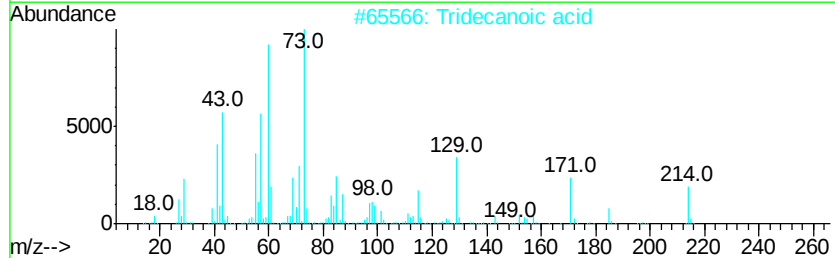
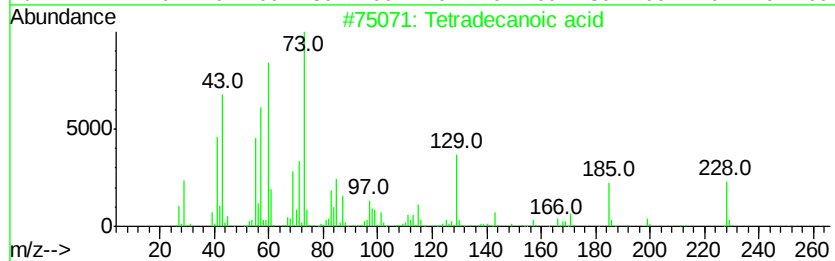
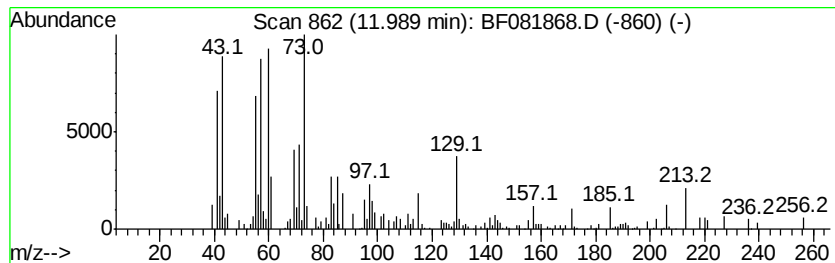
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Tetradecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.99	2.43 ng	106566	Phenanthrene-d10	11.46	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
2	Tridecanoic acid	214	C13H26O2	000638-53-9	64
3	Dodecanoic acid	200	C12H24O2	000143-07-7	58
4	n-Decanoic acid	172	C10H20O2	000334-48-5	52
5	Nonanoic acid	158	C9H18O2	000112-05-0	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092815\
Data File : BF081868.D
Acq On : 29 Sep 2015 6:03
Operator : UM/IZ
Sample : G3717-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

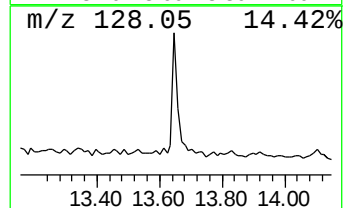
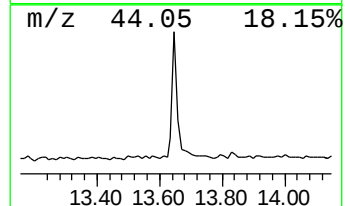
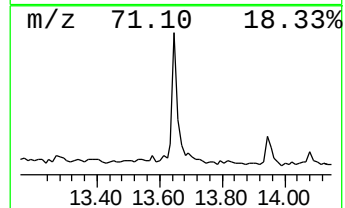
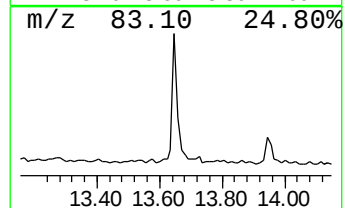
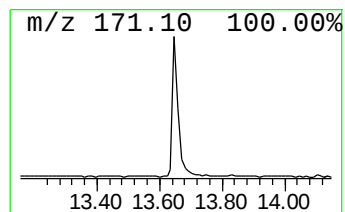
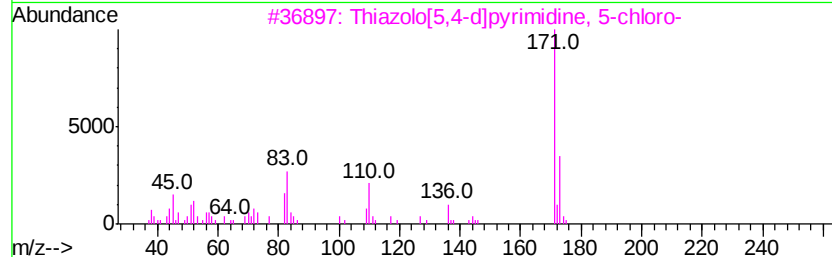
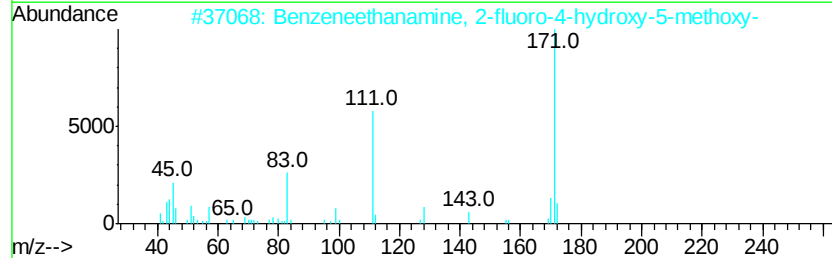
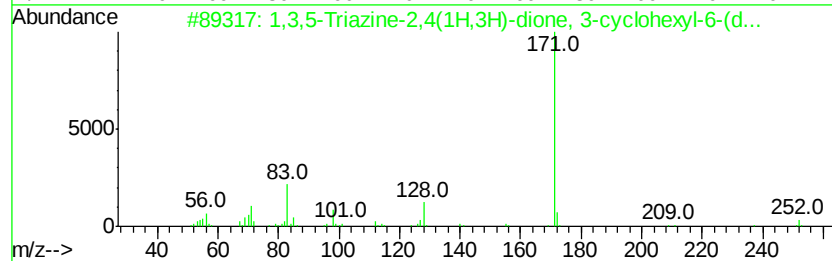
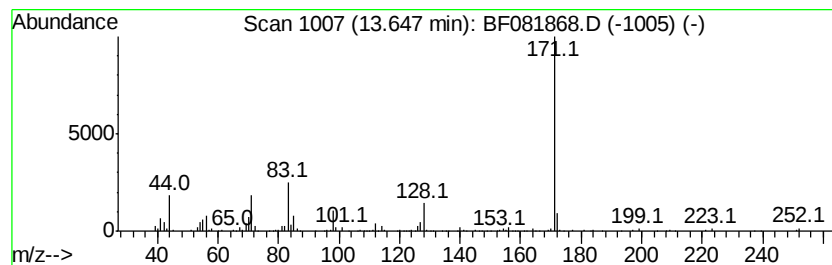
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1,3,5-Triazine-2,4(1H,3H)-d... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.65	2.96 ng	137068	Chrysene-d12	14.10
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1,3,5-Triazine-2,4(1H,3H)-dione,...	252 C12H20N4O2	051235-04-2	96
2	Benzeneethanamine, 2-fluoro-4-hy...	171 C8H10FN02	1000116-43-3	72
3	Thiazolo[5,4-d]pyrimidine, 5-chl...	171 C5H2ClN3S	013316-08-0	64
4	3-Cyanoquinoxaline 1-oxide	171 C9H5N3O	053688-02-1	50
5	4-Pyrrolylbenzaldehyde	171 C11H9NO	023351-05-5	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092815\
Data File : BF081868.D
Acq On : 29 Sep 2015 6:03
Operator : UM/IZ
Sample : G3717-07
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PZ-8-9-16-15

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	14.2	ng	416138	1	6.94	586388	20.0
unknown6.70	6.70	77.2	ng	2263790	1	6.94	586388	20.0
1-Adamantanol	8.75	2.3	ng	89996	2	8.22	788283	20.0
unknown8.95	8.95	3.9	ng	152758	2	8.22	788283	20.0
unknown9.38	9.38	2.4	ng	111084	3	9.98	939882	20.0
unknown10.23	10.23	2.1	ng	100669	3	9.98	939882	20.0
unknown10.31	10.31	2.0	ng	94276	3	9.98	939882	20.0
1-Isopropenyl-3-p...	11.34	3.2	ng	139097	4	11.46	876802	20.0
Tetradecanoic acid	11.99	2.4	ng	106566	4	11.46	876802	20.0
1,3,5-Triazine-2,...	13.65	3.0	ng	137068	5	14.10	926720	20.0