

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121415\
 Data File : BF083798.D
 Acq On : 15 Dec 2015 4:57
 Operator : UM/IZ
 Sample : G4779-13
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-38

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121315.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.082	260	262	266	rVB	217287	255721	3.95%	0.311%
2	5.504	296	299	301	rVB	1175821	1339823	20.71%	1.628%
3	6.064	345	348	350	rVB	79889	78726	1.22%	0.096%
4	6.522	386	388	390	rBV	1009274	876844	13.55%	1.065%
5	6.670	398	401	403	rBV	1921710	2373735	36.69%	2.884%
6	6.853	413	417	419	rBV3	33884	85178	1.32%	0.103%
7	6.899	419	421	423	rVB	801133	673266	10.41%	0.818%
8	7.059	432	435	436	rBV	2274217	2369018	36.62%	2.878%
9	7.150	441	443	446	rVB	403038	416338	6.44%	0.506%
10	7.242	448	451	453	rVB	231203	396274	6.13%	0.481%
11	7.299	453	456	458	rBV2	65578	83950	1.30%	0.102%
12	7.459	466	470	474	rBV3	1572274	3363808	51.99%	4.087%
13	7.550	476	478	480	rBV2	185905	202230	3.13%	0.246%
14	7.596	480	482	483	rVB	168767	146523	2.26%	0.178%
15	7.676	485	489	491	rVB	893570	895921	13.85%	1.089%
16	7.710	491	492	495	rBV2	39137	75934	1.17%	0.092%
17	7.790	497	499	501	rBV	133329	141537	2.19%	0.172%
18	7.859	501	505	508	rBV3	413007	716633	11.08%	0.871%
19	7.928	508	511	513	rBV	1995293	1985447	30.69%	2.412%
20	7.962	513	514	516	rBV2	82960	75166	1.16%	0.091%
21	8.030	516	520	521	rBV	1053636	1218385	18.83%	1.480%
22	8.053	521	522	524	rBV	84467	101889	1.57%	0.124%
23	8.122	526	528	530	rBV2	246021	261455	4.04%	0.318%
24	8.179	530	533	534	rBV	1012729	1224425	18.93%	1.488%
25	8.282	540	542	543	rBV2	94116	120109	1.86%	0.146%
26	8.316	543	545	546	rVB2	89004	79822	1.23%	0.097%
27	8.339	546	547	548	rBV	98729	98114	1.52%	0.119%
28	8.385	548	551	553	rVB2	273063	341951	5.29%	0.415%
29	8.442	553	556	560	rBV3	318121	817955	12.64%	0.994%
30	8.522	560	563	566	rVV5	61082	150048	2.32%	0.182%
31	8.568	566	567	568	rVV	206004	148239	2.29%	0.180%
32	8.659	571	575	576	rBV	236457	403689	6.24%	0.490%
33	8.693	576	578	580	rBV2	543818	613604	9.48%	0.746%
34	8.750	580	583	584	rVV3	72098	129804	2.01%	0.158%

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35	8.773	584	585	587	rVB2	160475	135260	2.09%	0.164%
36	8.819	587	589	590	rBV2	92899	170401	2.63%	0.207%
37	8.842	590	591	593	rVB	564574	752272	11.63%	0.914%
38	8.899	593	596	598	rBV	552251	654263	10.11%	0.795%
39	9.002	601	605	607	rVV	2268120	2717219	42.00%	3.301%
40	9.036	607	608	610	rVB2	94740	104161	1.61%	0.127%
41	9.093	610	613	614	rBV3	203421	331489	5.12%	0.403%
42	9.128	614	616	619	rVB2	450576	479809	7.42%	0.583%
43	9.185	619	621	622	rBV	701850	641462	9.92%	0.779%
44	9.208	622	623	625	rVV2	84846	99551	1.54%	0.121%
45	9.265	625	628	629	rVV	3404238	3227897	49.89%	3.922%
46	9.299	629	631	633	rVV	961863	846693	13.09%	1.029%
47	9.379	633	638	640	rVV2	1432379	1570971	24.28%	1.909%
48	9.413	640	641	644	rVV3	121328	264968	4.10%	0.322%
49	9.471	644	646	648	rVV3	269235	539487	8.34%	0.655%
50	9.528	648	651	652	rVV2	396920	763559	11.80%	0.928%
51	9.608	652	658	662	rVV3	2535786	6469525	100.00%	7.861%
52	9.665	662	663	664	rVV	622658	680948	10.53%	0.827%
53	9.699	664	666	667	rVV	918882	1120922	17.33%	1.362%
54	9.722	667	668	670	rVV	1181935	1232084	19.04%	1.497%
55	9.791	672	674	676	rVV2	220396	297769	4.60%	0.362%
56	9.848	676	679	681	rVV	1206340	1290924	19.95%	1.568%
57	9.905	681	684	685	rVV	1518255	1980466	30.61%	2.406%
58	9.939	685	687	690	rVV	1965338	2296848	35.50%	2.791%
59	10.019	690	694	697	rVV3	1490061	2674376	41.34%	3.249%
60	10.076	697	699	700	rVV2	376585	734599	11.35%	0.893%
61	10.099	700	701	706	rVV3	989590	2602361	40.22%	3.162%
62	10.225	709	712	713	rVV3	465081	967143	14.95%	1.175%
63	10.248	713	714	717	rVV3	1034204	1556068	24.05%	1.891%
64	10.328	717	721	723	rVV4	970843	2155205	33.31%	2.619%
65	10.373	723	725	726	rVV2	311054	454406	7.02%	0.552%
66	10.442	726	731	733	rVV4	1254770	2958360	45.73%	3.594%
67	10.476	733	734	736	rVV2	841175	950130	14.69%	1.154%
68	10.511	736	737	739	rVV	705565	718320	11.10%	0.873%
69	10.579	739	743	745	rVV4	409898	1102557	17.04%	1.340%
70	10.659	745	750	751	rVV3	482798	1239681	19.16%	1.506%
71	10.682	751	752	754	rVV2	557987	691925	10.70%	0.841%

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72	10.728	754	756	758	rVV2	1829529	2563886	39.63%	3.115%
73	10.774	758	760	761	rVV2	497801	781517	12.08%	0.950%
74	10.796	761	762	765	rVV3	661413	1127142	17.42%	1.369%
75	10.842	765	766	770	rVV3	353633	618926	9.57%	0.752%
76	10.934	770	774	778	rVB4	299341	755042	11.67%	0.917%
77	11.082	784	787	789	rVV4	124347	221280	3.42%	0.269%
78	11.116	789	790	797	rVB	333904	523897	8.10%	0.637%
79	11.219	797	799	800	rBV2	94014	133151	2.06%	0.162%
80	11.299	804	806	807	rBV2	88548	78852	1.22%	0.096%
81	11.425	814	817	818	rBV	699554	844689	13.06%	1.026%
82	11.528	825	826	833	rVB2	107705	158394	2.45%	0.192%
83	11.631	833	835	838	rVV2	77319	149142	2.31%	0.181%
84	11.814	849	851	855	rVV2	41276	78499	1.21%	0.095%
85	12.785	934	936	938	rVV	72019	72344	1.12%	0.088%
86	13.002	952	955	958	rBV	2770669	3083834	47.67%	3.747%
87	14.042	1044	1046	1049	rBV	695598	782529	12.10%	0.951%
88	15.483	1169	1172	1175	rBV	493807	595106	9.20%	0.723%

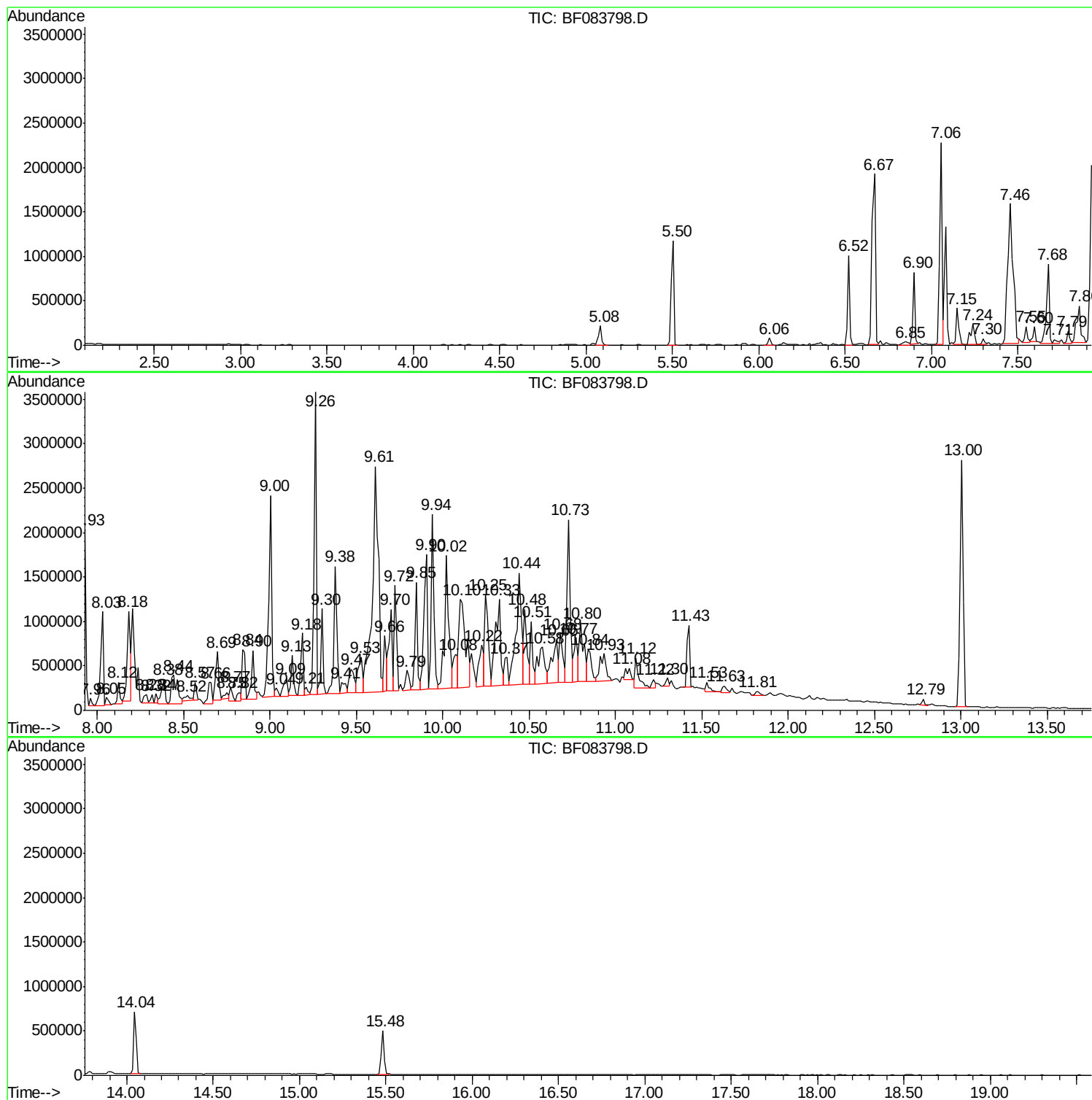
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W-38

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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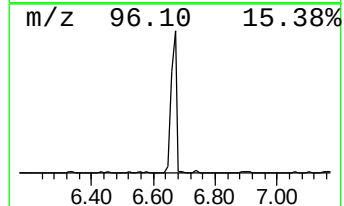
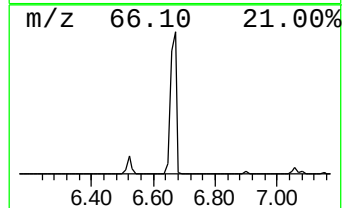
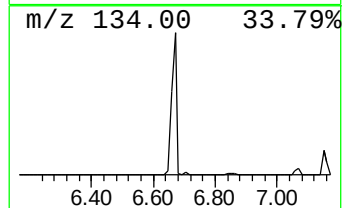
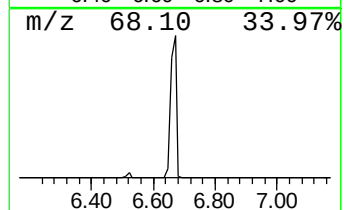
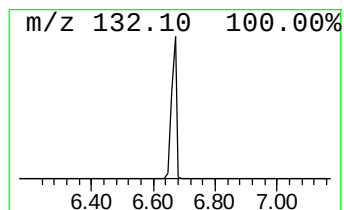
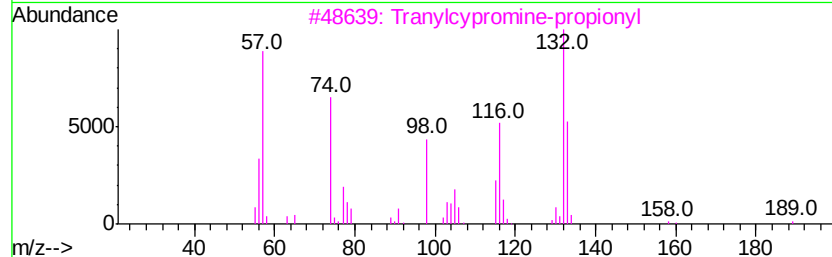
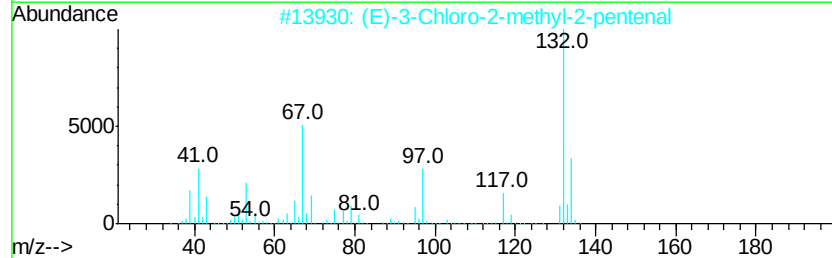
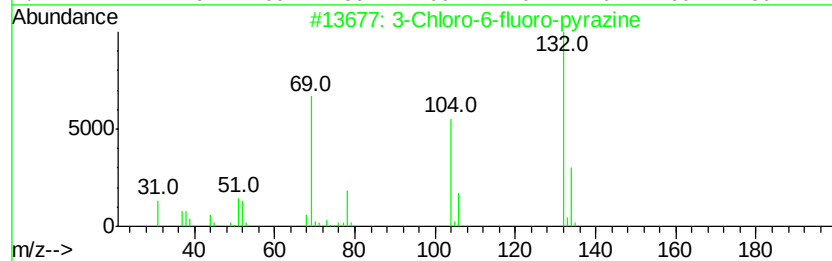
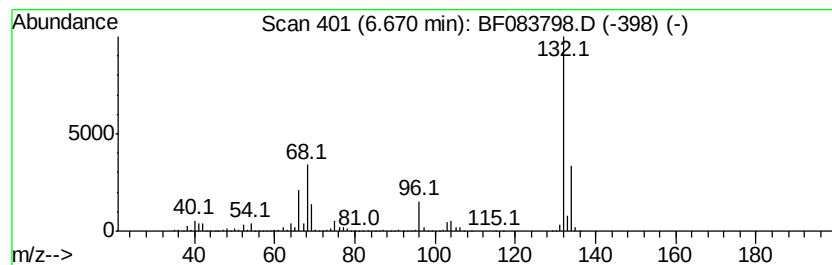
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121315.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	70.51 ng	2373740	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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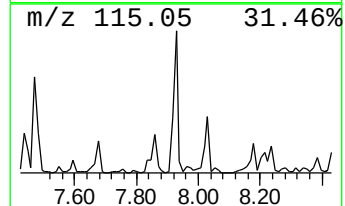
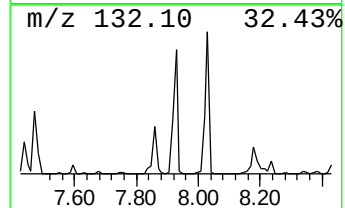
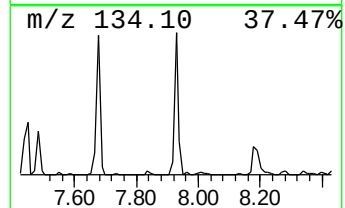
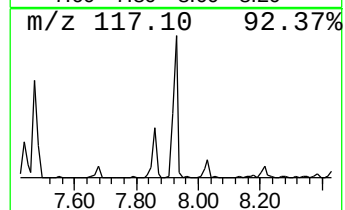
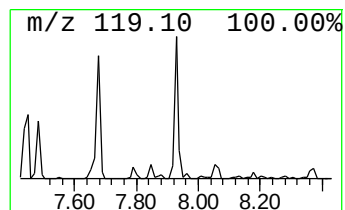
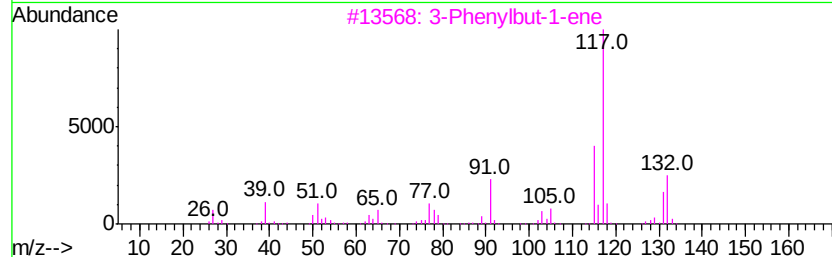
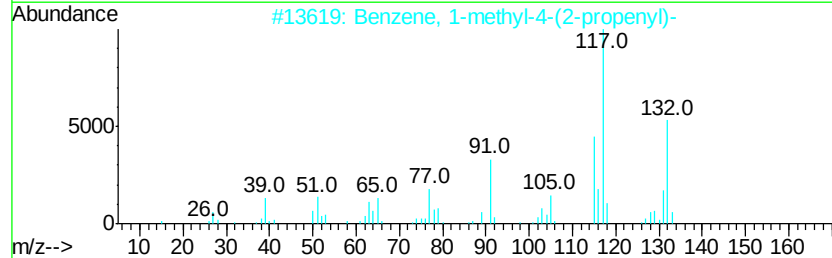
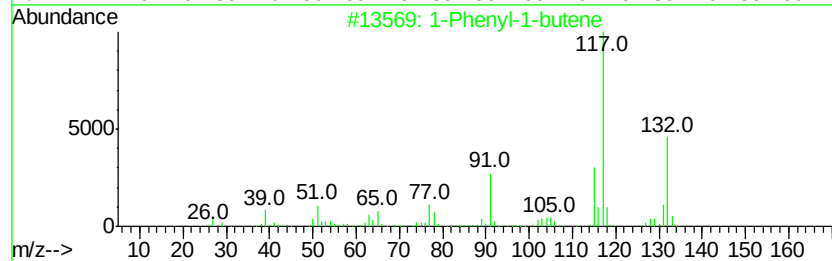
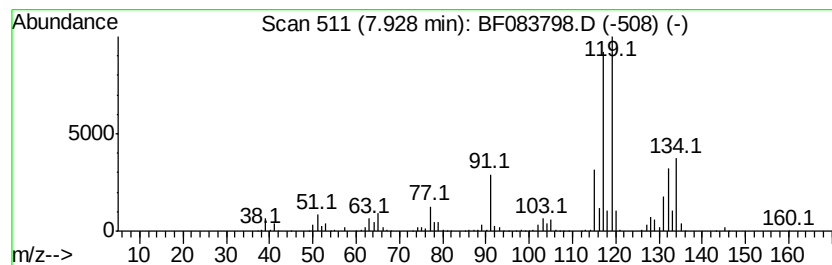
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Phenyl-1-butene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.93	32.43 ng	1985450	Naphthalene-d8	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	91
2		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	55
4		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	55
5		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	55



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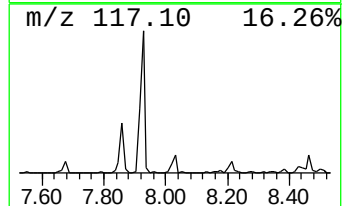
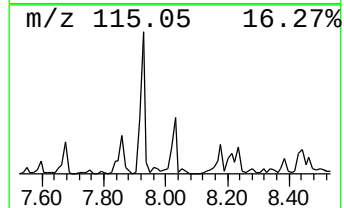
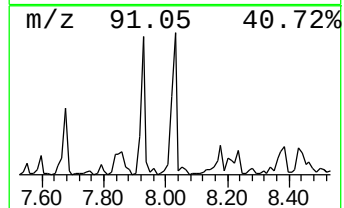
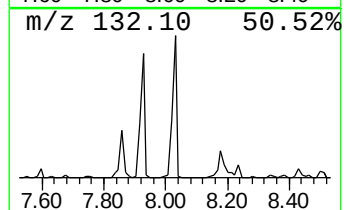
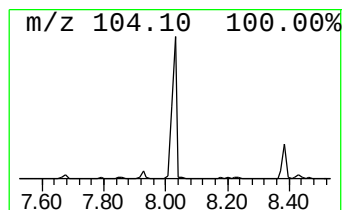
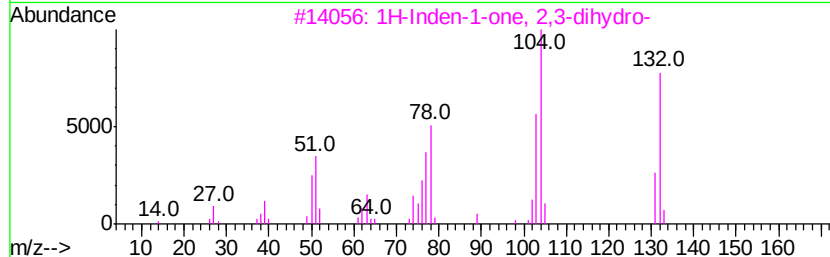
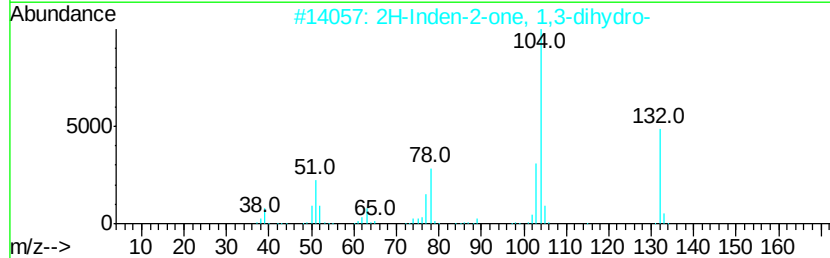
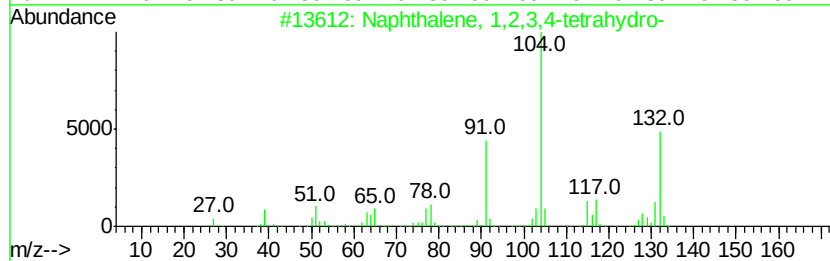
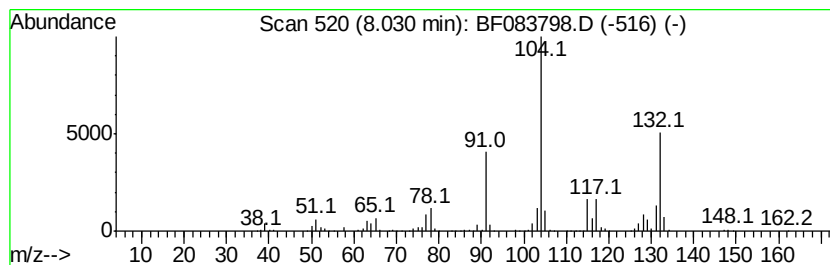
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.03	19.90 ng	1218390	Naphthalene-d8	8.18

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	97
2		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	53
3		1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	43
4		3(2H)-Furanone, dihydro-2,2-dime...	190	C12H14O2	063678-00-2	38
5		Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	38



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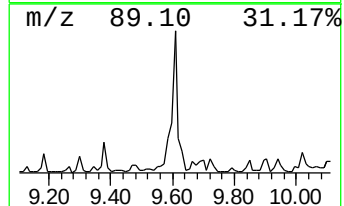
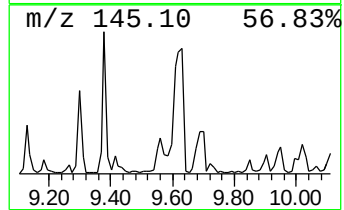
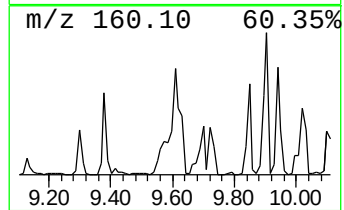
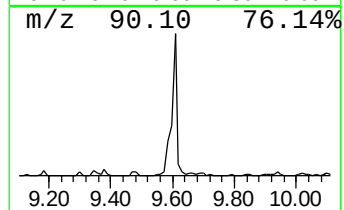
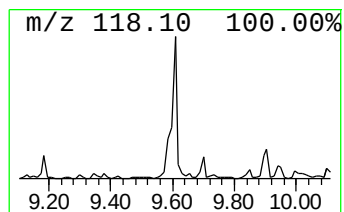
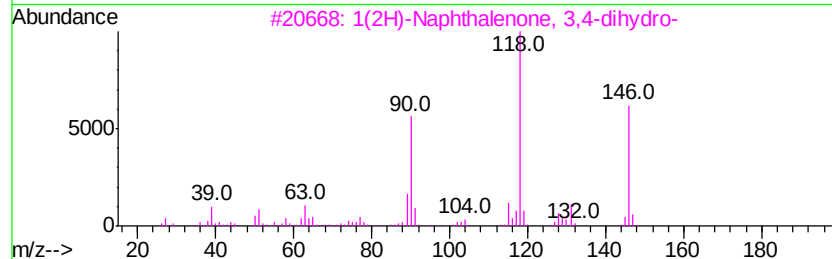
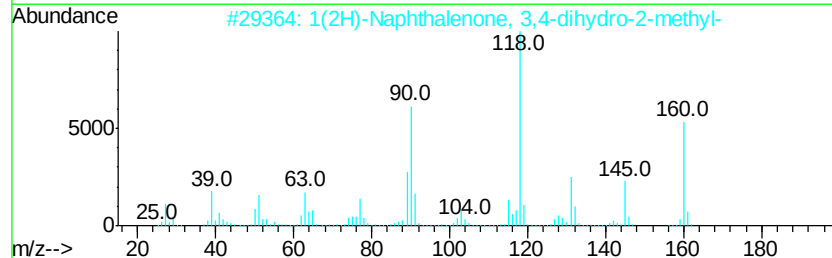
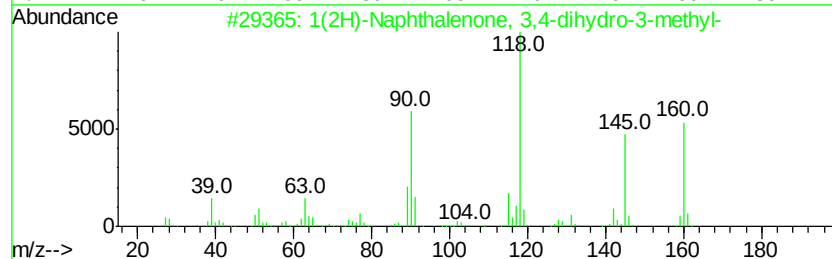
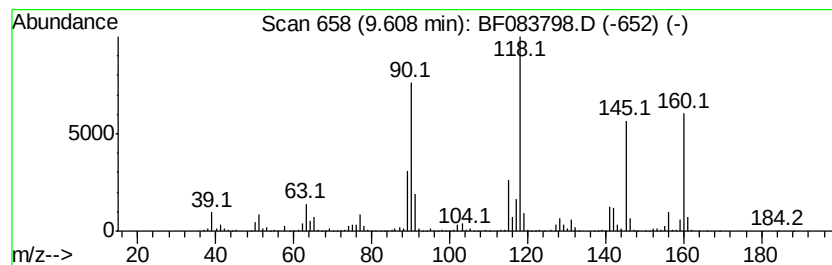
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 1(2H)-Naphthalenone, 3,4-di... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.61	56.33 ng	6469530	Acenaphthene-d10	9.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	014944-23-1	97
2		1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	001590-08-5	87
3		1(2H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000529-34-0	72
4		Homophthalimide	161	C9H7NO2	004456-77-3	68
5		1(2H)-Isoquinolinone	145	C9H7NO	000491-30-5	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF121415\
Data File : BF083798.D
Acq On : 15 Dec 2015 4:57
Operator : UM/IZ
Sample : G4779-13
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
W-38

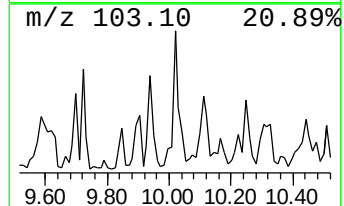
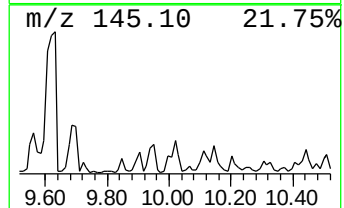
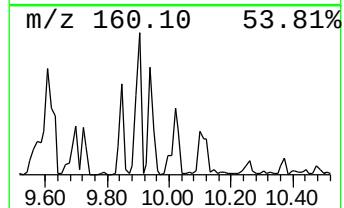
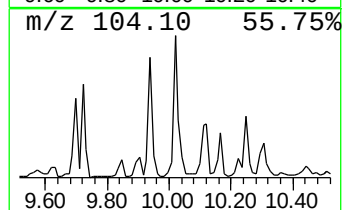
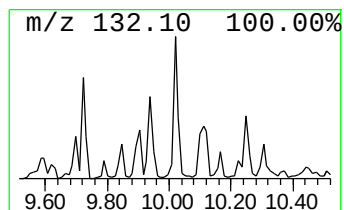
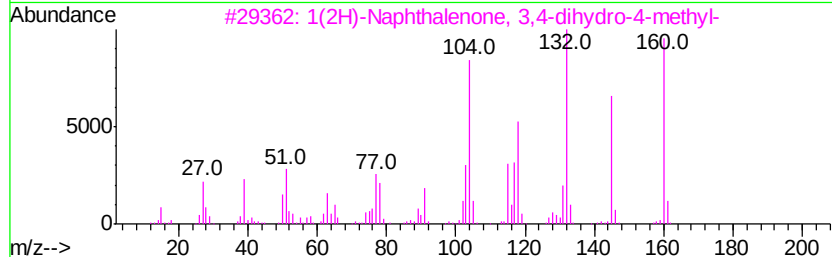
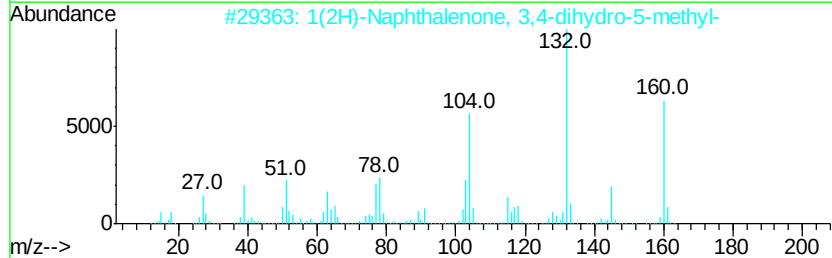
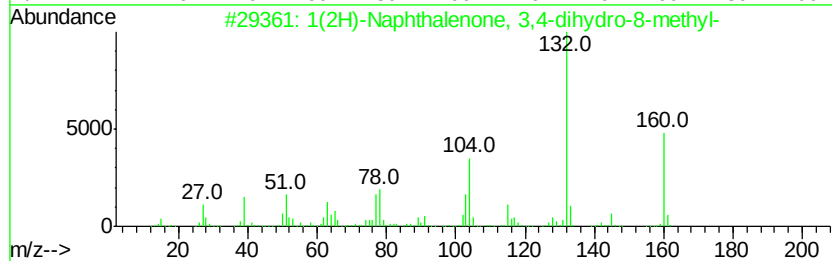
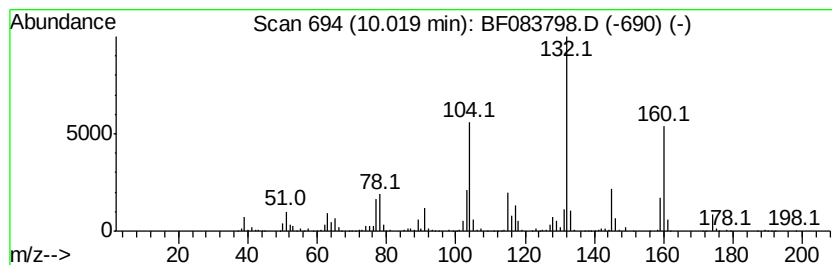
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1(2H)-Naphthalenone, 3,4-di... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.02	23.29 ng	2674380	Acenaphthene-d10	9.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	051015-28-2	92
2		1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	006939-35-1	90
3		1(2H)-Naphthalenone, 3,4-dihydro...	160	C11H12O	019832-98-5	58
4		Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	53
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121415\
Data File : BF083798.D
Acq On : 15 Dec 2015 4:57
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Sample : G4779-13
Misc :
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Instrument :
BNA_F
ClientSampleId :
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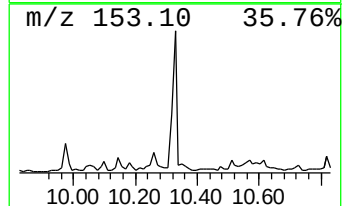
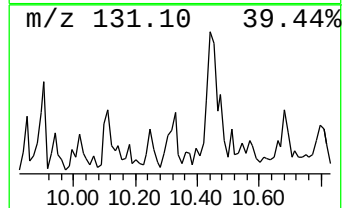
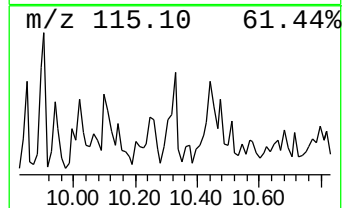
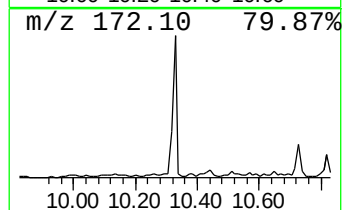
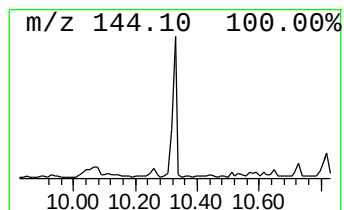
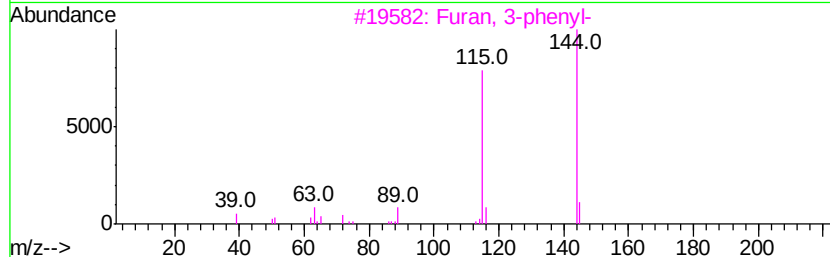
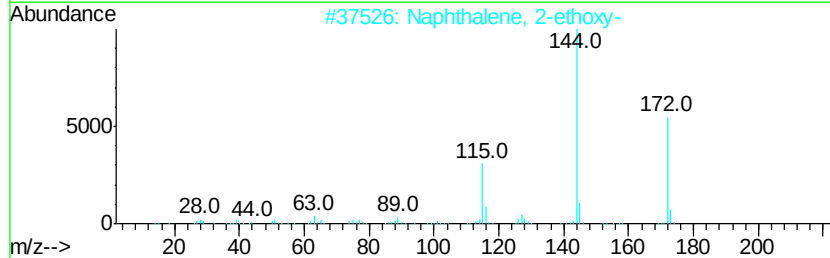
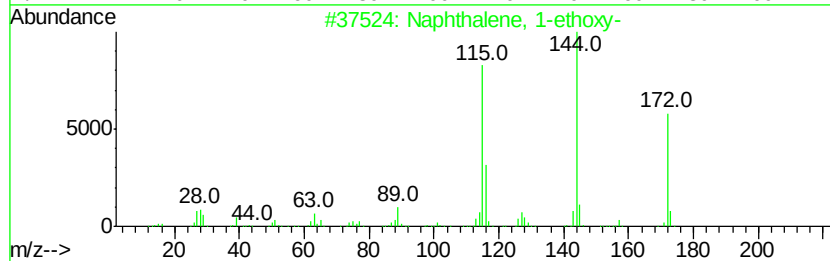
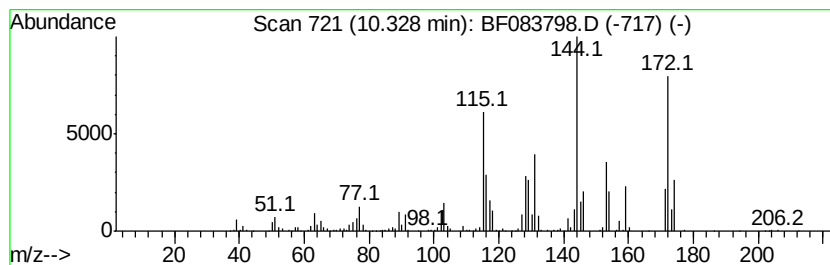
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown10.33 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.33	18.77 ng	2155210	Acenaphthene-d10	9.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-ethoxyv-	172	C12H12O	005328-01-8	46
2		Naphthalene, 2-ethoxyv-	172	C12H12O	000093-18-5	41
3		Furan, 3-phenyl-	144	C10H8O	013679-41-9	25
4		1-Naphthalenol	144	C10H8O	000090-15-3	25
5		Ethanone, 1-(1H-indol-3-yl)-	159	C10H9NO	000703-80-0	20



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121415\
Data File : BF083798.D
Acq On : 15 Dec 2015 4:57
Operator : UM/IZ
Sample : G4779-13
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
W-38

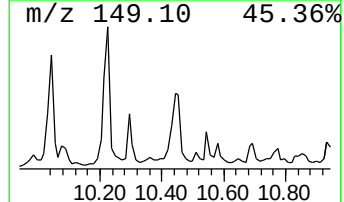
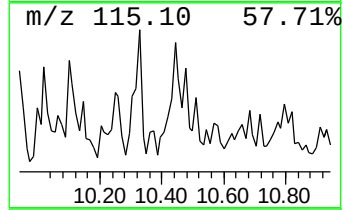
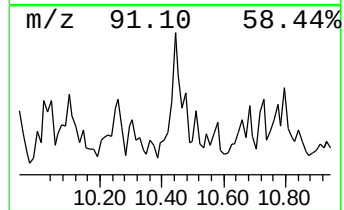
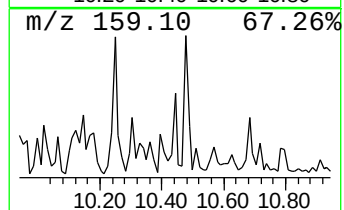
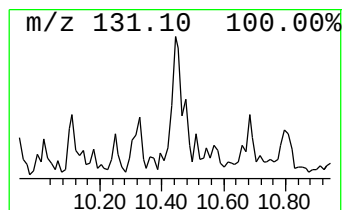
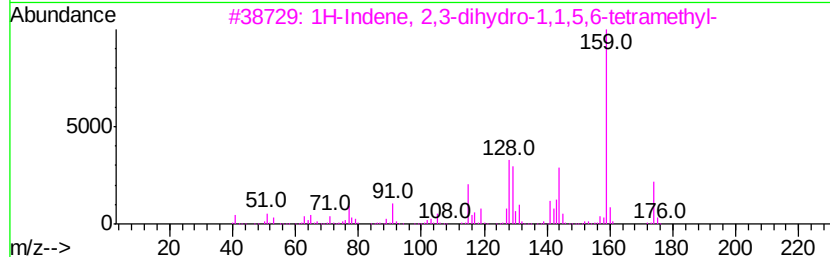
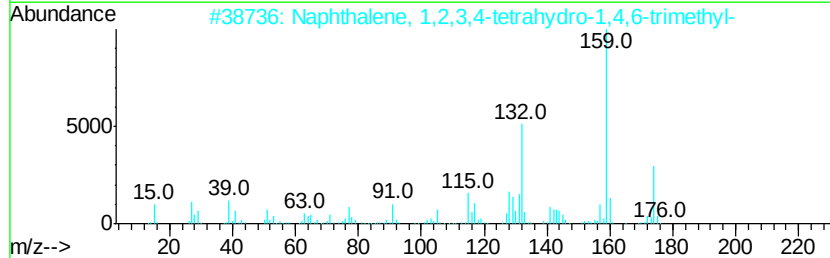
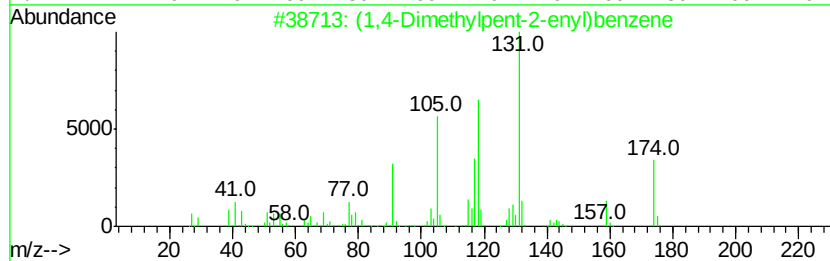
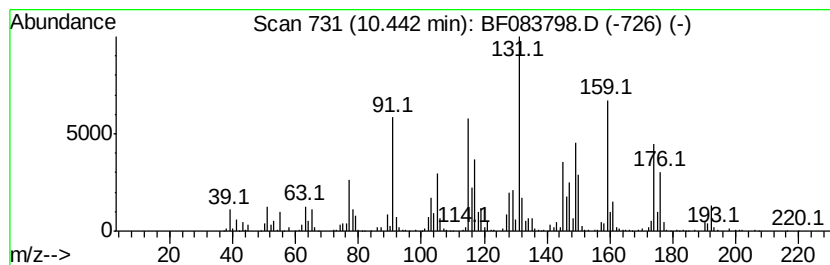
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown10.44 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.44	25.76 ng	2958360	Acenaphthene-d10	9.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(1,4-Dimethylpent-2-enyl)benzene	174	C13H18	1000184-97-9	41
2		Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	022824-32-4	38
3		1H-Indene, 2,3-dihydro-1,1,5,6-t...	174	C13H18	000942-43-8	35
4		Naphthalene, 1,2,3,4-tetrahydro-...	174	C13H18	021693-51-6	35
5		8-Quinolinemethanol	159	C10H9NO	016032-35-2	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121415\
Data File : BF083798.D
Acq On : 15 Dec 2015 4:57
Operator : UM/IZ
Sample : G4779-13
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
W-38

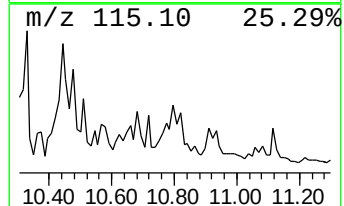
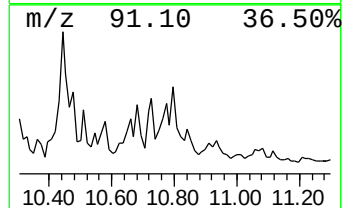
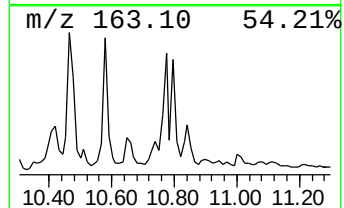
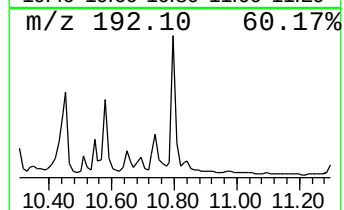
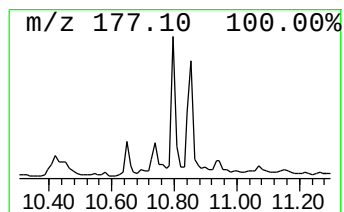
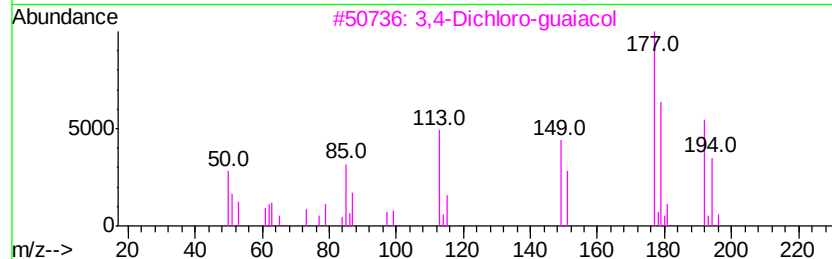
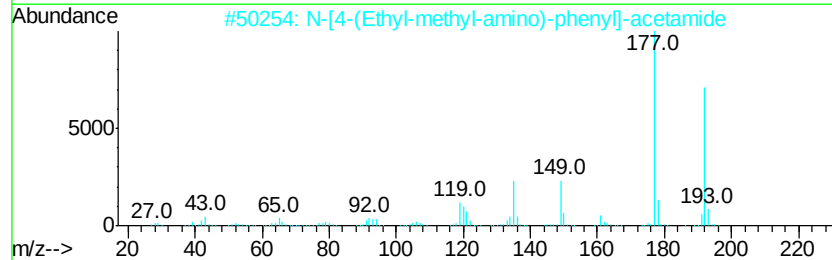
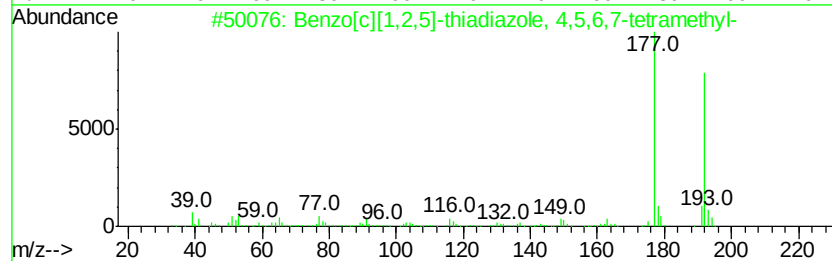
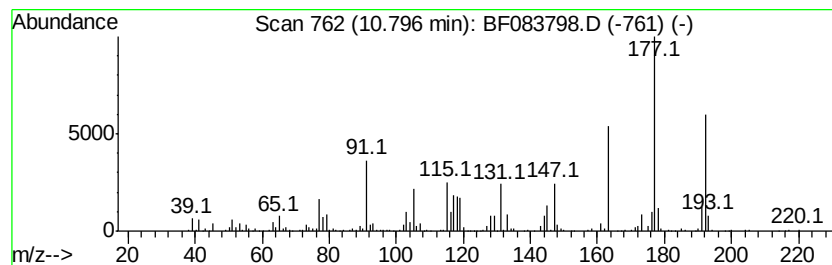
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown10.80 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.80	26.69 ng	1127140	Phenanthrene-d10	11.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[c][1,2,5]-thiadiazole, 4,5...	192	C10H12N2S	106148-64-5	46
2		N-[4-(Ethyl-methyl-amino)-phenyl]...	192	C11H16N2O	099981-45-0	46
3		3,4-Dichloro-quaiacol	192	C7H6Cl2O2	077102-94-4	43
4		3-Buten-1-one, 4-[2,6,6-trimethy...	192	C13H20O	085949-43-5	38
5		2-Propanamine, N-[(3-nitrophenyl...	192	C10H12N2O2	027895-80-3	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121415\
Data File : BF083798.D
Acq On : 15 Dec 2015 4:57
Operator : UM/IZ
Sample : G4779-13
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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1-Phenyl-1-butene	7.93	32.4	ng	1985450	2	8.18	1224430	20.0
Naphthalene, 1,2,...	8.03	19.9	ng	1218390	2	8.18	1224430	20.0
1(2H)-Naphthaleno...	9.61	56.3	ng	6469530	3	9.94	2296850	20.0
1(2H)-Naphthaleno...	10.02	23.3	ng	2674380	3	9.94	2296850	20.0
unknown10.33	10.33	18.8	ng	2155210	3	9.94	2296850	20.0
unknown10.44	10.44	25.8	ng	2958360	3	9.94	2296850	20.0
unknown10.80	10.80	26.7	ng	1127140	4	11.43	844689	20.0