

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102715\
 Data File : BF082564.D
 Acq On : 27 Oct 2015 23:38
 Operator : UM/IZ
 Sample : G4168-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CWC-2-20151026

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-BF102615.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	4.880	236	240	250	rBV	670317	1245559	46.22%	3.746%
2	5.326	276	279	281	rBV	1626226	2254862	83.68%	6.781%
3	6.377	368	371	378	rBV	1928911	2240808	83.15%	6.739%
4	6.480	378	380	383	rBV	1780746	2344697	87.01%	7.052%
5	6.709	398	400	405	rBV	506744	471362	17.49%	1.418%
6	6.869	411	414	416	rBV	1613955	1934497	71.79%	5.818%
7	7.292	448	451	453	rBV	1346551	1580421	58.65%	4.753%
8	8.000	511	513	519	rBV	666661	692614	25.70%	2.083%
9	8.720	574	576	581	rBV	27824	35784	1.33%	0.108%
10	8.812	581	584	588	rBV	37071	39178	1.45%	0.118%
11	9.075	604	607	610	rBV	2159629	2694761	100.00%	8.104%
12	9.417	634	637	638	rBV	26846	35367	1.31%	0.106%
13	9.475	640	642	645	rVV	498280	605295	22.46%	1.820%
14	9.532	645	647	652	rVB2	18785	29383	1.09%	0.088%
15	9.612	652	654	656	rBV	60199	64319	2.39%	0.193%
16	9.749	664	666	668	rBV	701207	697803	25.89%	2.099%
17	9.783	668	669	673	rVB	146538	126599	4.70%	0.381%
18	9.955	682	684	686	rBV	63466	82677	3.07%	0.249%
19	10.149	699	701	702	rBV	20648	27032	1.00%	0.081%
20	10.183	702	704	705	rVB	32245	35530	1.32%	0.107%
21	10.298	712	714	717	rVB	119955	119827	4.45%	0.360%
22	10.389	717	722	727	rBV3	29041	101105	3.75%	0.304%
23	10.549	733	736	738	rBV	1364168	1552234	57.60%	4.668%
24	10.652	743	745	749	rVB2	56348	88004	3.27%	0.265%
25	10.755	749	754	756	rBV2	21946	37347	1.39%	0.112%
26	10.846	759	762	764	rBV2	35790	66822	2.48%	0.201%
27	10.926	766	769	771	rBV4	14667	30928	1.15%	0.093%
28	10.972	771	773	778	rBV3	25306	64024	2.38%	0.193%
29	11.052	778	780	782	rBV	34965	41417	1.54%	0.125%
30	11.098	782	784	785	rBV2	62730	64940	2.41%	0.195%
31	11.132	785	787	791	rVB	66993	83795	3.11%	0.252%
32	11.258	794	798	801	rBV2	1001006	1735757	64.41%	5.220%
33	11.315	801	803	805	rVV	245936	217087	8.06%	0.653%
34	11.486	815	818	820	rBV2	75767	147400	5.47%	0.443%

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35	11.521	820	821	823	rVB2	38967	49709	1.84%	0.149%
36	11.578	823	826	829	rBV4	30016	67028	2.49%	0.202%
37	11.738	838	840	841	rBV	152290	139610	5.18%	0.420%
38	11.772	841	843	845	rVV	336695	385525	14.31%	1.159%
39	11.818	845	847	848	rVV	82864	95689	3.55%	0.288%
40	11.852	848	850	855	rVB2	211350	352011	13.06%	1.059%
41	11.932	855	857	858	rBV2	36556	42274	1.57%	0.127%
42	12.035	864	866	868	rVB	87283	80202	2.98%	0.241%
43	12.069	868	869	871	rBV2	52488	68960	2.56%	0.207%
44	12.183	877	879	880	rBV	35510	40085	1.49%	0.121%
45	12.286	886	888	890	rBV	115048	148385	5.51%	0.446%
46	12.321	890	891	894	rVV3	57533	84491	3.14%	0.254%
47	12.378	894	896	900	rVB2	50527	89333	3.32%	0.269%
48	12.446	900	902	905	rBV	1038019	1188073	44.09%	3.573%
49	12.561	910	912	914	rVV2	75300	115376	4.28%	0.347%
50	12.606	914	916	917	rVV	62618	88419	3.28%	0.266%
51	12.675	919	922	925	rVV	960788	1223076	45.39%	3.678%
52	12.732	925	927	929	rVB	56690	74109	2.75%	0.223%
53	12.823	932	935	939	rVV	2137997	2164925	80.34%	6.511%
54	12.903	939	942	944	rVV2	67935	128145	4.76%	0.385%
55	13.006	947	951	953	rVV2	101548	203455	7.55%	0.612%
56	13.075	956	957	958	rVV	63305	60378	2.24%	0.182%
57	13.109	958	960	964	rVB2	95967	162274	6.02%	0.488%
58	13.521	994	996	1000	rVB	64769	117979	4.38%	0.355%
59	13.635	1004	1006	1009	rVV2	78634	179931	6.68%	0.541%
60	13.715	1011	1013	1019	rVB3	143257	286669	10.64%	0.862%
61	13.875	1024	1027	1029	rBV	684925	1168439	43.36%	3.514%
62	14.755	1102	1104	1106	rBV	67545	80283	2.98%	0.241%
63	14.892	1114	1116	1123	rVB	357349	842580	31.27%	2.534%
64	14.995	1123	1125	1127	rBV	76975	107632	3.99%	0.324%
65	15.178	1138	1141	1143	rBV	225957	267673	9.93%	0.805%
66	15.235	1144	1146	1149	rVB	335311	379271	14.07%	1.141%
67	15.292	1149	1151	1153	rBV	301671	368120	13.66%	1.107%
68	16.275	1234	1237	1242	rVB2	74957	165338	6.14%	0.497%
69	16.652	1267	1270	1275	rVB	164408	341163	12.66%	1.026%
70	17.052	1302	1305	1311	rVB	141175	304996	11.32%	0.917%

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

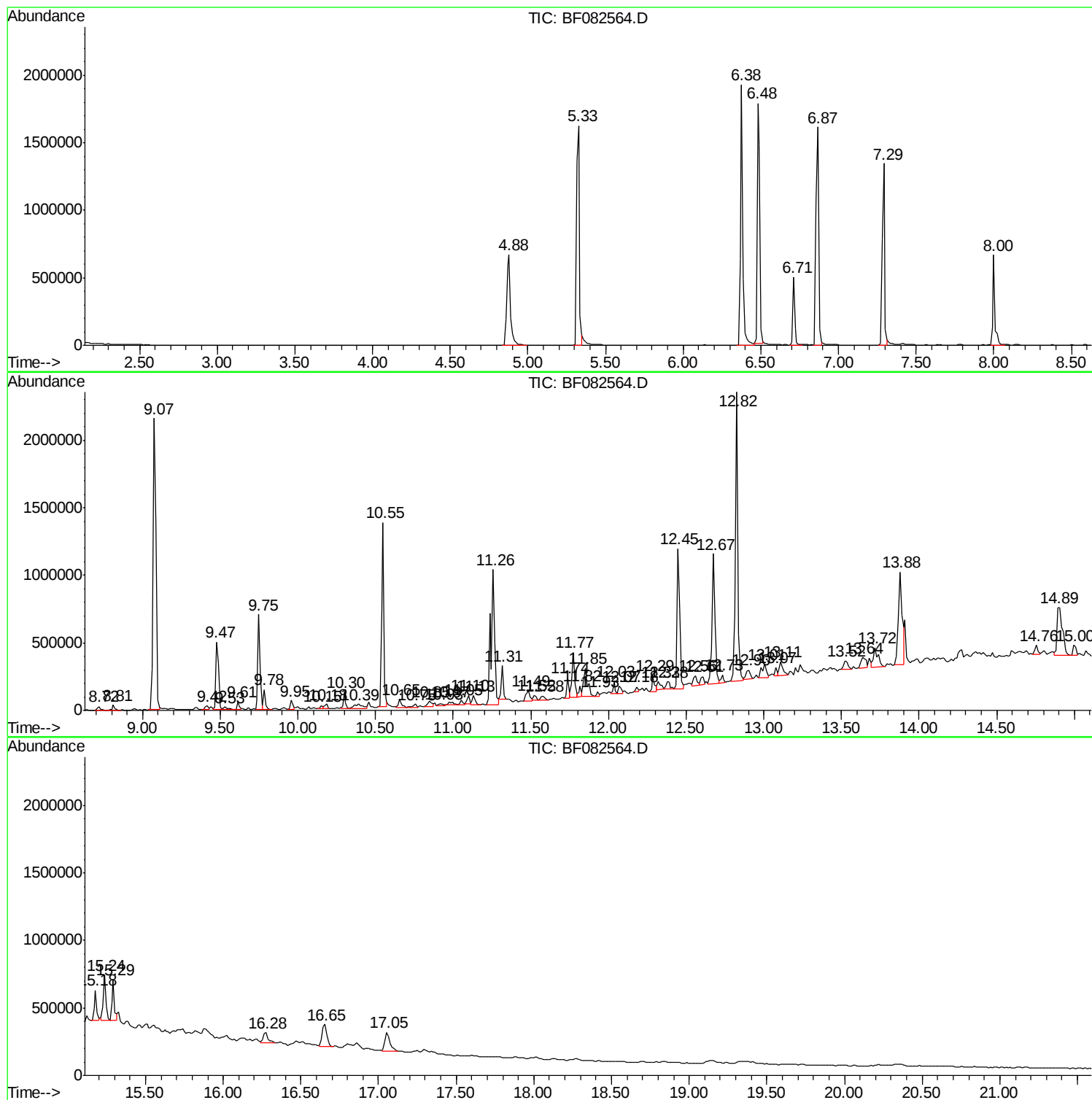
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Misc :
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Instrument :
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ClientSampleId :
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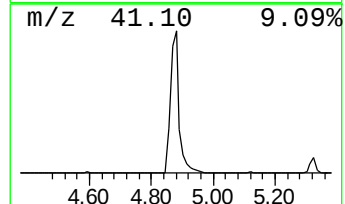
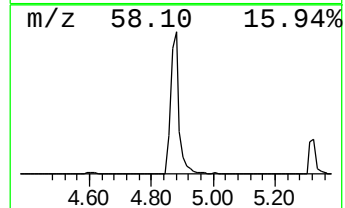
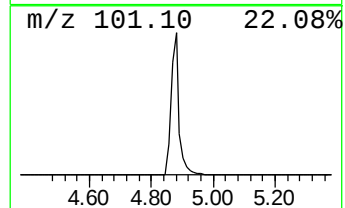
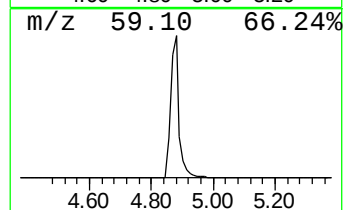
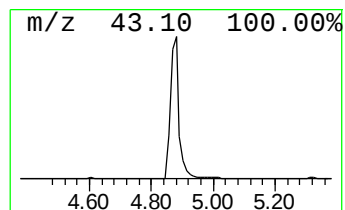
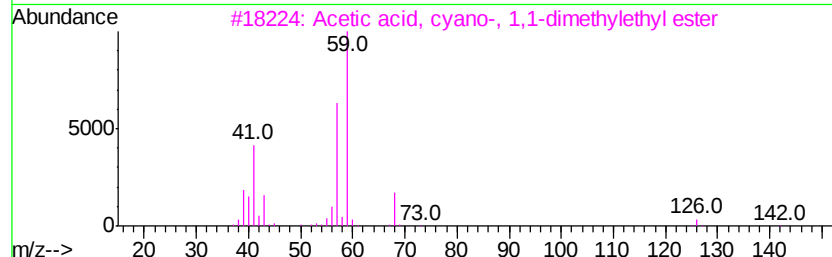
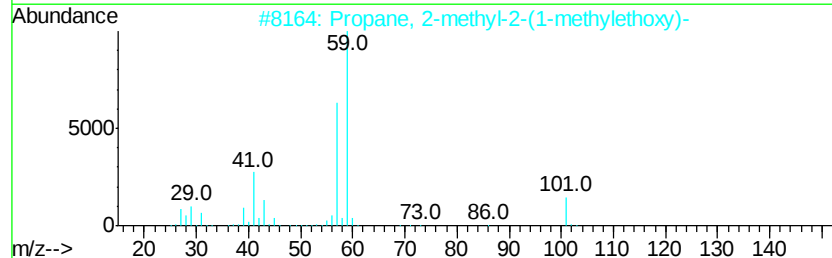
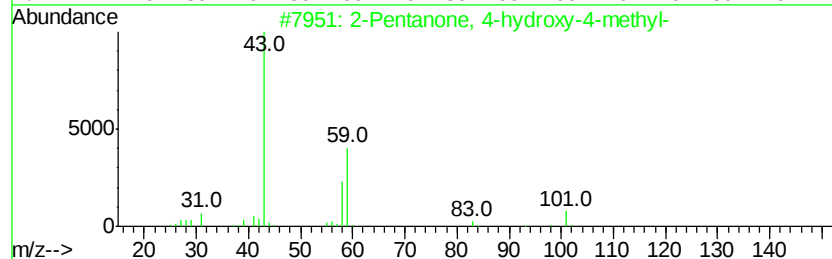
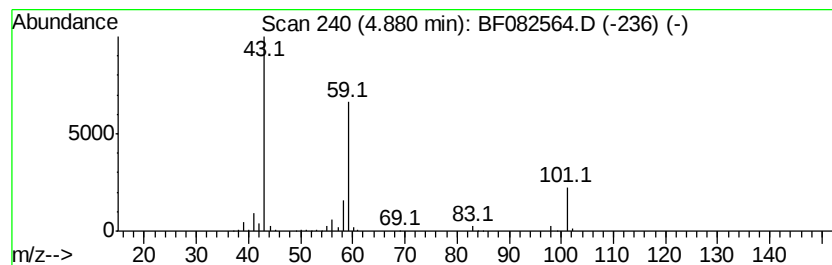
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102615.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.88	52.85 ng	1245560	1,4-Dichlorobenzene-d4	6.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	53
3		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	23
4		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5		Guanidine	59	CH5N3	000113-00-8	9



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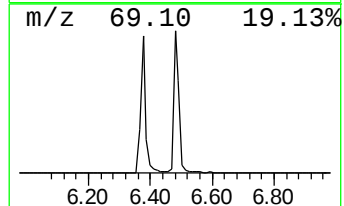
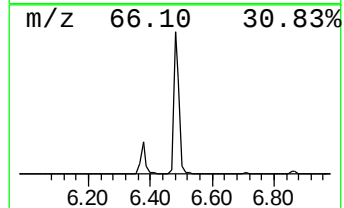
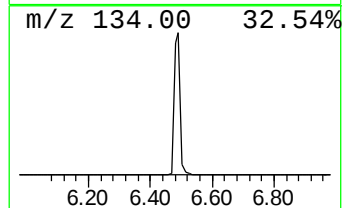
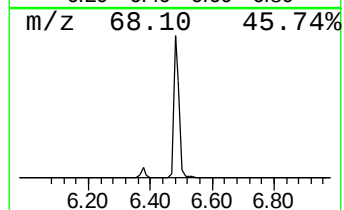
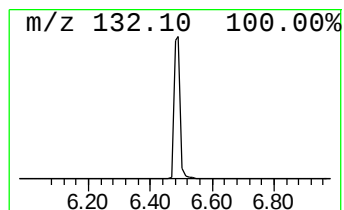
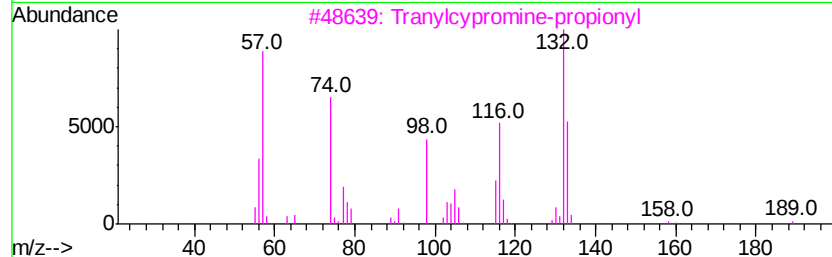
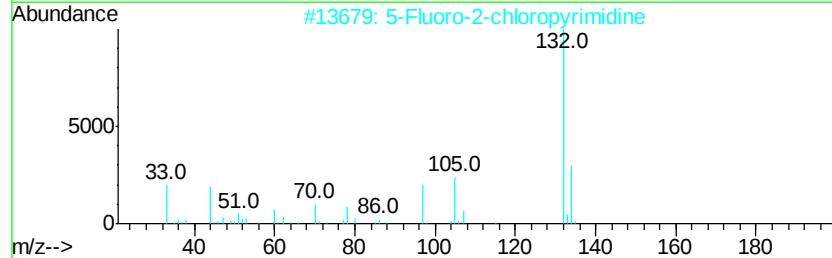
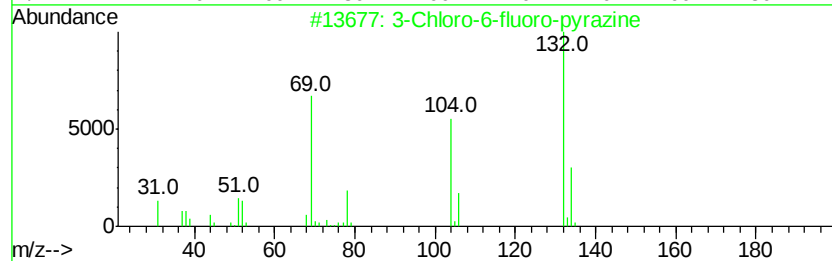
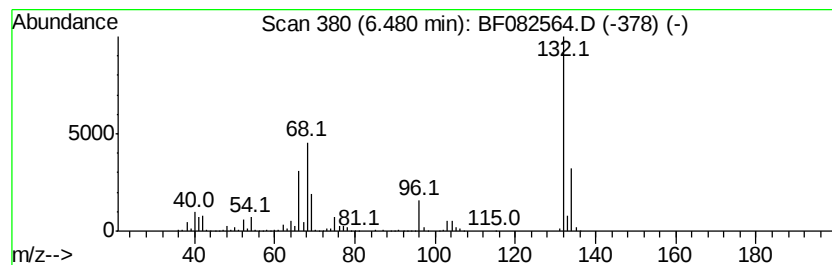
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.48	99.49 ng	2344700	1,4-Dichlorobenzene-d4	6.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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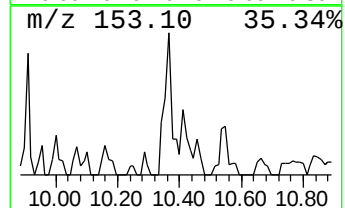
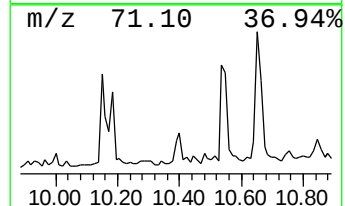
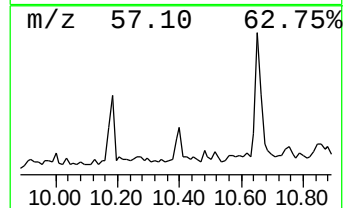
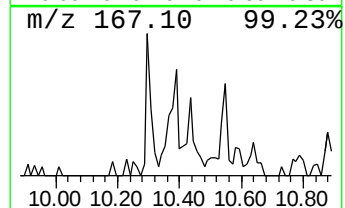
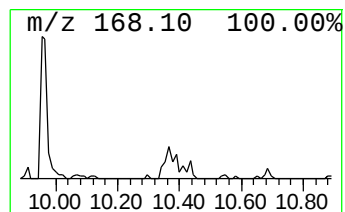
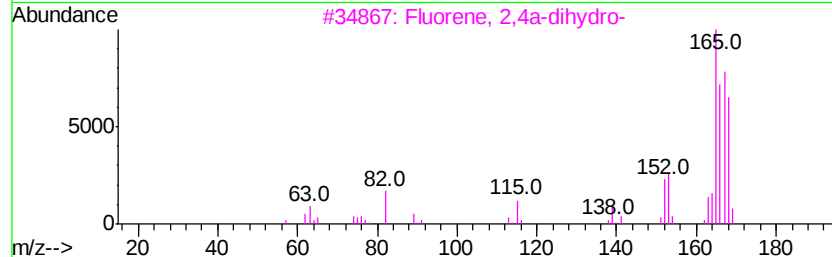
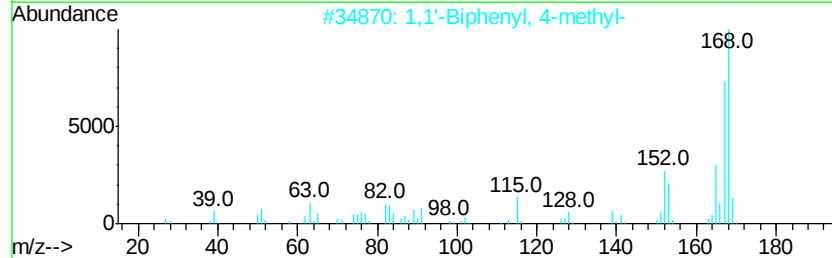
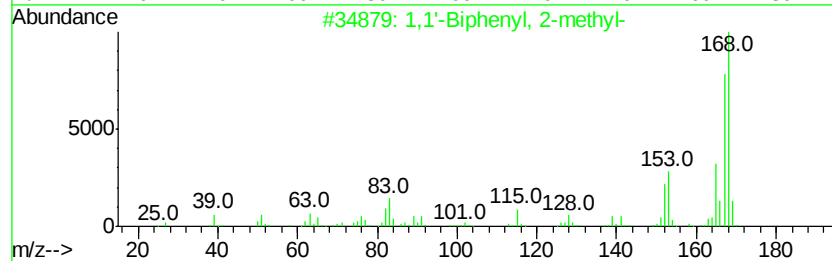
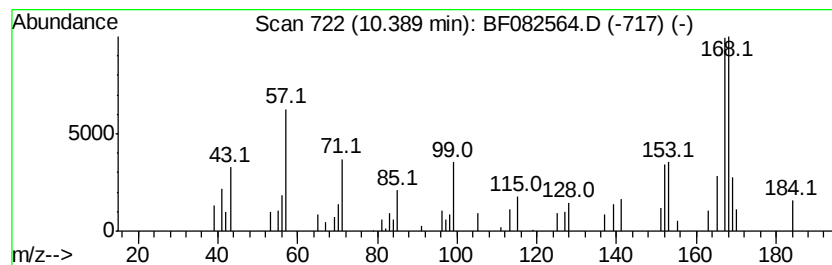
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 1,1'-Biphenyl, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.39	2.90 ng	101105	Acenaphthene-d10		9.75
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	70
2	1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	58
3	Fluorene, 2,4a-dihydro-	168	C13H12	059247-36-8	58
4	Fluorene, 1,4-dihydro-	168	C13H12	041593-21-9	58
5	Diphenylmethane	168	C13H12	000101-81-5	58



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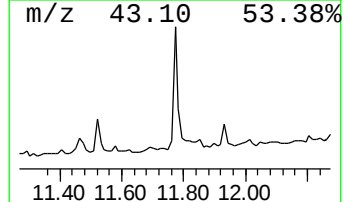
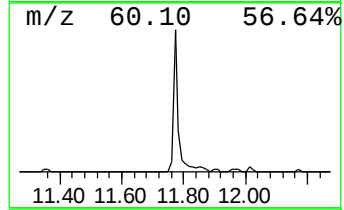
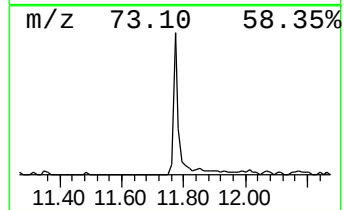
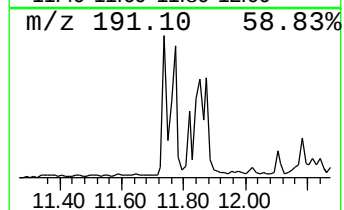
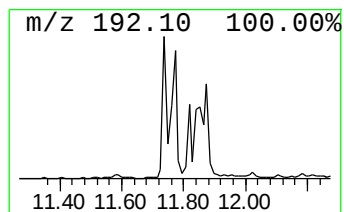
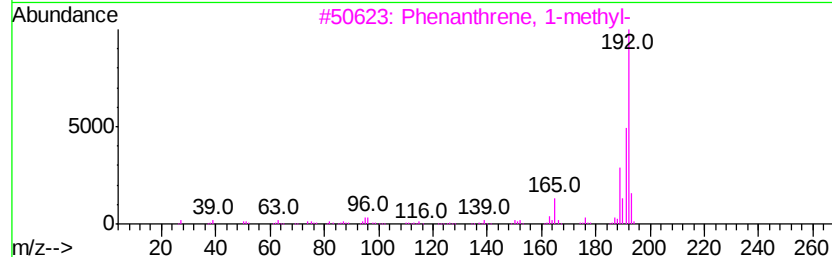
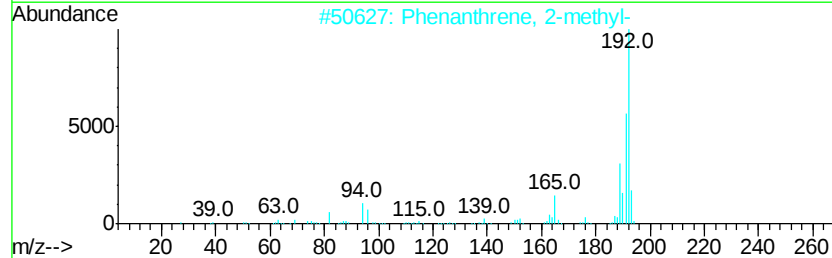
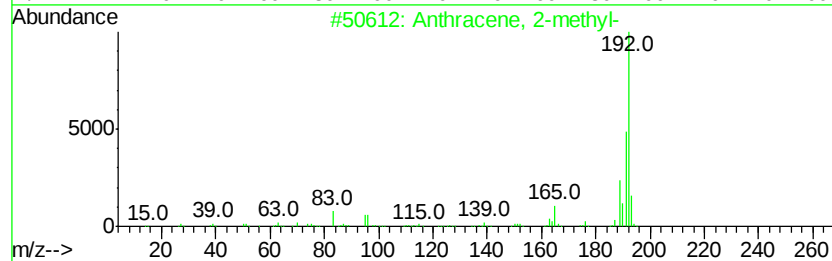
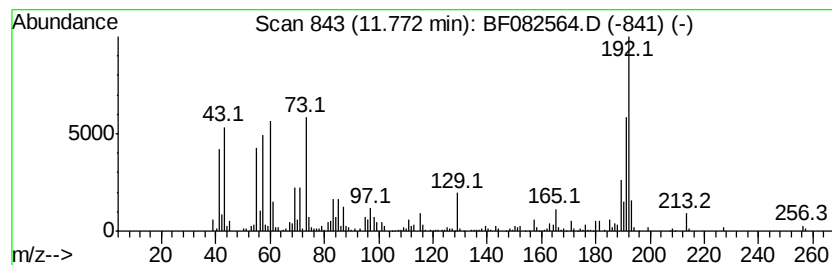
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Anthracene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.77	4.44 ng	385525	Phenanthrene-d10		11.23	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	92
5		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102715\
Data File : BF082564.D
Acq On : 27 Oct 2015 23:38
Operator : UM/IZ
Sample : G4168-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

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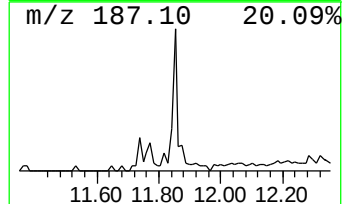
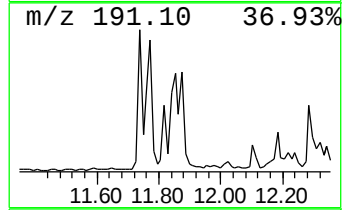
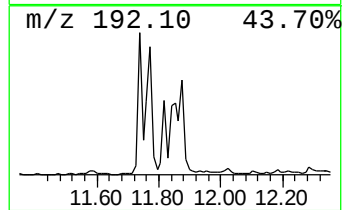
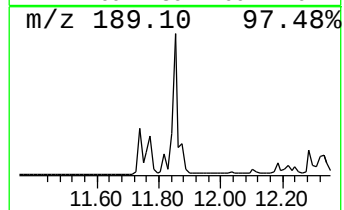
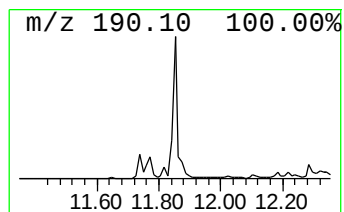
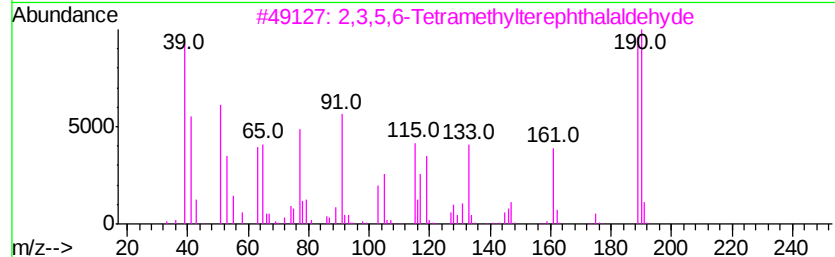
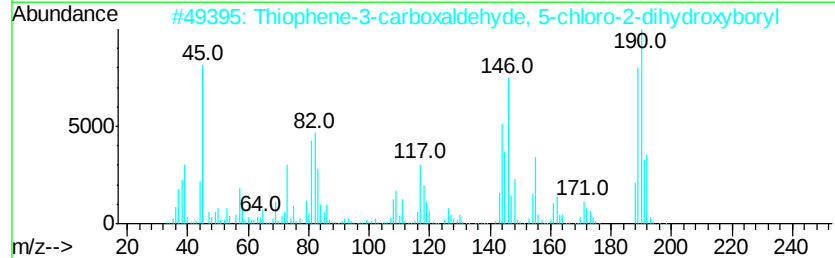
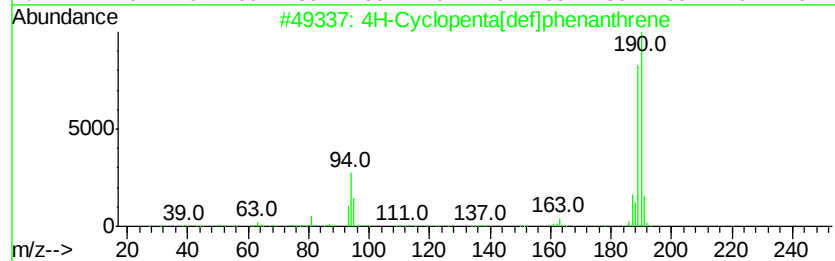
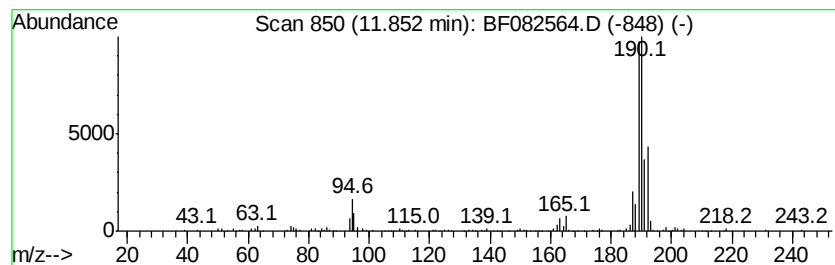
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.85	4.06 ng	352011	Phenanthrene-d10	11.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
2		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	50
3		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	47
4		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40
5		3H-Indazole-3-carbonitrile, 3-(4...	253	C14H8ClN3	056630-97-8	17



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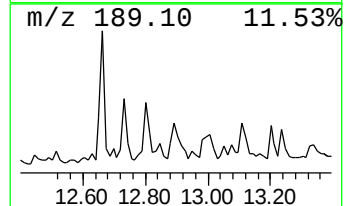
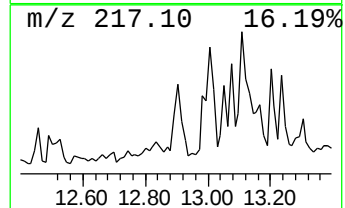
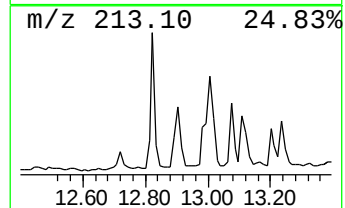
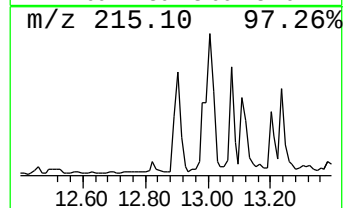
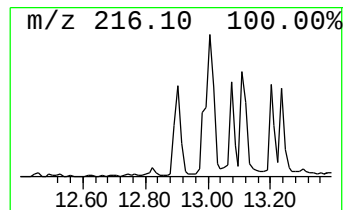
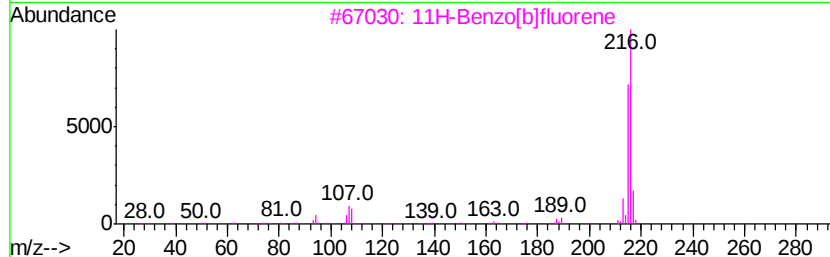
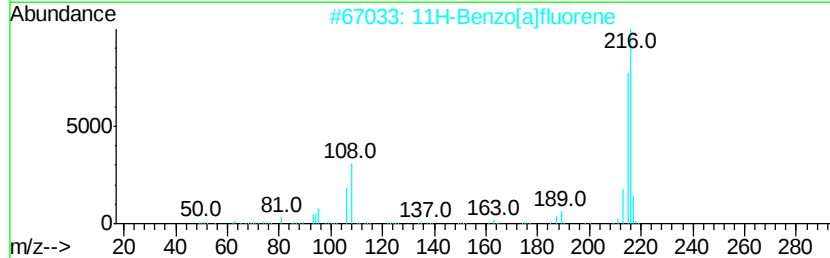
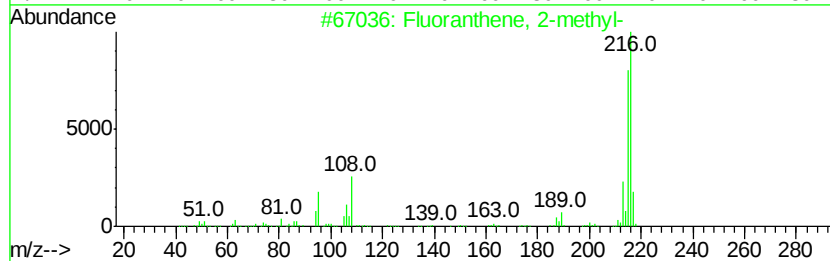
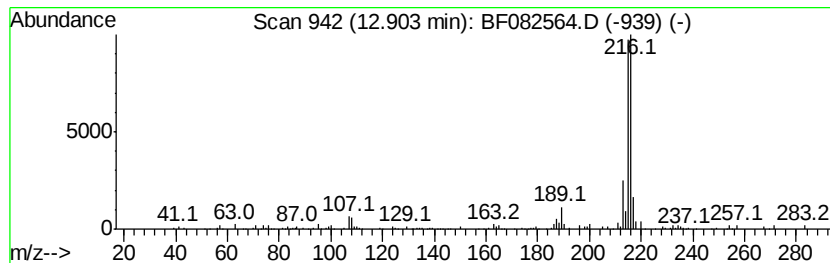
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Fluoranthene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.90	2.19 ng	128145	Chrysene-d12	13.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
3		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	86
4		Pvrene, 1-methyl-	216	C17H12	002381-21-7	70
5		Pyridin-3-ol, 2-(4-nitrophenyl)-	216	C11H8N2O3	030820-89-4	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102715\
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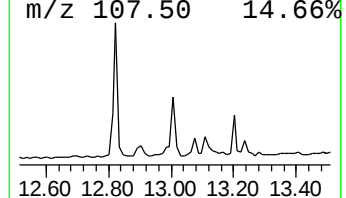
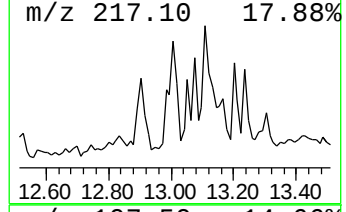
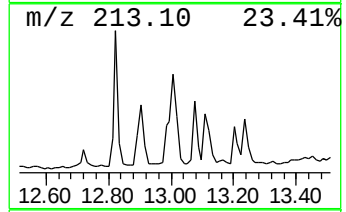
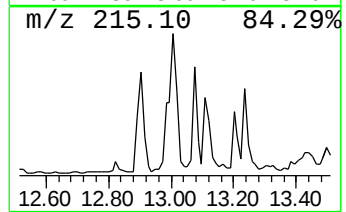
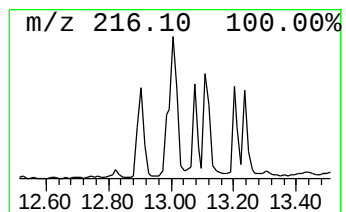
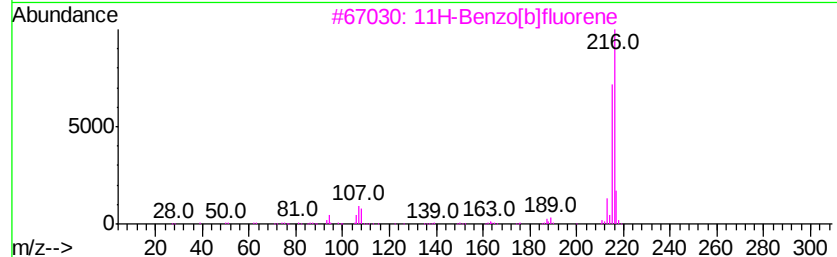
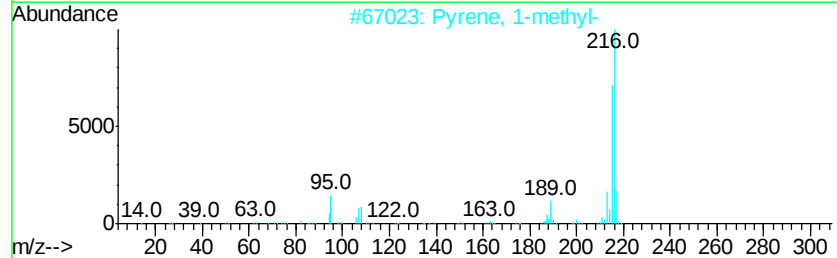
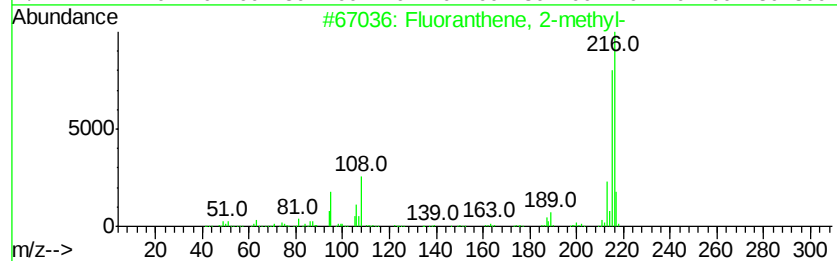
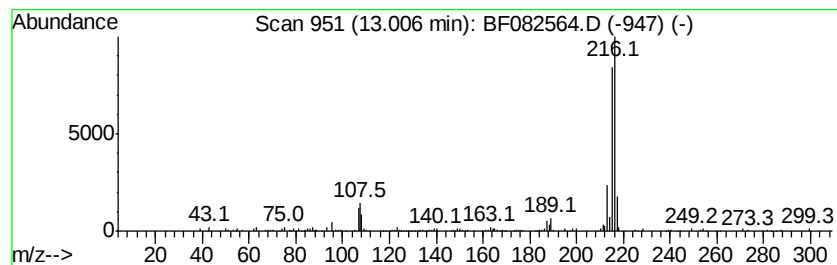
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Pyrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.01	3.48 ng	203455	Chrysene-d12	13.88

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



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102715\
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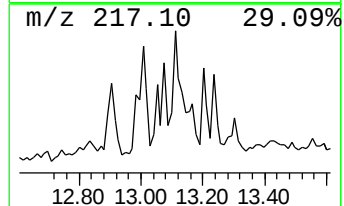
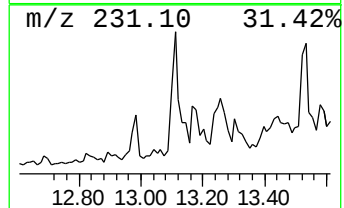
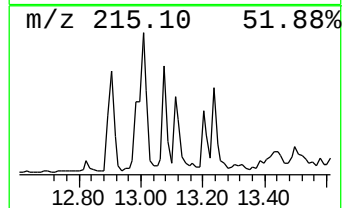
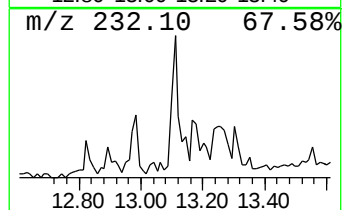
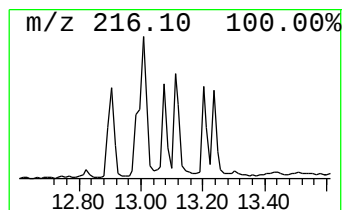
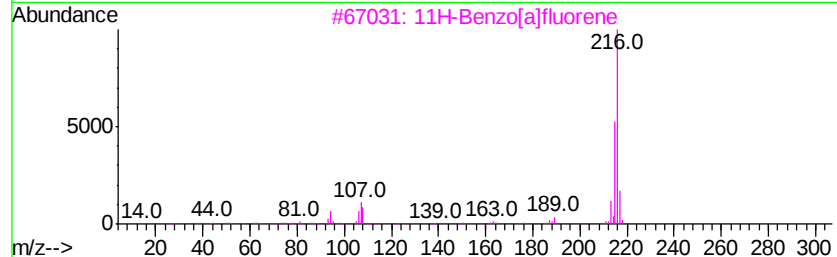
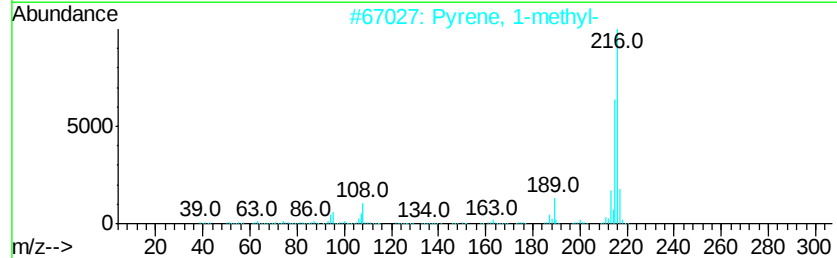
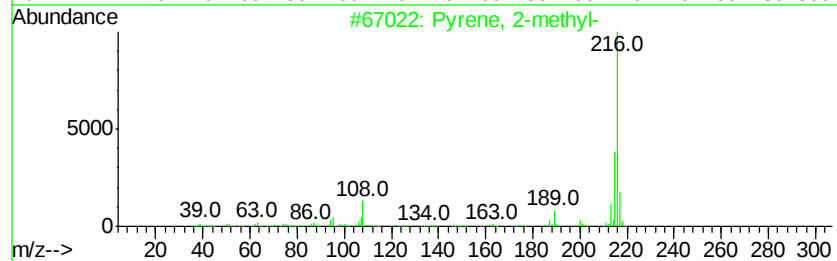
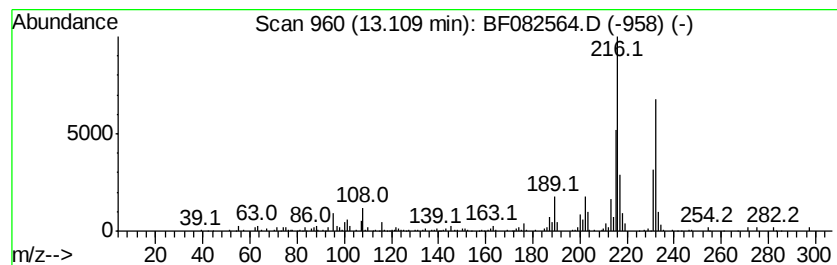
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Pyrene, 2-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.11	2.78 ng	162274	Chrysene-d12	13.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	91
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	70
3		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	60
4		Pyrene, 4-methyl-	216	C17H12	003353-12-6	60
5		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	46



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102715\
Data File : BF082564.D
Acq On : 27 Oct 2015 23:38
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ClientSampleId :
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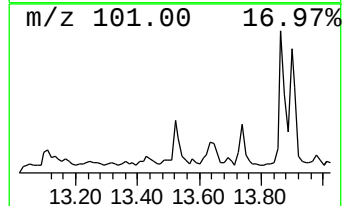
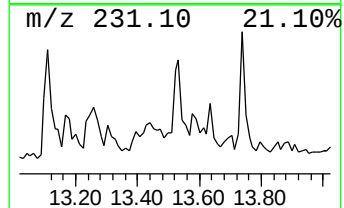
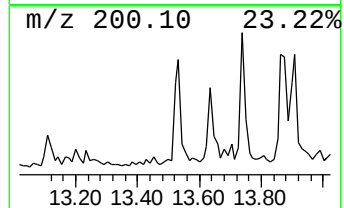
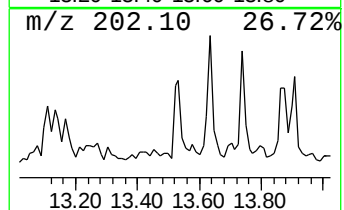
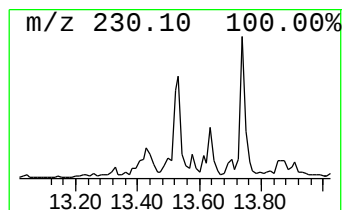
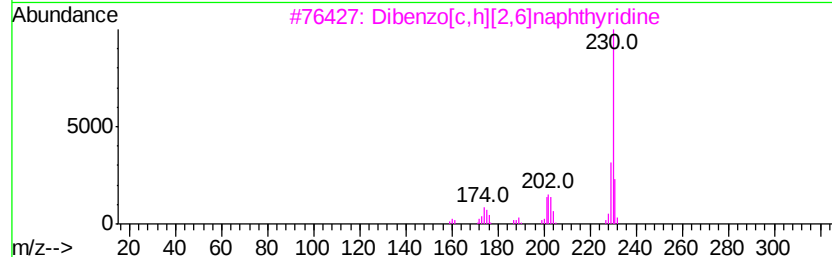
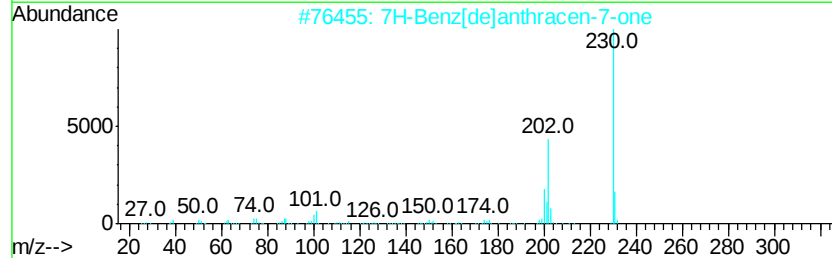
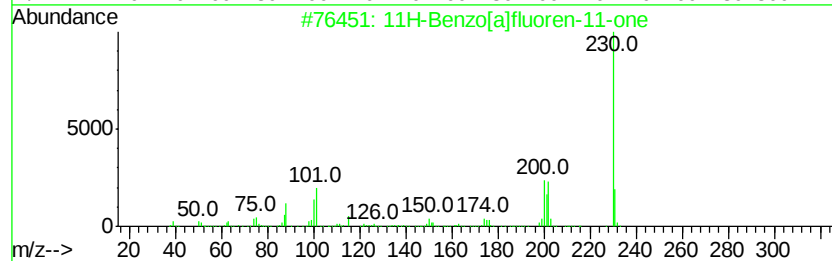
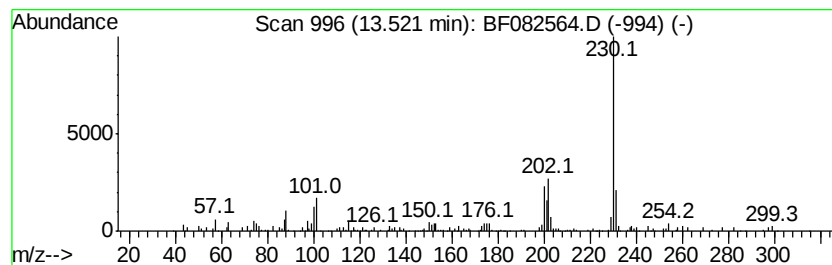
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 11H-Benzo[a]fluoren-11-one Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	2.02 ng	117979	Chrysene-d12	13.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	99
2		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	91
3		Dibenzo[c,h][2,6]naphthvridine	230	C16H10N2	000218-30-4	59
4		o-Terphenyl	230	C18H14	000084-15-1	53
5		5,6-Dihydrochrysene	230	C18H14	002091-92-1	53



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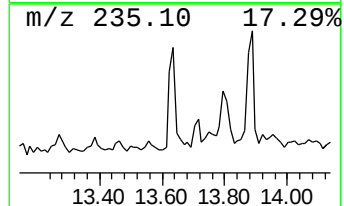
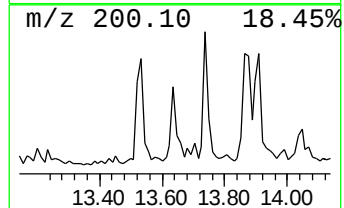
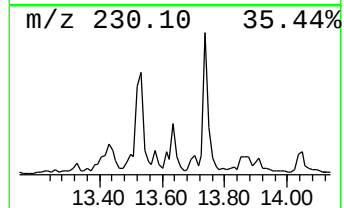
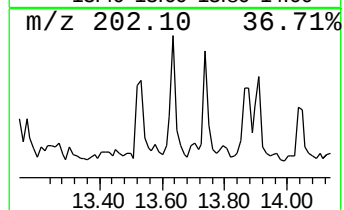
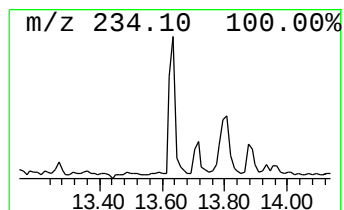
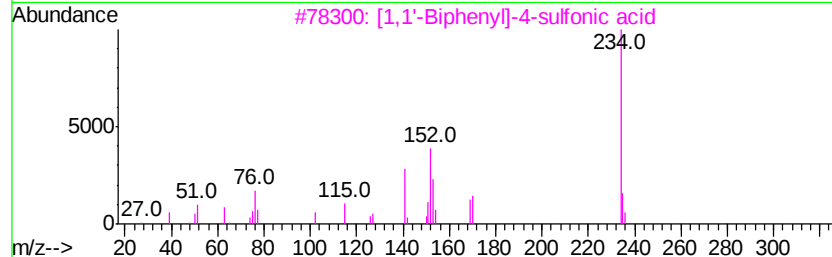
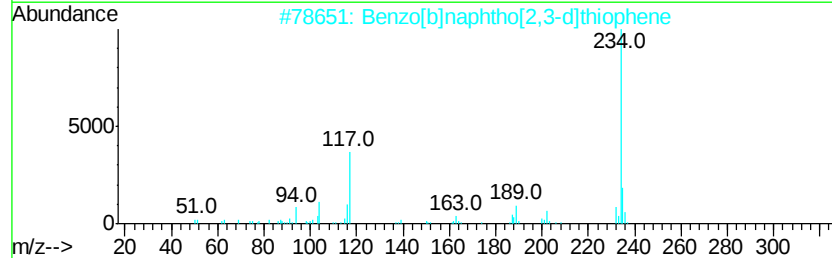
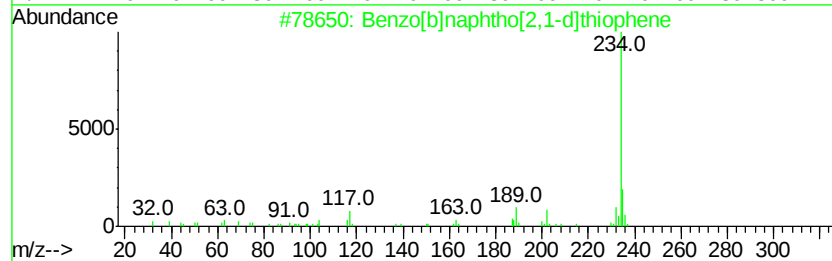
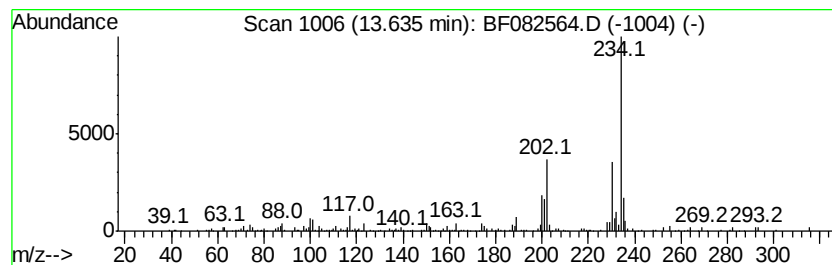
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.64	3.08 ng	179931	Chrysene-d12		13.88	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	70
2		Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	58
3		[1,1'-Biphenyl]-4-sulfonic acid	234	C12H10O3S	002113-68-0	46
4		Oxantrene-2,3-dicarbonitrile	234	C14H6N2O2	309735-17-9	43
5		Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	40



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102715\
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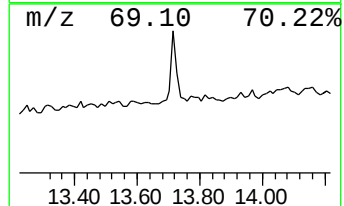
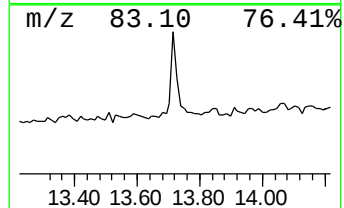
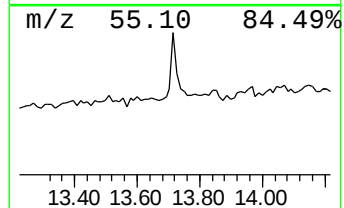
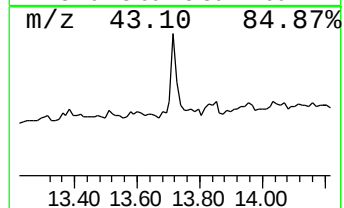
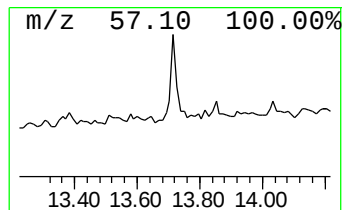
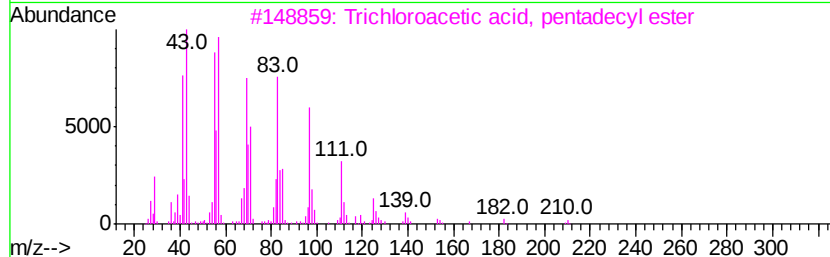
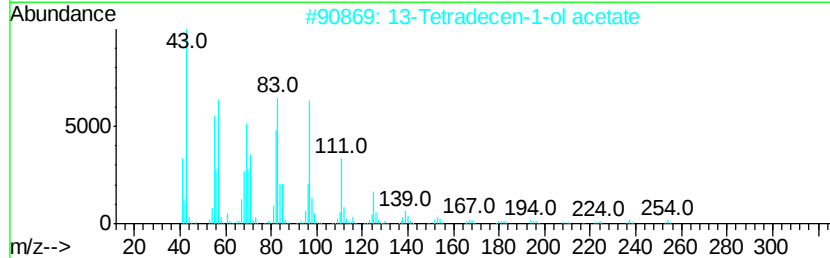
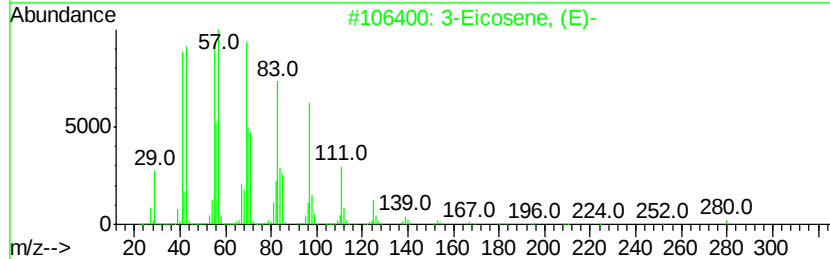
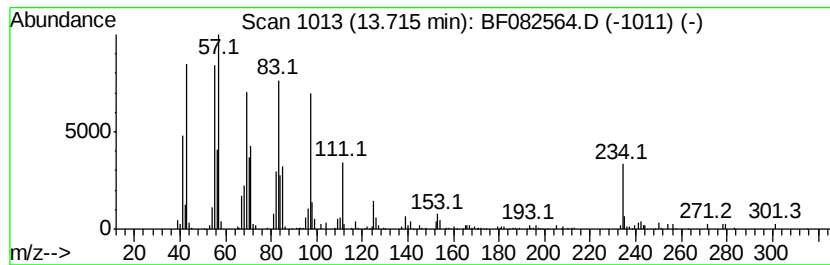
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 3-Eicosene, (E)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	4.91 ng	286669	Chrysene-d12	13.88

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Eicosene, (E)-	280	C20H40	074685-33-9	90
2			13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	87
3			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	87
4			Dichloroacetic acid, tetradecyl ...	324	C16H30Cl2O2	083005-02-1	87
5			Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102715\
Data File : BF082564.D
Acq On : 27 Oct 2015 23:38
Operator : UM/IZ
Sample : G4168-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CWC-2-20151026

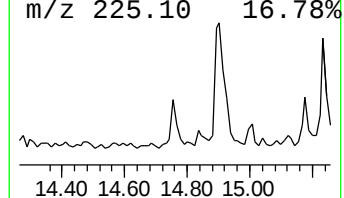
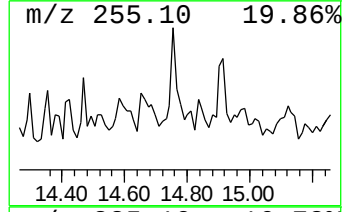
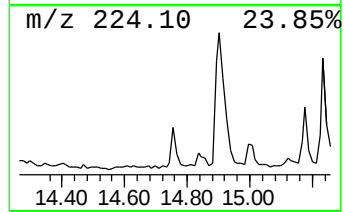
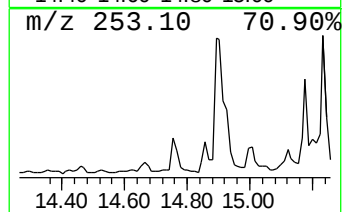
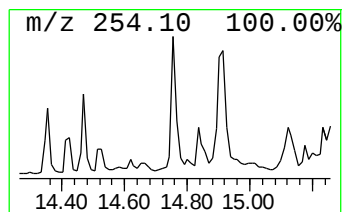
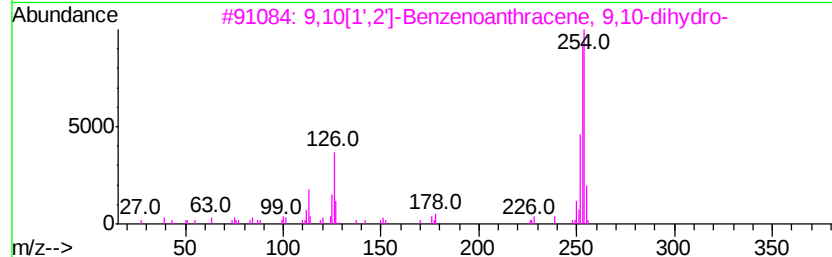
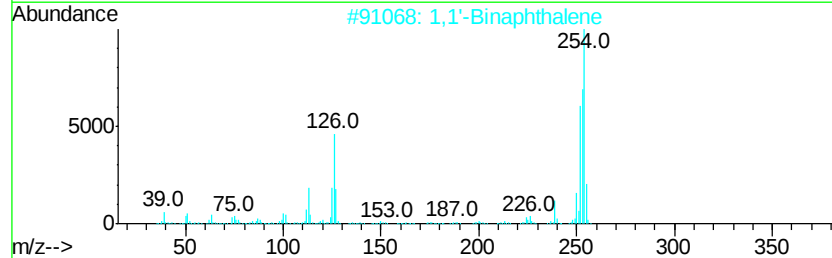
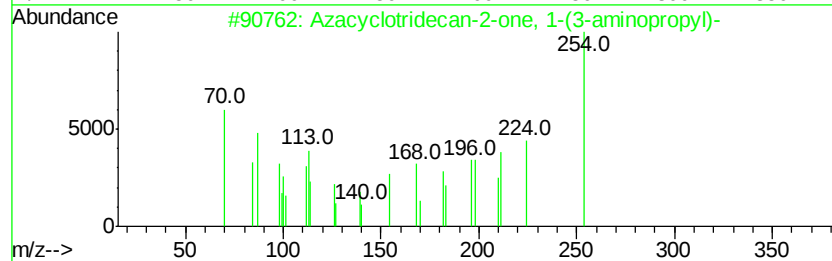
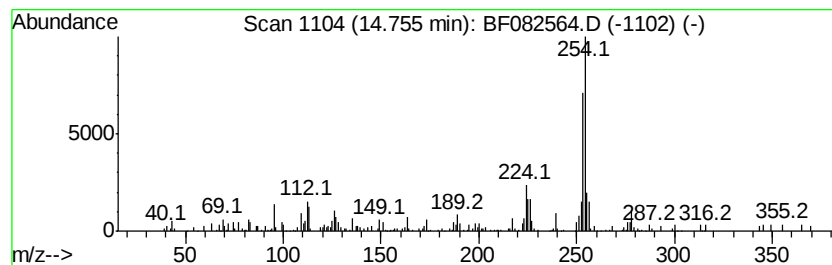
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Azacyclotridecan-2-one, 1-(... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	4.36 ng	80283	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Azacyclotridecan-2-one, 1-(3-ami...	254	C15H30N2O	064414-61-5	89
2		1,1'-Binaphthalene	254	C20H14	000604-53-5	70
3		9,10[1',2'-Benzenoanthracene, 9...	254	C20H14	000477-75-8	62
4		Benzenamine, 4-(2-benzothiazolyl...	254	C15H14N2S	010205-56-8	58
5		1,2'-Binaphthalene	254	C20H14	004325-74-0	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102715\
Data File : BF082564.D
Acq On : 27 Oct 2015 23:38
Operator : UM/IZ
Sample : G4168-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CWC-2-20151026

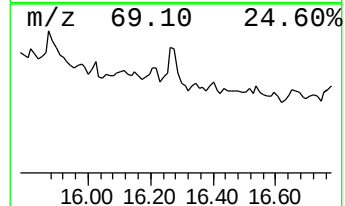
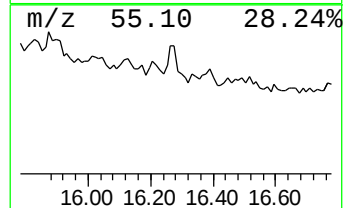
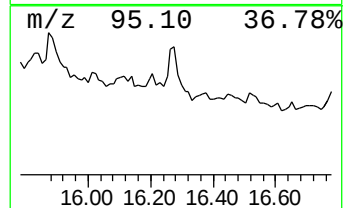
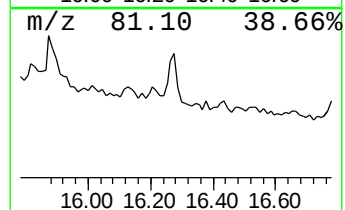
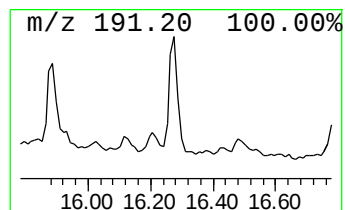
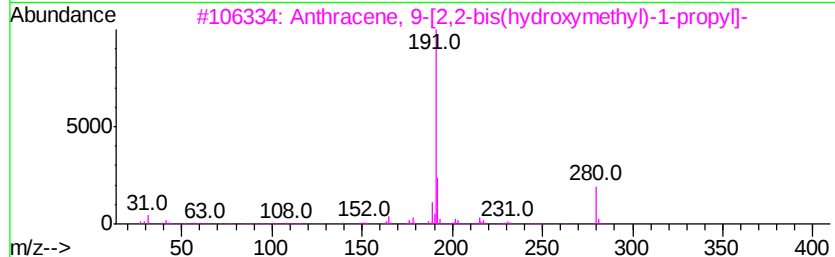
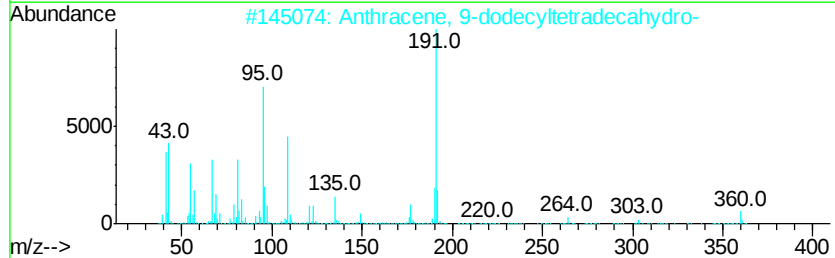
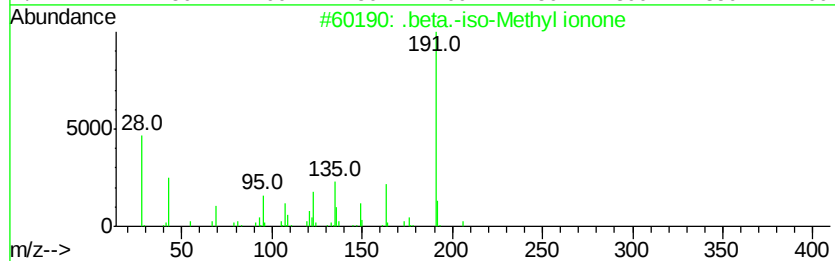
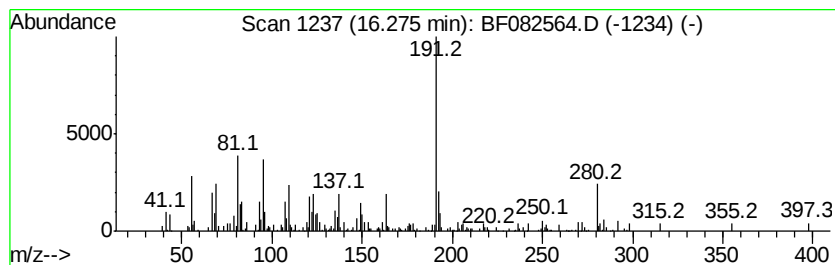
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 .beta.-iso-Methyl ionone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.28	8.98 ng	165338	Perylene-d12	15.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	55
2		Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	47
3		Anthracene, 9-[2,2-bis(hydroxyme...	280	C19H20O2	144000-85-1	46
4		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	45
5		3-Buten-2-one, 4-(2,5,6,6-tetram...	206	C14H22O	000079-70-9	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF102715\
 Data File : BF082564.D
 Acq On : 27 Oct 2015 23:38
 Operator : UM/IZ
 Sample : G4168-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CWC-2-20151026

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102615.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...	4.88	52.9	ng	1245560	1	6.71	471362	20.0
unknown6.48	6.48	99.5	ng	2344700	1	6.71	471362	20.0
1,1'-Biphenyl, 2-...	10.39	2.9	ng	101105	3	9.75	697803	20.0
Anthracene, 2-met...	11.77	4.4	ng	385525	4	11.23	1735760	20.0
4H-Cyclopenta[def...	11.85	4.1	ng	352011	4	11.23	1735760	20.0
Fluoranthene, 2-m...	12.90	2.2	ng	128145	5	13.88	1168440	20.0
Pyrene, 1-methyl-	13.01	3.5	ng	203455	5	13.88	1168440	20.0
Pyrene, 2-methyl-	13.11	2.8	ng	162274	5	13.88	1168440	20.0
11H-Benzo[a]fluor...	13.52	2.0	ng	117979	5	13.88	1168440	20.0
Benzo[b]naphtho[2...	13.64	3.1	ng	179931	5	13.88	1168440	20.0
3-Eicosene, (E)-	13.72	4.9	ng	286669	5	13.88	1168440	20.0
Azacyclotridecan-...	14.76	4.4	ng	80283	6	15.29	368120	20.0
.beta.-iso-Methyl...	16.28	9.0	ng	165338	6	15.29	368120	20.0