

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060816\
 Data File : BF087820.D
 Acq On : 8 Jun 2016 20:19
 Operator : UM/SJ
 Sample : H3379-15 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-10

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-BF060716.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	1.758	14	15	24	rVV	127571	230604	7.27%	1.139%
2	1.884	24	26	38	rVB	1997492	3171394	100.00%	15.668%
3	4.993	295	298	302	rVB	57338	83226	2.62%	0.411%
4	5.416	332	335	337	rBV	993942	935190	29.49%	4.620%
5	6.456	424	426	428	rVB	1054123	999136	31.50%	4.936%
6	6.593	435	438	440	rBV	1008988	1000702	31.55%	4.944%
7	6.822	456	458	461	rBV	607367	592223	18.67%	2.926%
8	6.982	469	472	474	rVB	916529	745967	23.52%	3.685%
9	7.382	505	507	510	rBV	472786	521923	16.46%	2.579%
10	8.113	568	571	572	rBV	964287	871030	27.47%	4.303%
11	8.136	572	573	575	rVB	80768	55728	1.76%	0.275%
12	8.822	631	633	635	rVB	96909	76371	2.41%	0.377%
13	8.925	639	642	644	rBB	81678	70033	2.21%	0.346%
14	9.188	663	665	668	rBV	928760	834005	26.30%	4.120%
15	9.450	686	688	690	rBB	40624	49526	1.56%	0.245%
16	9.530	692	695	696	rBV	59143	64430	2.03%	0.318%
17	9.588	698	700	702	rVV	120607	95679	3.02%	0.473%
18	9.645	703	705	710	rVB	30709	31971	1.01%	0.158%
19	9.873	722	725	726	rBV	709213	817885	25.79%	4.041%
20	9.896	726	727	729	rVB	74266	70829	2.23%	0.350%
21	10.068	740	742	744	rVB	69709	66871	2.11%	0.330%
22	10.411	771	772	775	rVB	101717	90576	2.86%	0.447%
23	10.651	791	793	796	rBV	551523	537871	16.96%	2.657%
24	10.959	817	820	822	rBV2	21021	33252	1.05%	0.164%
25	11.245	843	845	849	rVB2	47196	61777	1.95%	0.305%
26	11.348	852	854	855	rBV	572518	645284	20.35%	3.188%
27	11.371	855	856	858	rVV	459177	565184	17.82%	2.792%
28	11.428	858	861	863	rVB	176720	158260	4.99%	0.782%
29	11.576	871	874	876	rBV2	44644	63484	2.00%	0.314%
30	11.851	896	898	899	rBV	75472	69852	2.20%	0.345%
31	11.885	899	901	903	rVB	94091	98958	3.12%	0.489%
32	11.931	903	905	906	rBV	41416	39291	1.24%	0.194%
33	11.965	906	908	912	rVB2	119767	168956	5.33%	0.835%
34	12.148	923	924	928	rVB2	49246	65462	2.06%	0.323%

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35	12.399	943	946	948	rBV	37671	49070	1.55%	0.242%
36	12.434	948	949	952	rVV2	20189	40287	1.27%	0.199%
37	12.479	952	953	958	rVB3	20790	34723	1.09%	0.172%
38	12.559	958	960	963	rBV	519351	495710	15.63%	2.449%
39	12.708	970	973	977	rBV4	18583	39031	1.23%	0.193%
40	12.788	977	980	983	rVV	485580	525780	16.58%	2.598%
41	12.937	991	993	996	rVB	714903	654534	20.64%	3.234%
42	13.005	996	999	1003	rBV3	32151	65381	2.06%	0.323%
43	13.119	1003	1009	1011	rBV	71239	120660	3.80%	0.596%
44	13.188	1011	1015	1016	rBV	48291	53224	1.68%	0.263%
45	13.222	1016	1018	1020	rBV	63204	70228	2.21%	0.347%
46	13.348	1027	1029	1031	rVV2	28392	31937	1.01%	0.158%
47	13.519	1040	1044	1050	rBV5	18820	59137	1.86%	0.292%
48	13.622	1050	1053	1056	rBV	48740	61730	1.95%	0.305%
49	13.737	1060	1063	1064	rBV2	47641	72932	2.30%	0.360%
50	13.988	1082	1085	1092	rVB2	838062	1181573	37.26%	5.838%
51	14.377	1114	1119	1120	rBV	60826	102258	3.22%	0.505%
52	14.514	1128	1131	1133	rVB	188633	223109	7.04%	1.102%
53	15.005	1171	1174	1179	rBV	168863	345339	10.89%	1.706%
54	15.108	1180	1183	1185	rVV	90899	146548	4.62%	0.724%
55	15.188	1188	1190	1194	rVV3	26450	69365	2.19%	0.343%
56	15.291	1196	1199	1202	rVV	134879	210524	6.64%	1.040%
57	15.348	1202	1204	1207	rVV	120907	173470	5.47%	0.857%
58	15.417	1207	1210	1217	rVV	453647	668728	21.09%	3.304%
59	15.771	1237	1241	1244	rVB	72120	150723	4.75%	0.745%
60	15.885	1248	1251	1253	rBV2	27417	60817	1.92%	0.300%
61	16.343	1287	1291	1297	rVB2	55999	170723	5.38%	0.843%
62	16.800	1328	1331	1335	rVB2	50530	105627	3.33%	0.522%
63	17.028	1347	1351	1357	rVB3	41020	148138	4.67%	0.732%
64	17.211	1364	1367	1370	rBV	30853	61111	1.93%	0.302%
65	17.851	1421	1423	1433	rVB4	18117	65470	2.06%	0.323%

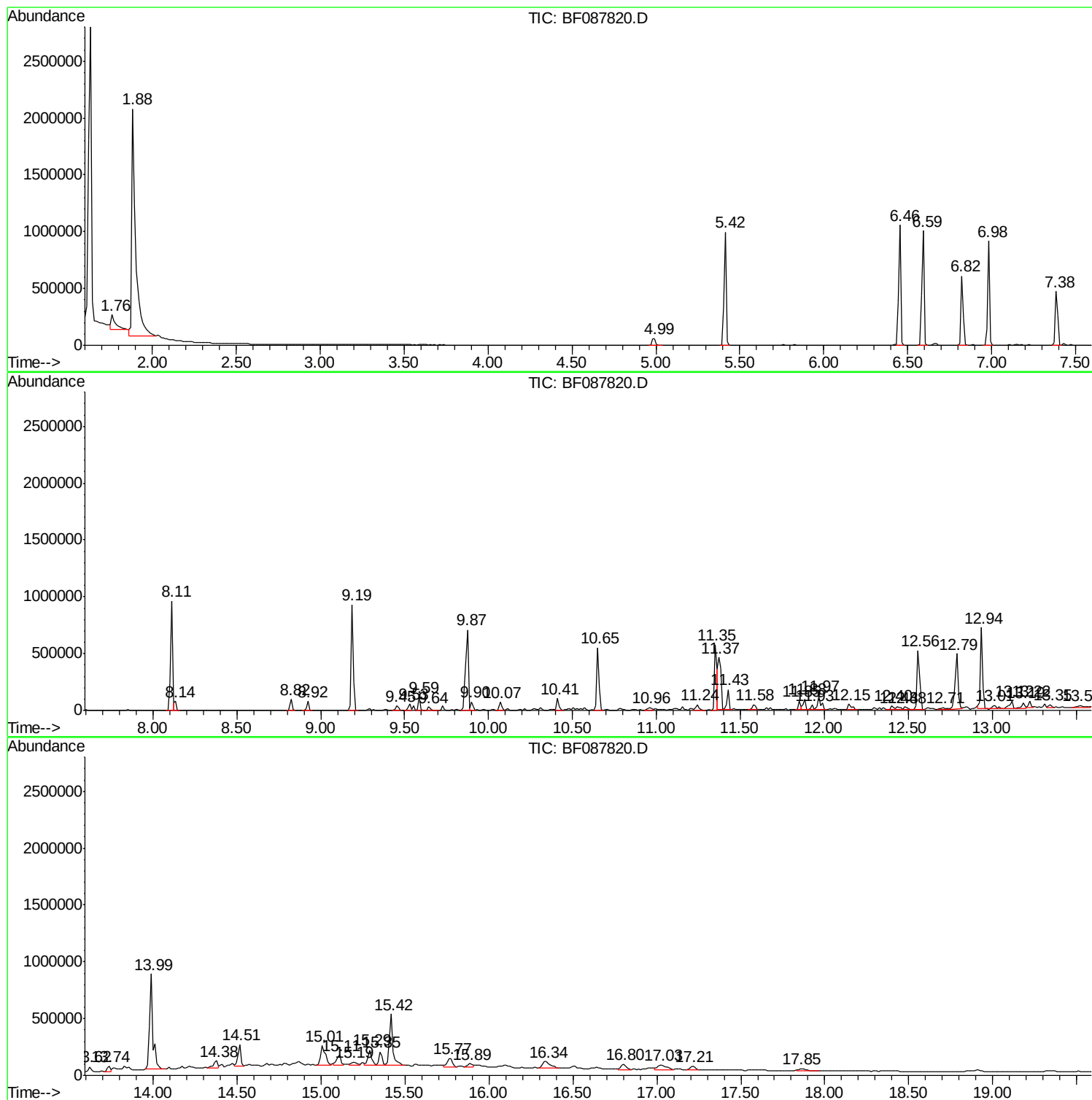
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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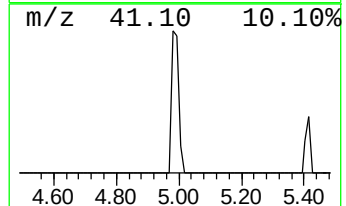
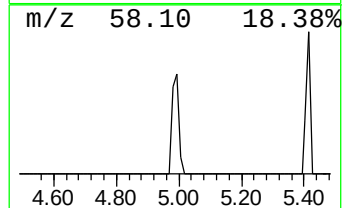
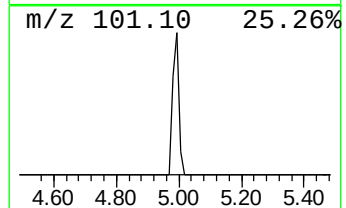
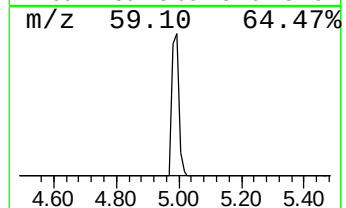
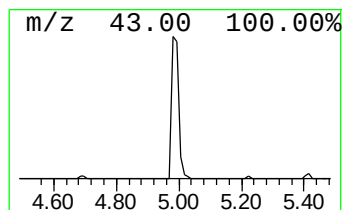
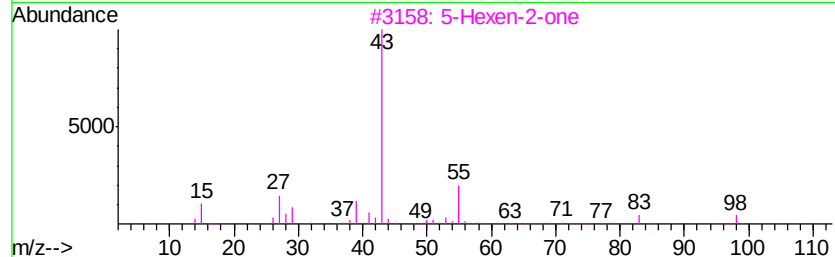
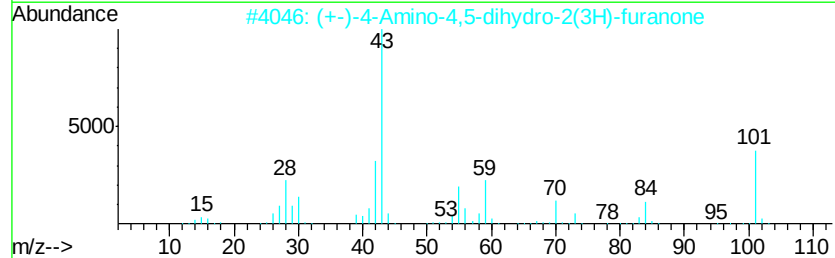
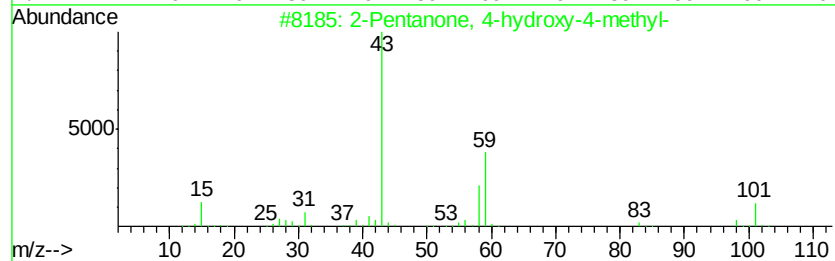
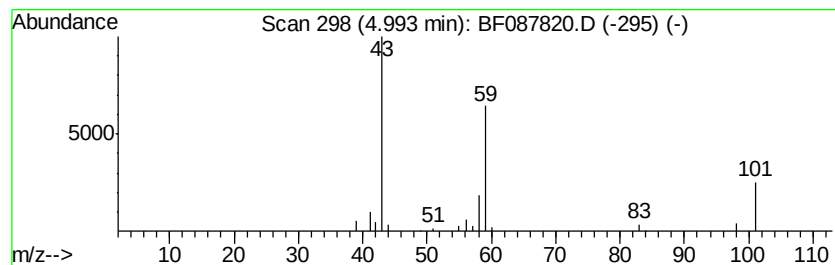
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.99	2.81 ng	83226	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	25
3		5-Hexen-2-one	98	C6H10O	000109-49-9	22
4		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9
5		3-Hexanol	102	C6H14O	000623-37-0	9



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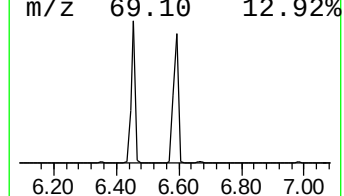
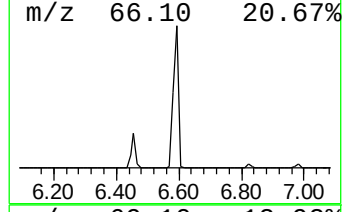
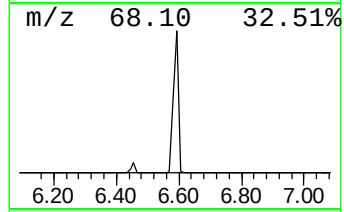
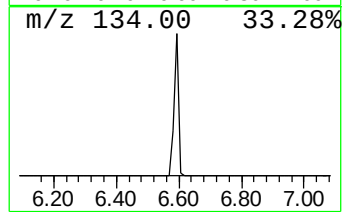
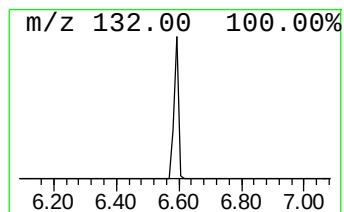
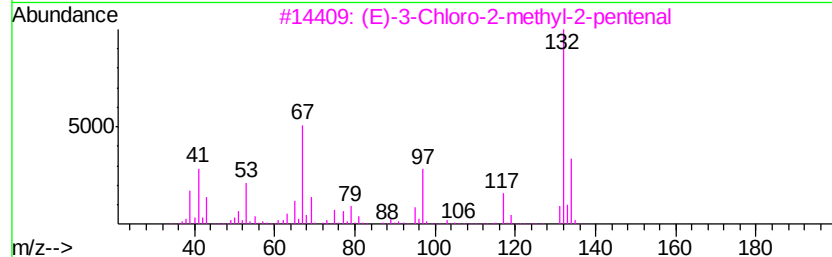
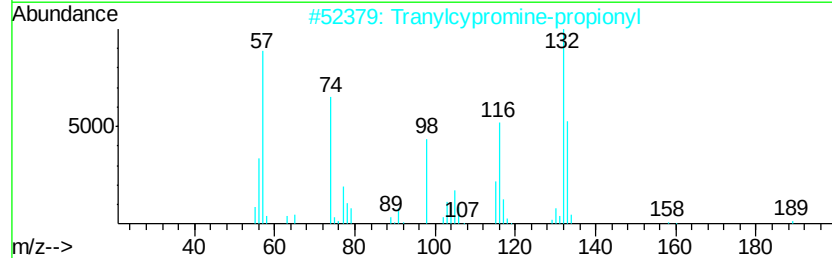
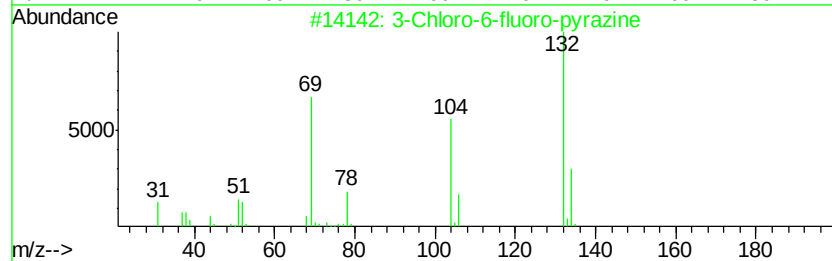
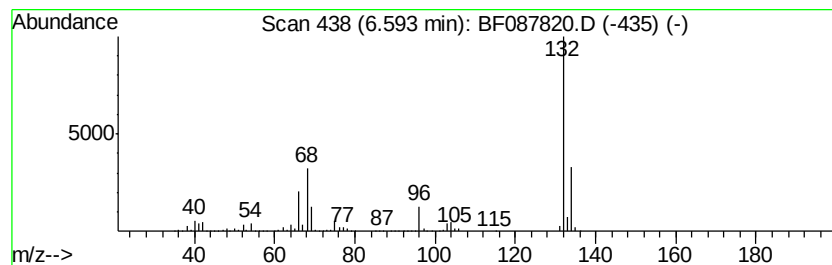
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.59 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	33.79 ng	1000700	1,4-Dichlorobenzene-d4	6.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	25
2	Tranylcypromine-propionyl			189	C12H15NO	1000123-86-3	12
3	(E)-3-Chloro-2-methyl-2-pentenal			132	C6H9ClO	031357-76-3	12
4	Xenon			132	Xe	007440-63-3	9
5	4-Chloro-6-fluoro-pyrimidine			132	C4H2ClFN2	051422-01-6	9



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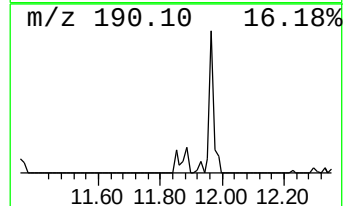
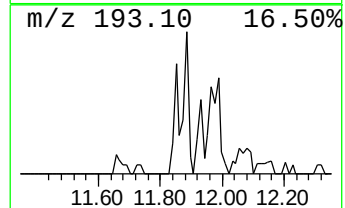
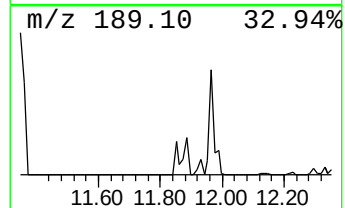
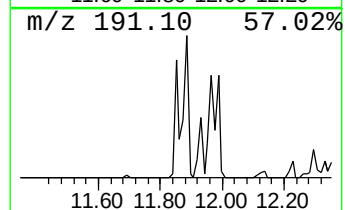
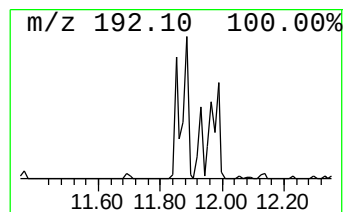
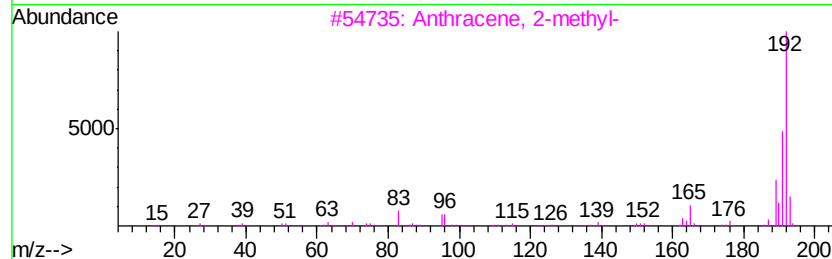
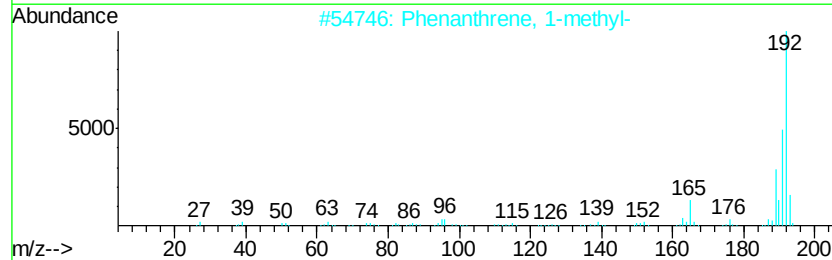
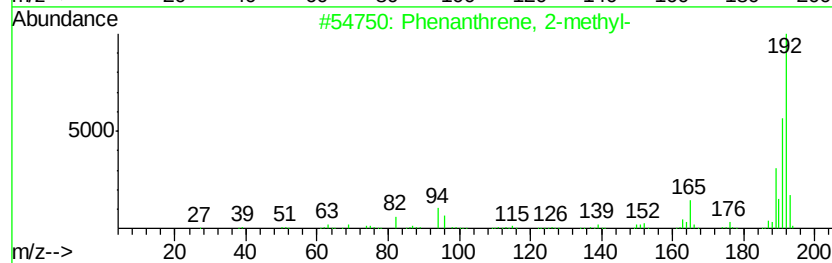
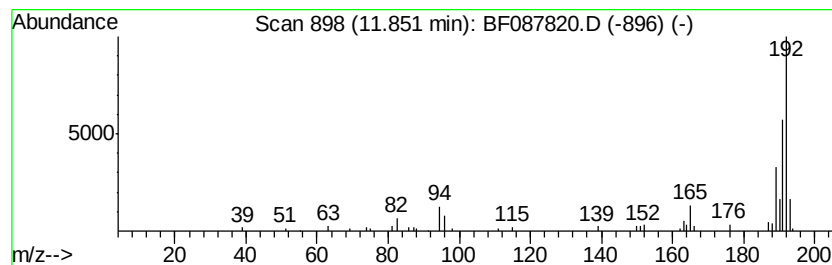
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TIC Library : C:\Database\NIST11.L
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Peak Number 7 Phenanthrene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD		R.T.
11.85	2.17 ng	69852	Phenanthrene-d10		11.35
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4	1H-Cyclopropa[1,1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	95



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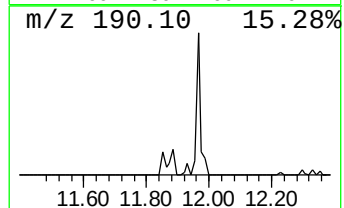
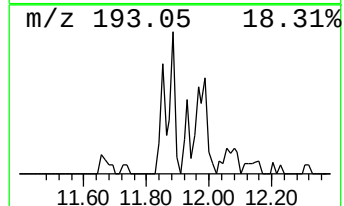
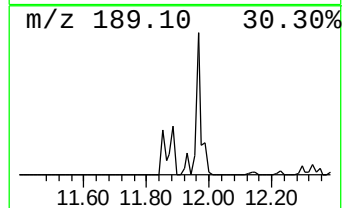
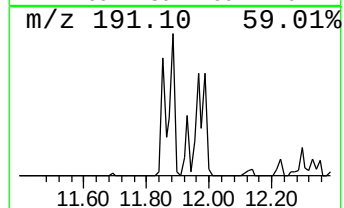
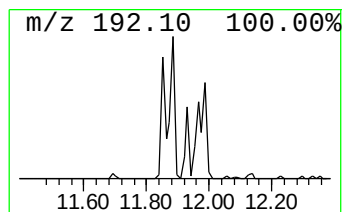
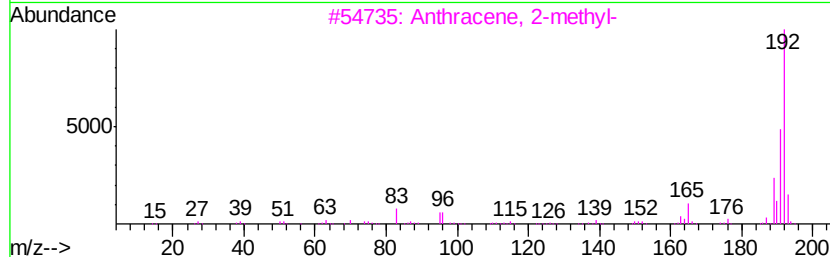
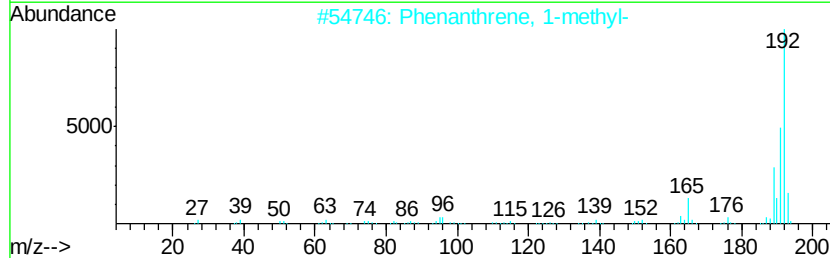
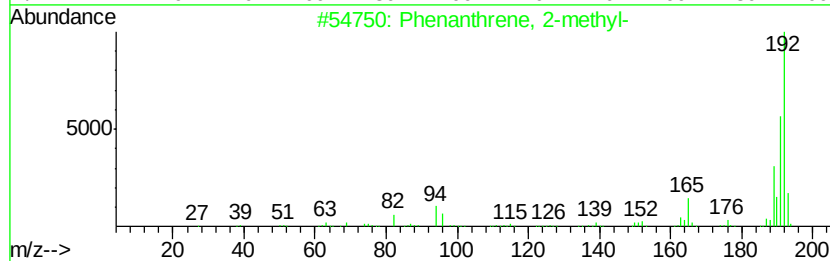
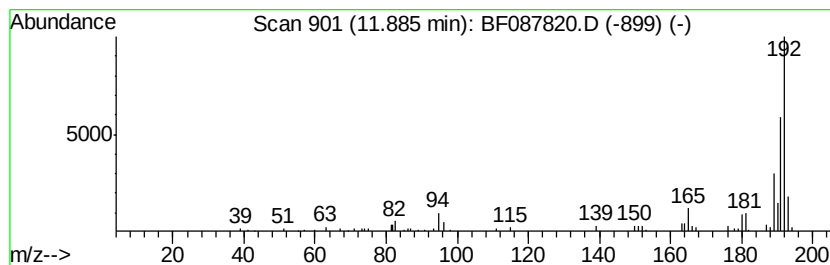
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Peak Number 8 Phenanthrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.89	3.07 ng	98958	Phenanthrene-d10	11.35	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	89
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	86



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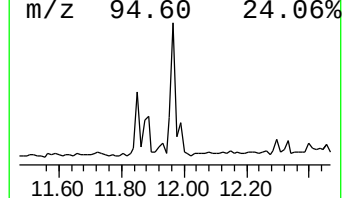
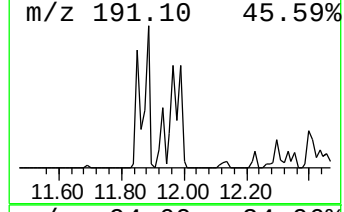
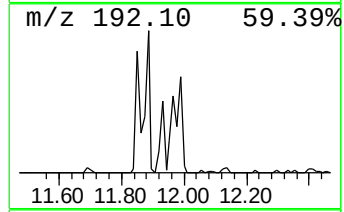
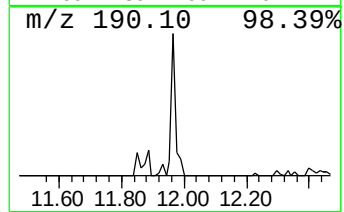
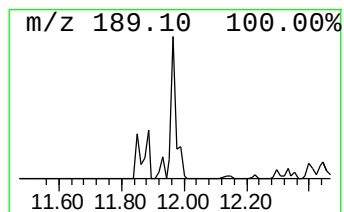
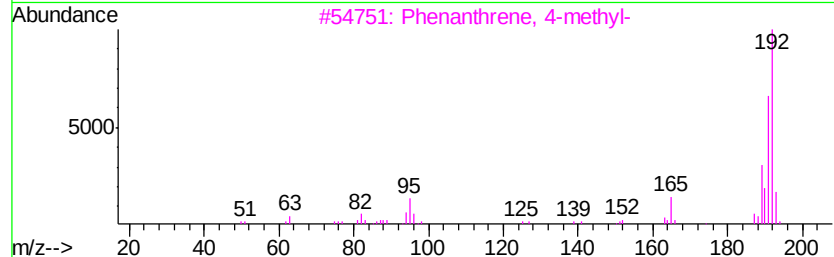
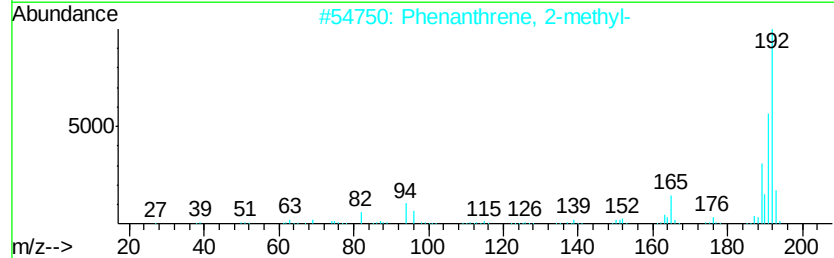
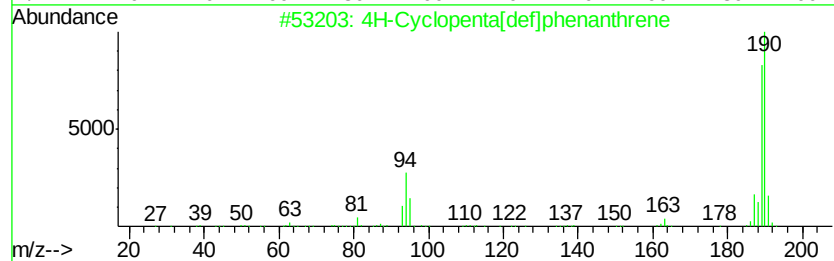
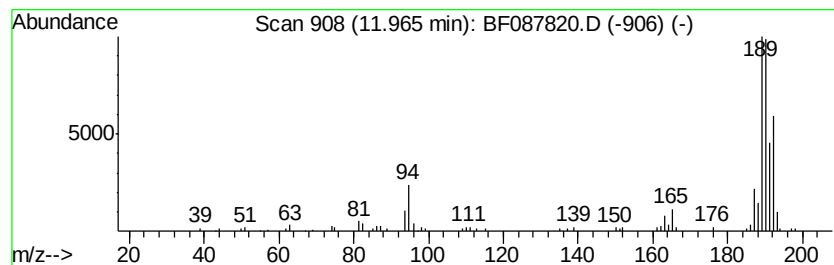
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.97	5.24 ng	168956	Phenanthrene-d10	11.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	92
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	38
3	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	38
4	Methyl diselenide	190	C2H6Se2	007101-31-7	38
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060816\
Data File : BF087820.D
Acq On : 8 Jun 2016 20:19
Operator : UM/SJ
Sample : H3379-15 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SS-10

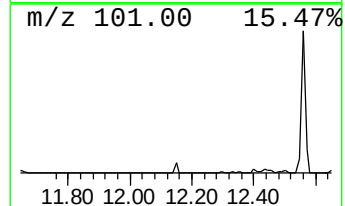
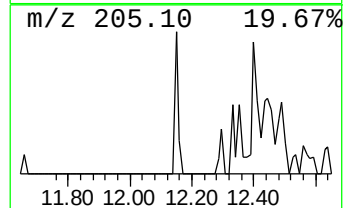
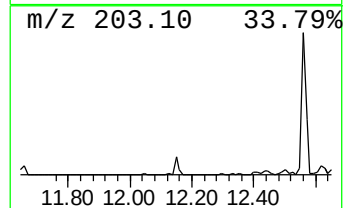
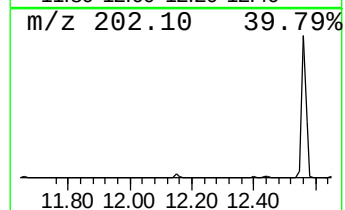
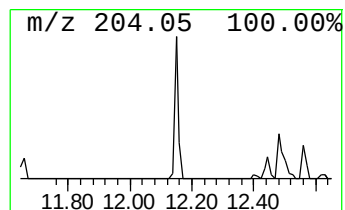
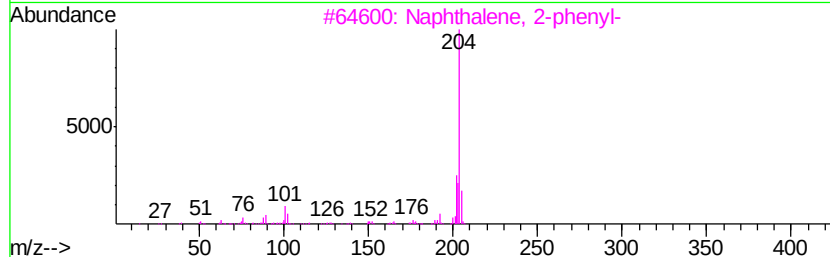
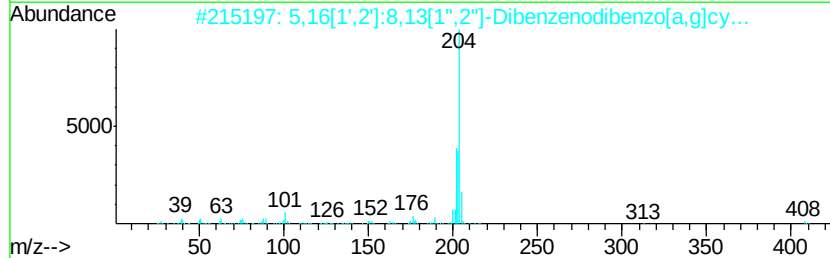
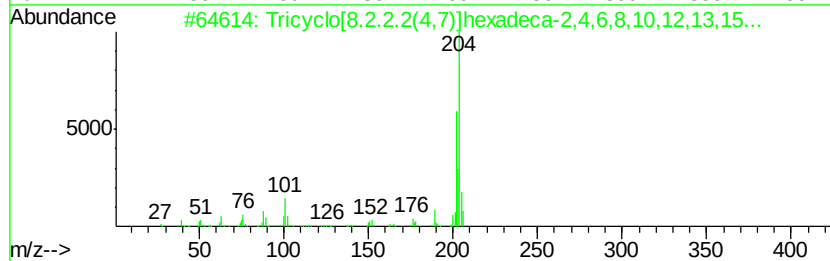
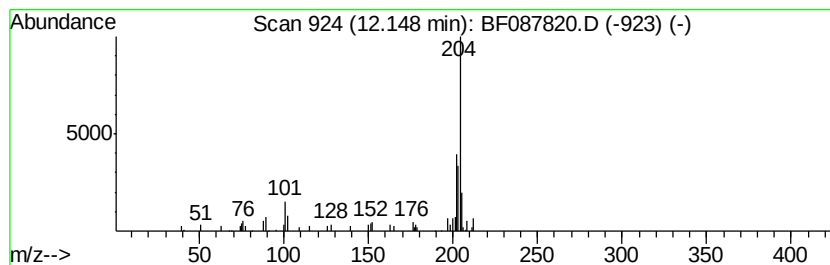
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Tricyclo[8.2.2.2(4,7)]hexad... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	2.03 ng	65462	Phenanthrene-d10	11.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	86
2		5,16[1',2'][:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	86
3		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	76
4		9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	68
5		6-Cyanophenanthridine	204	C14H8N2	002176-28-5	55



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060816\
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Acq On : 8 Jun 2016 20:19
Operator : UM/SJ
Sample : H3379-15 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

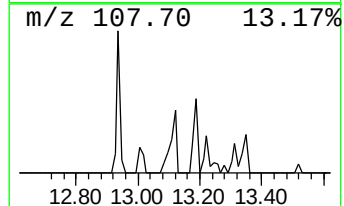
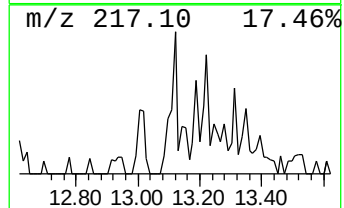
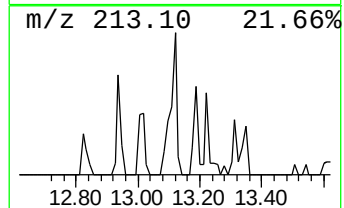
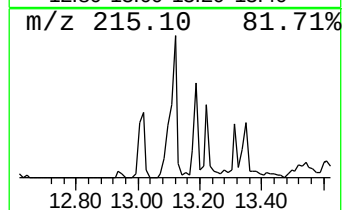
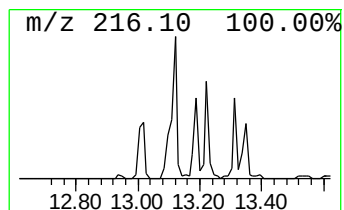
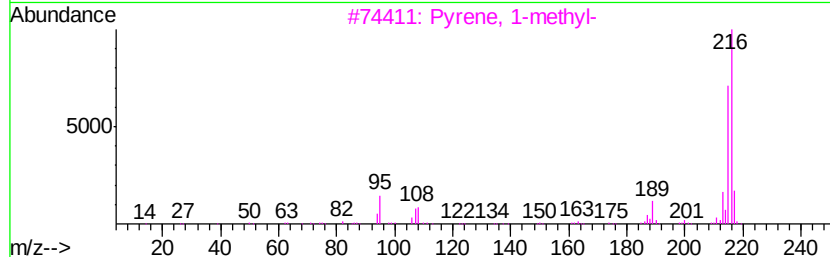
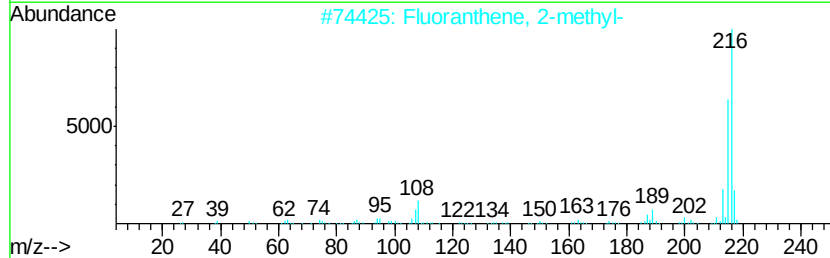
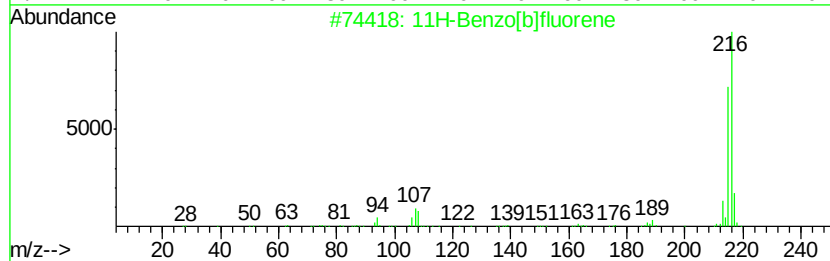
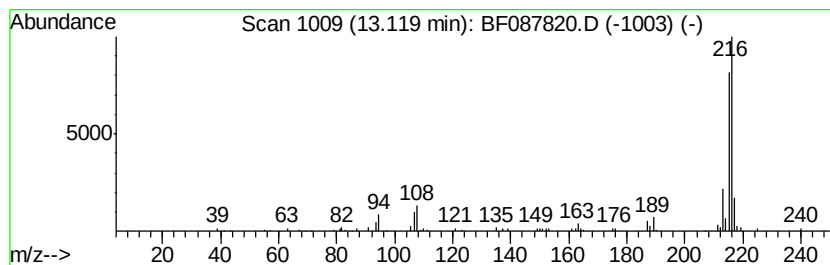
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 11H-Benzo[b]fluorene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.12	2.04 ng	120660	Chrysene-d12	13.99	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	92
5	Pyrene, 2-methyl-	216	C17H12	003442-78-2	86



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060816\
Data File : BF087820.D
Acq On : 8 Jun 2016 20:19
Operator : UM/SJ
Sample : H3379-15 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SS-10

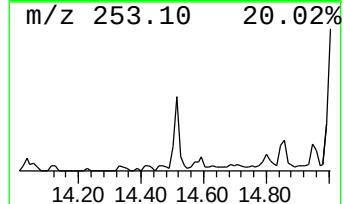
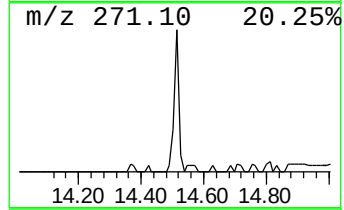
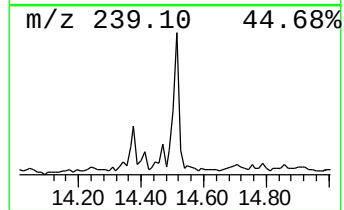
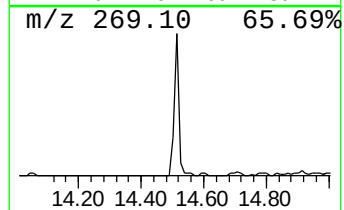
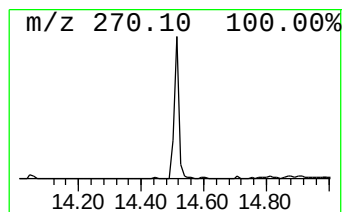
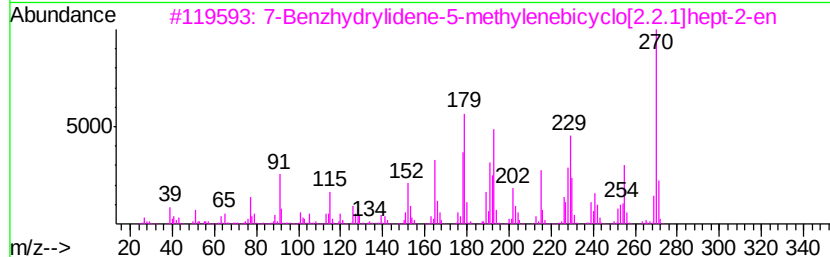
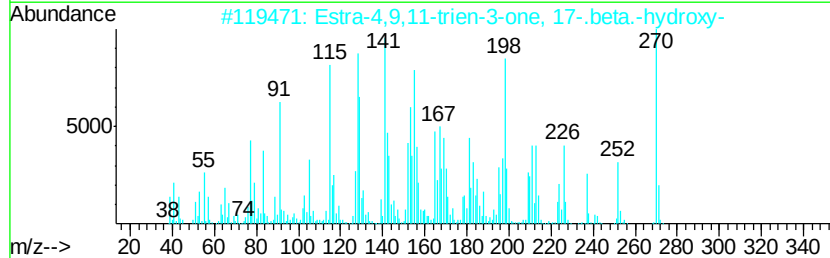
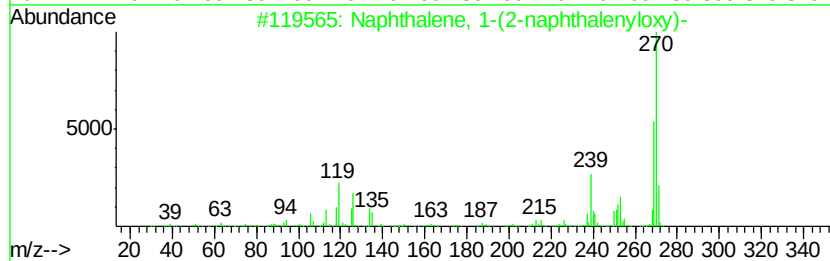
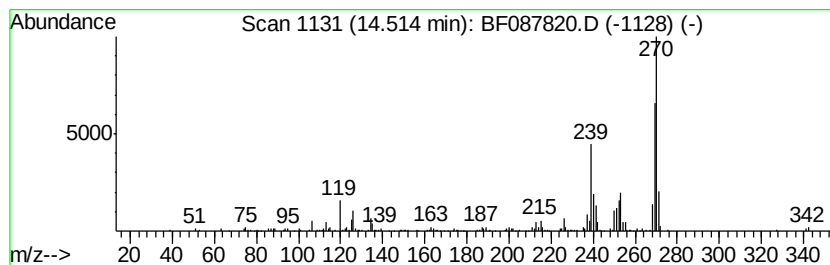
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Naphthalene, 1-(2-naphthale... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.51	3.78 ng	223109	Chrysene-d12	13.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-(2-naphthalenyloxy)-	270	C20H14O	000611-49-4	97
2		Estra-4,9,11-trien-3-one, 17-.be...	270	C18H22O2	010161-33-8	91
3		7-Benzhydrylidene-5-methylenebic...	270	C21H18	094371-60-5	83
4		9-Benzylidenexanthene	270	C20H14O	027980-52-5	53
5		Benz[a]anthracene-7-carboxaldehy...	270	C20H14O	013345-61-4	52



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060816\
Data File : BF087820.D
Acq On : 8 Jun 2016 20:19
Operator : UM/SJ
Sample : H3379-15 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampled :
SS-10

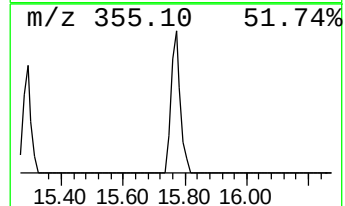
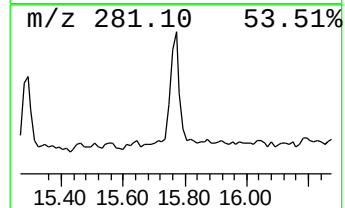
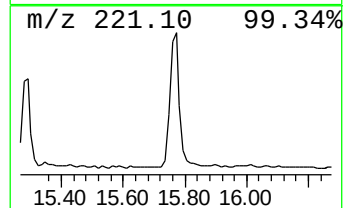
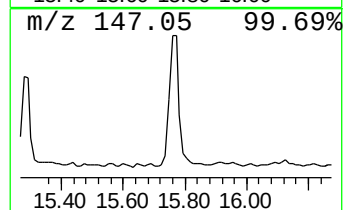
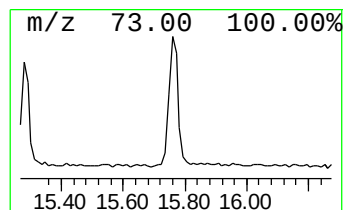
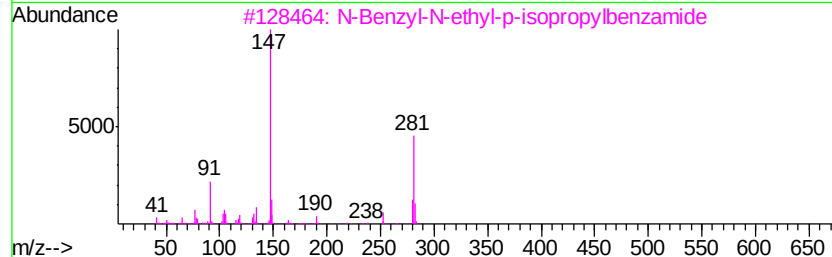
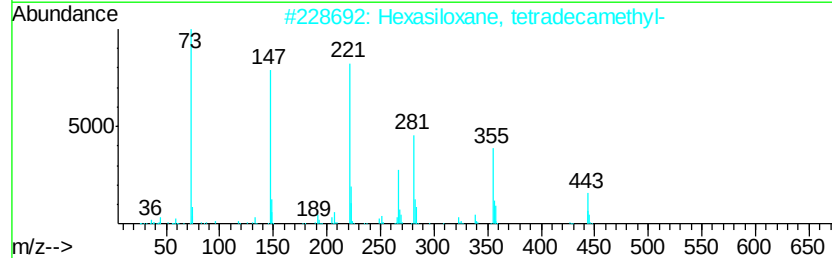
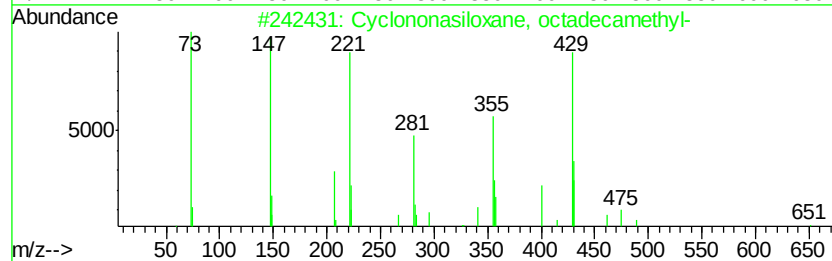
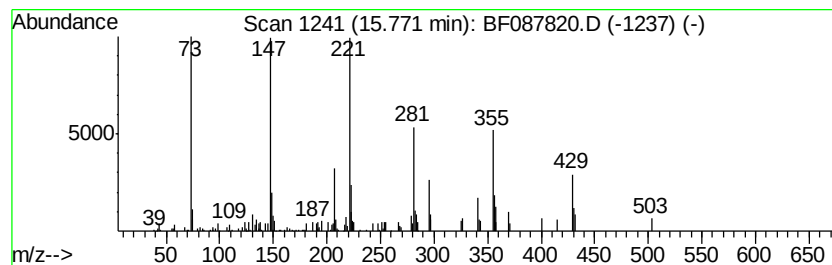
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Cyclononasiloxane, octadeca... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.77	4.51 ng	150723	Perylene-d12	15.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclononasiloxane, octadecamethyl-	666	C18H5409Si9	000556-71-8	87
2			Hexasiloxane, tetradecamethyl-	458	C14H4205Si6	000107-52-8	59
3			N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	22
4			4-Hydroxybenzyl alcohol, bis(ter...	352	C19H3602Si2	1000364-43-9	16
5			Heptasiloxane, hexadecamethyl-	532	C16H4806Si7	000541-01-5	14



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060816\
Data File : BF087820.D
Acq On : 8 Jun 2016 20:19
Operator : UM/SJ
Sample : H3379-15 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SS-10

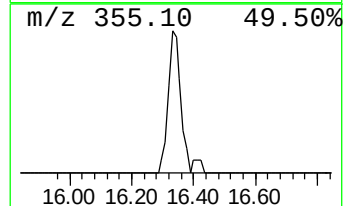
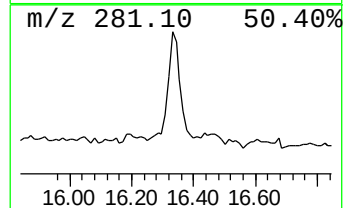
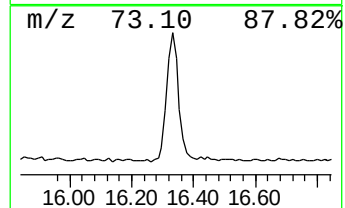
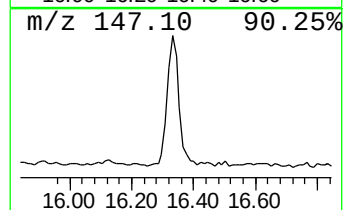
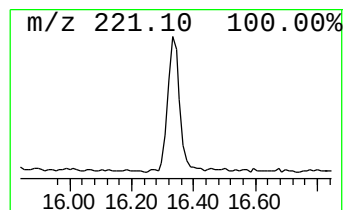
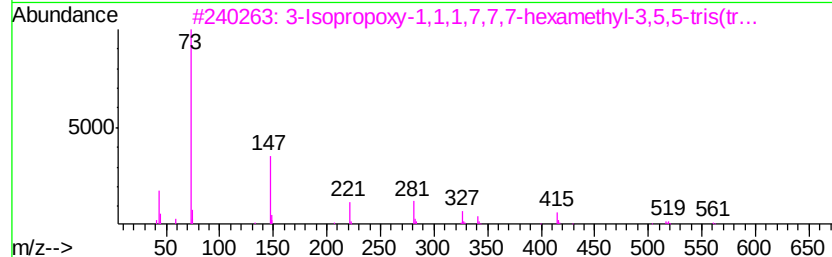
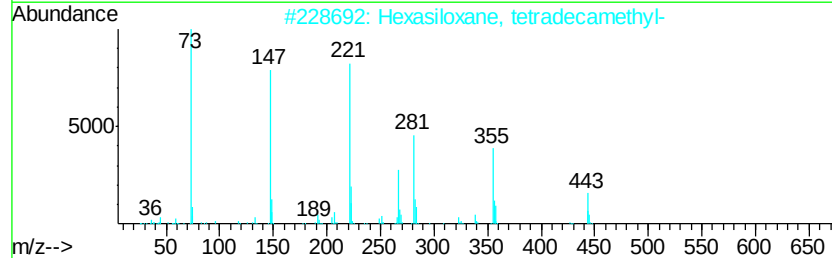
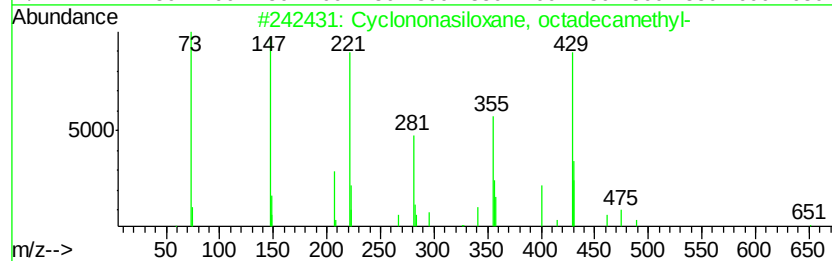
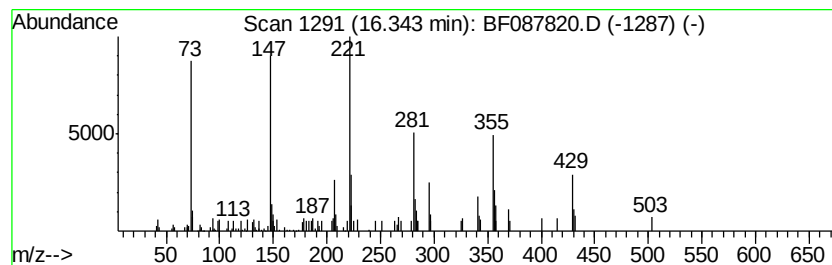
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Hexasiloxane, tetradecamethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.34	5.11 ng	170723	Perylene-d12	15.42

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	86
2			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	68
3			3-Isopropoxy-1,1,1,7,7,7-hexamet...	576	C18H52O7Si7	071579-69-6	32
4			N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	25
5			Pentasiloxane, dodecamethyl-	384	C12H36O4Si5	000141-63-9	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060816\
Data File : BF087820.D
Acq On : 8 Jun 2016 20:19
Operator : UM/SJ
Sample : H3379-15 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SS-10

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown6.59	6.59	33.8	ng	1000700	1	6.82	592223	20.0
Phenanthrene, 2-m...	11.85	2.2	ng	69852	4	11.35	645284	20.0
Phenanthrene, 1-m...	11.89	3.1	ng	98958	4	11.35	645284	20.0
4H-Cyclopenta[def...	11.97	5.2	ng	168956	4	11.35	645284	20.0
Tricyclo[8.2.2.2(...	12.15	2.0	ng	65462	4	11.35	645284	20.0
11H-Benzo[b]fluorene	13.12	2.0	ng	120660	5	13.99	1181570	20.0
Naphthalene, 1-(2...	14.51	3.8	ng	223109	5	13.99	1181570	20.0
Cyclononasiloxane...	15.77	4.5	ng	150723	6	15.42	668728	20.0
Hexasiloxane, tet...	16.34	5.1	ng	170723	6	15.42	668728	20.0
Piperidine, 1-(5-...	17.03	4.4	ng	148138	6	15.42	668728	20.0