

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
 Data File : BF086123.D
 Acq On : 7 Apr 2016 3:13
 Operator : UM/SJ
 Sample : H2191-02 5X
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-B1COMP

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-BF040216.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	2.304	17	19	27	rVB	42723	71131	4.96%	0.321%
2	4.087	173	175	177	rBV2	217809	420688	29.35%	1.897%
3	4.121	177	178	181	rVV3	209595	453586	31.65%	2.046%
4	4.167	181	182	183	rVV	195749	244926	17.09%	1.105%
5	4.201	183	185	187	rVB2	221949	337948	23.58%	1.524%
6	4.247	187	189	190	rBV	67400	90578	6.32%	0.409%
7	4.316	190	195	197	rVV5	109672	332212	23.18%	1.498%
8	4.361	197	199	200	rVV	68834	88698	6.19%	0.400%
9	4.407	202	203	205	rVV	40392	63861	4.46%	0.288%
10	4.464	205	208	211	rVV2	99973	257823	17.99%	1.163%
11	4.510	211	212	213	rVV	95795	98744	6.89%	0.445%
12	4.544	213	215	217	rVB3	109641	185575	12.95%	0.837%
13	5.333	283	284	293	rBV	379233	555516	38.76%	2.505%
14	5.676	312	314	317	rVB2	963804	1433317	100.00%	6.464%
15	6.396	375	377	383	rVV	399255	541666	37.79%	2.443%
16	6.510	386	387	390	rVV	399128	581888	40.60%	2.624%
17	6.567	390	392	395	rVB4	35781	53853	3.76%	0.243%
18	6.739	405	407	411	rVV	524055	629576	43.92%	2.839%
19	6.796	411	412	413	rVV	55612	53311	3.72%	0.240%
20	6.830	413	415	416	rVV2	76025	78539	5.48%	0.354%
21	6.899	416	421	423	rVV	540694	598084	41.73%	2.697%
22	7.013	426	431	434	rVV3	63999	142696	9.96%	0.644%
23	7.070	434	436	437	rVV2	59117	72189	5.04%	0.326%
24	7.184	443	446	448	rVB3	26211	64814	4.52%	0.292%
25	7.310	455	457	461	rVB	324831	313585	21.88%	1.414%
26	7.390	463	464	468	rBV4	51814	74650	5.21%	0.337%
27	7.493	471	473	475	rVB3	87732	106291	7.42%	0.479%
28	7.550	475	478	480	rVB4	62860	120925	8.44%	0.545%
29	7.607	480	483	485	rBV4	81570	104134	7.27%	0.470%
30	7.687	487	490	496	rBV4	40463	157289	10.97%	0.709%
31	7.859	502	505	507	rVB2	44239	75679	5.28%	0.341%
32	8.030	518	520	522	rBV	764985	661250	46.13%	2.982%
33	8.419	551	554	558	rVB6	29282	69856	4.87%	0.315%
34	9.105	612	614	616	rBV	717173	585820	40.87%	2.642%

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	9.242	624	626	630	rVB5	54476	54673	3.81%	0.247%
36	9.505	648	649	653	rVB	69725	73546	5.13%	0.332%
37	9.642	659	661	664	rBV	46462	60652	4.23%	0.274%
38	9.779	671	673	675	rVV	616089	586497	40.92%	2.645%
39	9.813	675	676	678	rVB	68857	59931	4.18%	0.270%
40	9.985	689	691	693	rVV	43257	54462	3.80%	0.246%
41	10.190	705	709	710	rBV3	23561	58187	4.06%	0.262%
42	10.328	720	721	723	rBV	85450	78706	5.49%	0.355%
43	10.568	740	742	745	rBV	263539	288742	20.15%	1.302%
44	10.865	766	768	774	rBV5	42149	135494	9.45%	0.611%
45	11.036	780	783	786	rVB5	39195	74260	5.18%	0.335%
46	11.265	801	803	804	rBV	672738	599941	41.86%	2.706%
47	11.345	807	810	811	rBV	99426	99158	6.92%	0.447%
48	11.505	821	824	828	rBV	54428	76284	5.32%	0.344%
49	11.791	844	849	851	rBV	126411	270806	18.89%	1.221%
50	11.871	855	856	861	rVB	99314	184562	12.88%	0.832%
51	12.065	870	873	878	rVB3	56114	162972	11.37%	0.735%
52	12.202	882	885	887	rBV4	42496	110323	7.70%	0.498%
53	12.316	892	895	897	rBV	63924	110536	7.71%	0.499%
54	12.351	897	898	901	rVB2	72084	87446	6.10%	0.394%
55	12.408	901	903	906	rBV	42852	68416	4.77%	0.309%
56	12.476	907	909	911	rBV	640694	530703	37.03%	2.393%
57	12.511	911	912	914	rVB2	89460	88107	6.15%	0.397%
58	12.625	920	922	924	rBV	32672	54719	3.82%	0.247%
59	12.705	926	929	931	rBV	581627	653234	45.57%	2.946%
60	12.762	932	934	937	rVB	50272	78423	5.47%	0.354%
61	12.842	939	941	945	rVB	521589	582309	40.63%	2.626%
62	12.934	945	949	952	rBV4	49311	153377	10.70%	0.692%
63	13.036	954	958	960	rBV2	77607	131426	9.17%	0.593%
64	13.082	960	962	965	rBV2	57714	131731	9.19%	0.594%
65	13.139	965	967	970	rVB2	48481	68998	4.81%	0.311%
66	13.231	973	975	976	rBV	40246	55254	3.85%	0.249%
67	13.254	976	977	979	rVB	390130	284159	19.83%	1.282%
68	13.322	981	983	985	rBV	380754	326526	22.78%	1.473%
69	13.425	990	992	993	rBV	53032	55401	3.87%	0.250%
70	13.551	1001	1003	1004	rBV	53934	77920	5.44%	0.351%
71	13.677	1010	1014	1015	rBV2	78749	156164	10.90%	0.704%

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Sampling : 1 Min Area: 3 % of largest Peak
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72	13.699	1015	1016	1019	rVV2	94822	136012	9.49%	0.613%
73	13.757	1019	1021	1024	rVB2	81885	124056	8.66%	0.559%
74	13.905	1029	1034	1035	rBV2	585050	1156768	80.71%	5.217%
75	14.065	1046	1048	1050	rBV2	92102	96896	6.76%	0.437%
76	14.248	1059	1064	1066	rBV6	73891	179391	12.52%	0.809%
77	14.374	1072	1075	1077	rBV3	68582	113366	7.91%	0.511%
78	14.419	1077	1079	1081	rBV3	121958	202743	14.15%	0.914%
79	14.648	1096	1099	1103	rBV2	132734	306377	21.38%	1.382%
80	14.911	1120	1122	1126	rBV	156986	252174	17.59%	1.137%
81	15.037	1130	1133	1135	rBV	60534	174024	12.14%	0.785%
82	15.185	1145	1146	1150	rVB	106582	162862	11.36%	0.734%
83	15.242	1150	1151	1154	rVV	108183	166800	11.64%	0.752%
84	15.300	1154	1156	1164	rVV	335265	614157	42.85%	2.770%
85	15.437	1167	1168	1170	rVB	194239	160963	11.23%	0.726%
86	15.711	1189	1192	1193	rBV	64913	126782	8.85%	0.572%
87	16.648	1272	1274	1282	rVB2	47833	140929	9.83%	0.636%
88	16.877	1292	1294	1298	rVB4	44312	59833	4.17%	0.270%
89	17.700	1364	1366	1368	rBV2	53661	115474	8.06%	0.521%
90	18.271	1413	1416	1417	rBV3	66236	85388	5.96%	0.385%
91	18.317	1417	1420	1425	rVB6	70815	245621	17.14%	1.108%
92	18.408	1425	1428	1429	rBV2	69941	157379	10.98%	0.710%
93	18.488	1433	1435	1438	rVB3	79814	147891	10.32%	0.667%
94	18.603	1443	1445	1450	rBV5	63753	179742	12.54%	0.811%
95	18.683	1450	1452	1454	rBV3	56259	110203	7.69%	0.497%
96	18.763	1456	1459	1461	rBV3	75488	170126	11.87%	0.767%
97	18.968	1475	1477	1478	rBV	64879	86892	6.06%	0.392%
98	19.174	1493	1495	1497	rBV2	51696	67666	4.72%	0.305%
99	19.357	1509	1511	1512	rBV2	49114	66753	4.66%	0.301%
100	19.414	1512	1516	1518	rVV5	36067	121175	8.45%	0.546%

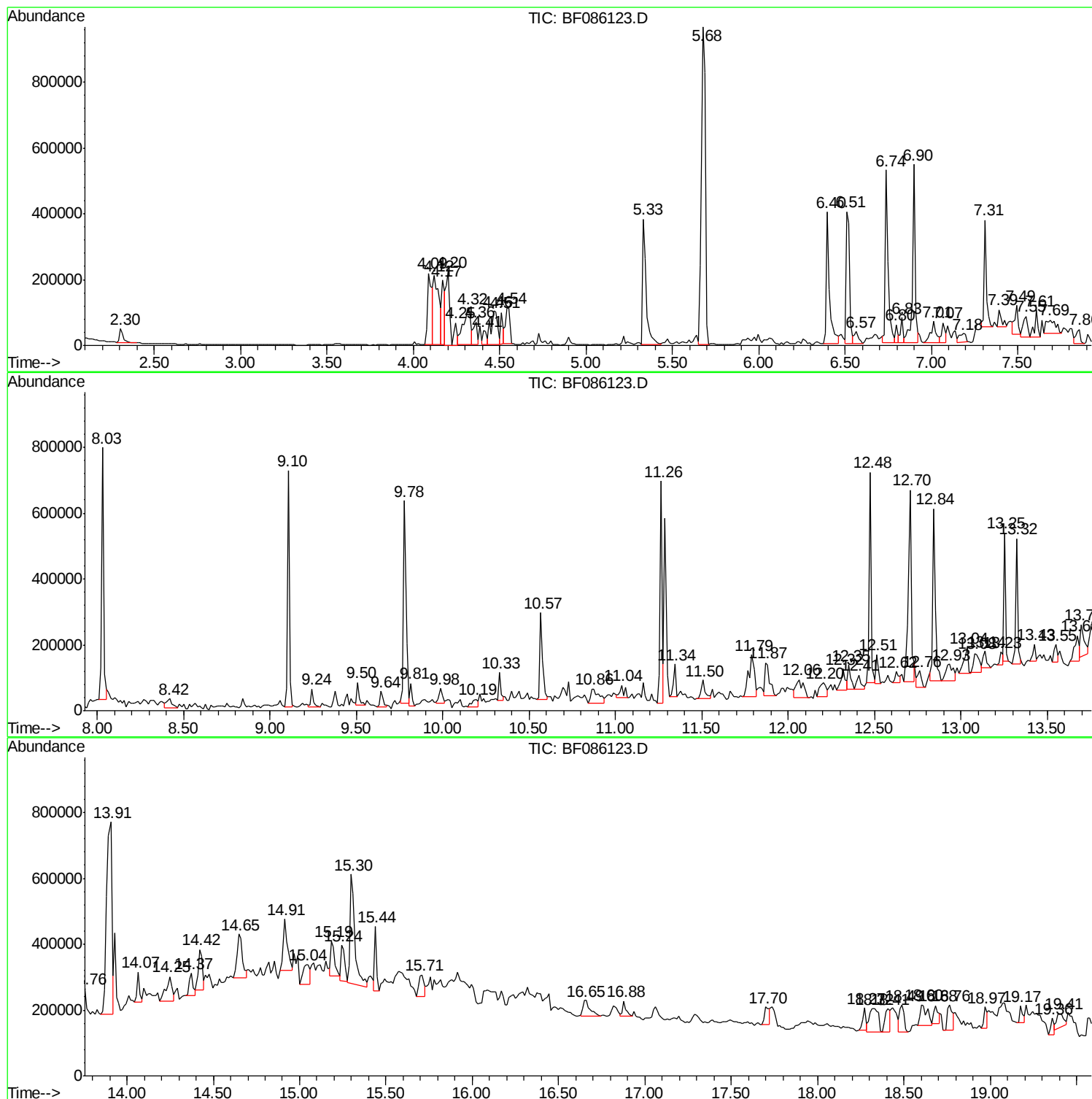
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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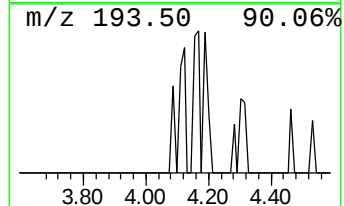
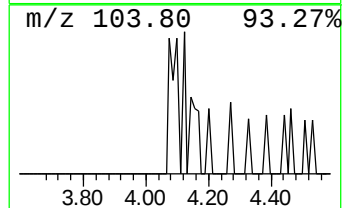
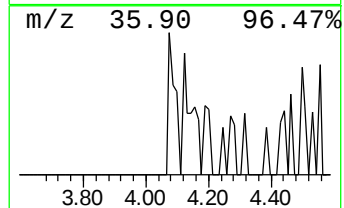
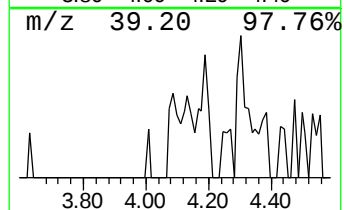
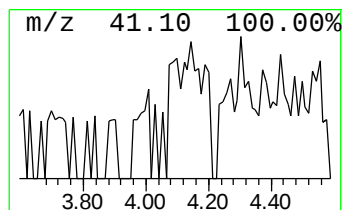
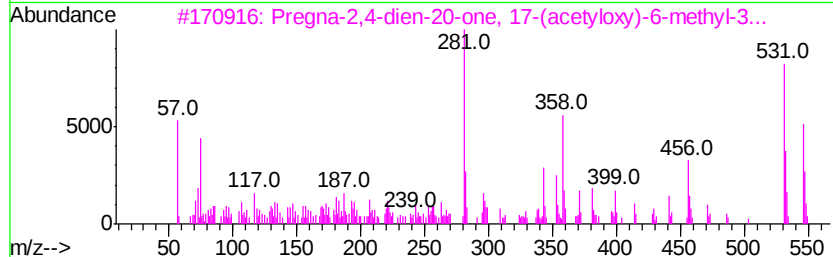
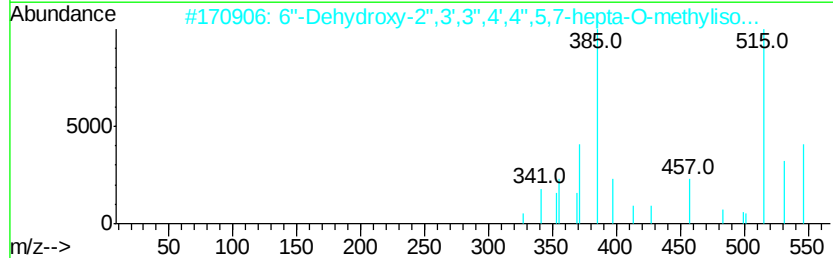
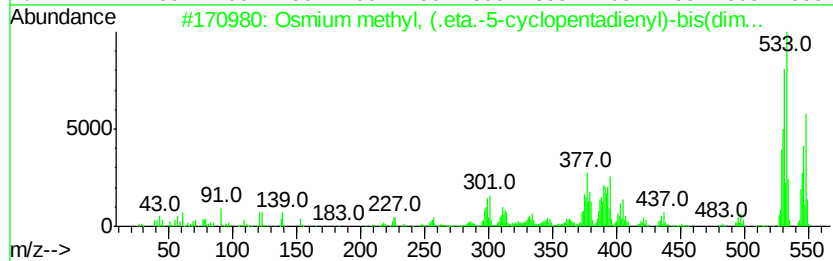
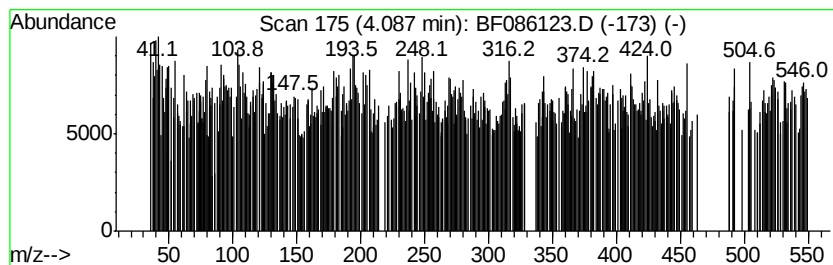
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown4.09 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.09	13.36 ng	420688	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Osmium methyl, (.eta.-5-cyclopenten...	548	C22H300sP2	1000158-92-4	20
2		6"-Dehydroxy-2",3',3",4',4",5,7-...	546	C28H34011	1000064-97-6	15
3		Predna-2,4-dien-20-one, 17-(acet...	546	C30H5005Si2	069833-57-4	15
4		Heptamethylethyloctasilsesquioxane	550	C9H26012Si8	077626-17-6	14
5		3ALPHA-METHOXY-3BETA-(METHANESUL...	521	C30H51N04S	1000241-18-6	11



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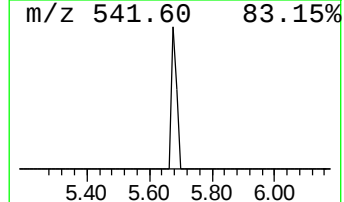
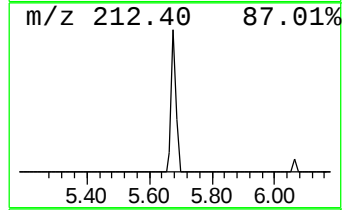
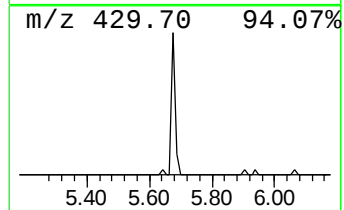
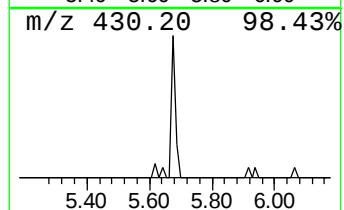
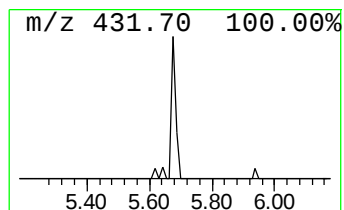
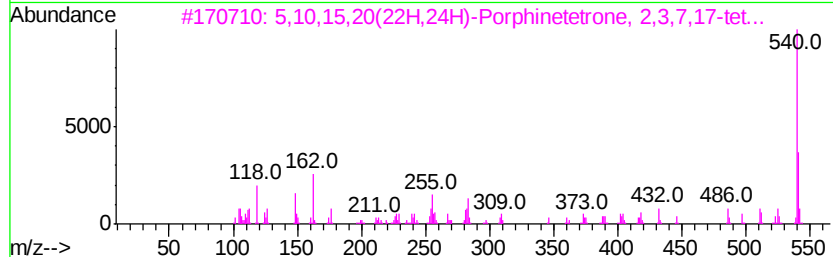
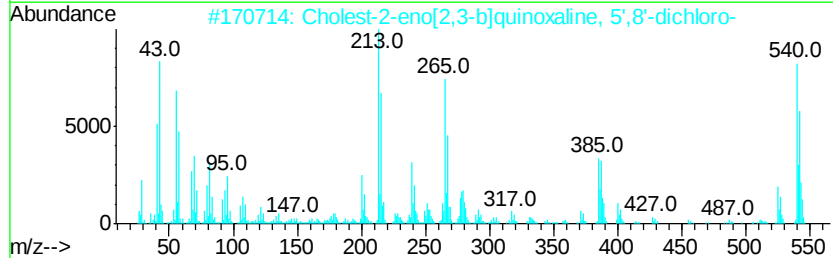
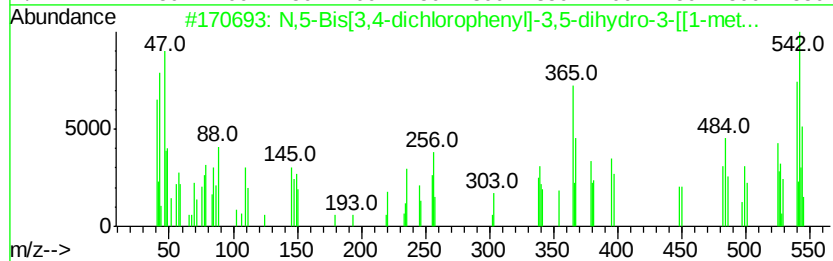
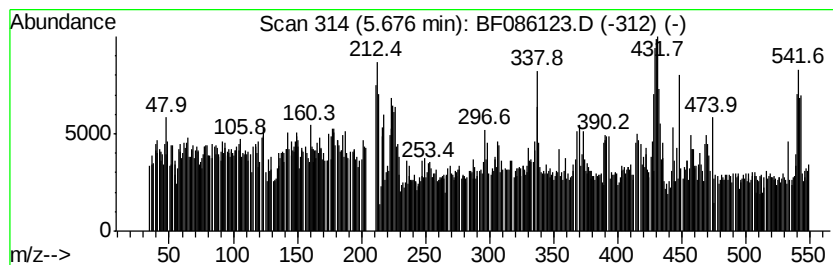
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 N,5-Bis[3,4-dichlorophenyl]... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.68	45.53 ng	1433320	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	N,5-Bis[3,4-dichlorophenyl]-3,5-...	540	C27H20Cl4N4	333399-40-9	56
2		Cholest-2-eno[2,3-b]quinoxaline,...	540	C33H46Cl2N2	1000210-84-2	38
3		5,10,15,20(22H,24H)-Porphinetetr...	540	C32H36N4O4	027800-02-8	25
4		Distannoxane, hexabutyl-	598	C24H54O2Sn2	000056-35-9	12
5		Nickel[II] bis(0,0'-dibutyldithi...	540	C16H36NiO4P2S4	037913-22-7	10



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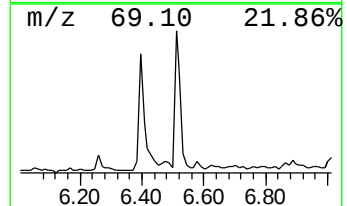
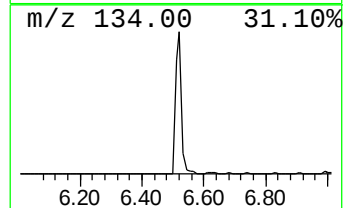
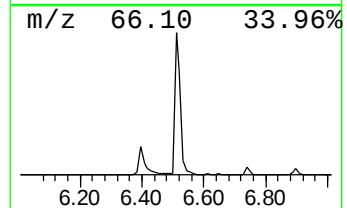
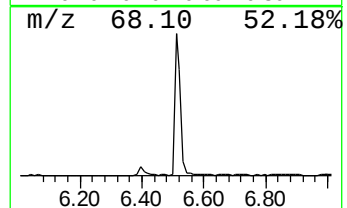
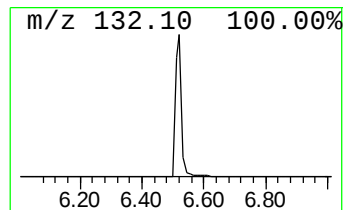
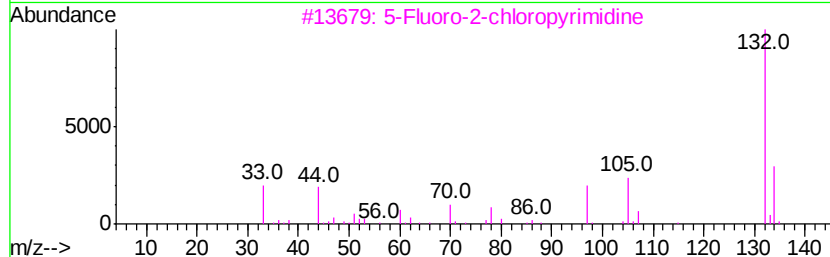
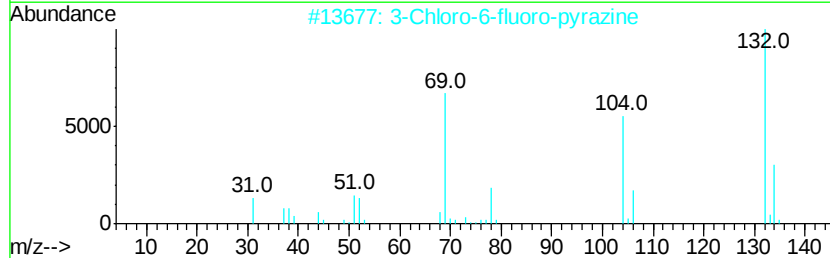
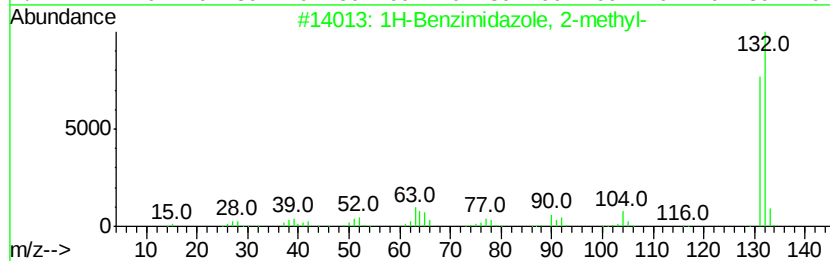
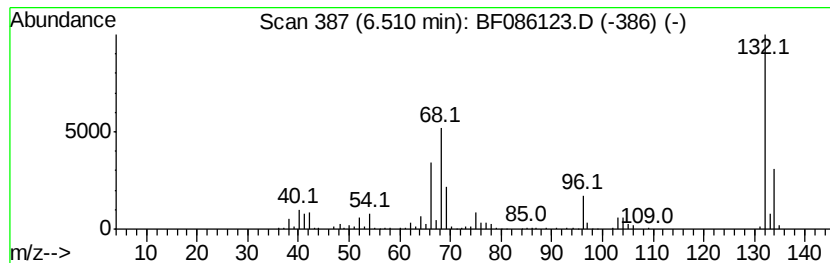
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown6.51 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	18.49 ng	581888	1,4-Dichlorobenzene-d4	6.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086123.D
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Operator : UM/SJ
Sample : H2191-02 5X
Misc :
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ClientSampleId :
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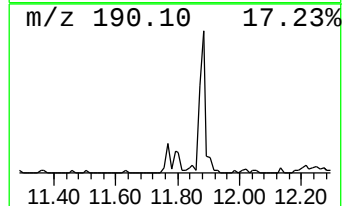
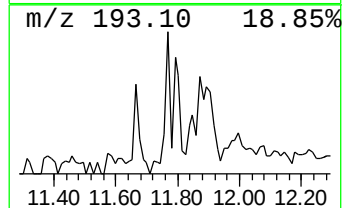
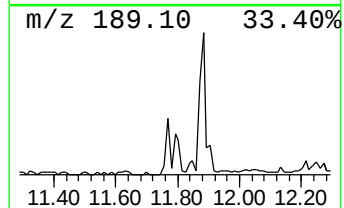
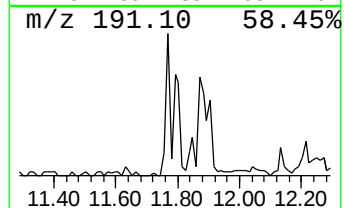
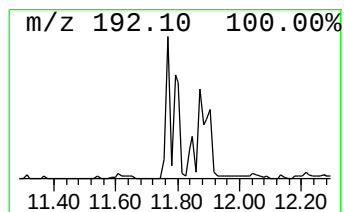
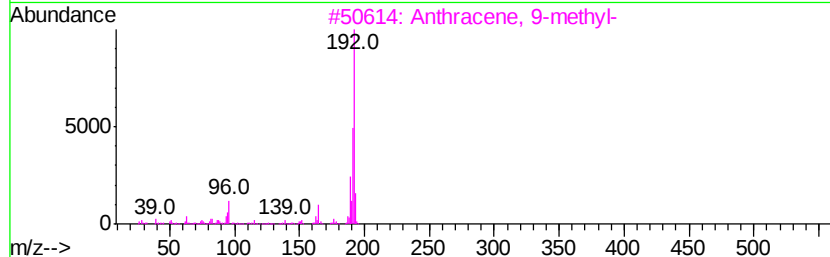
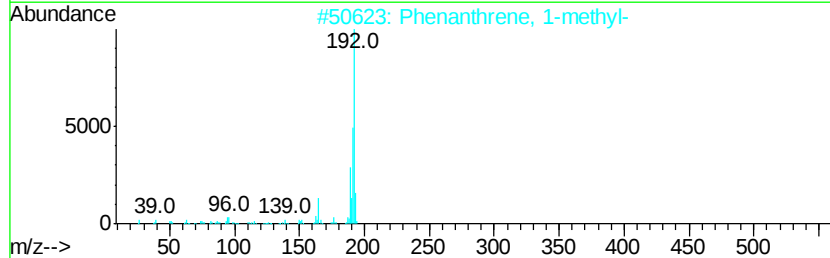
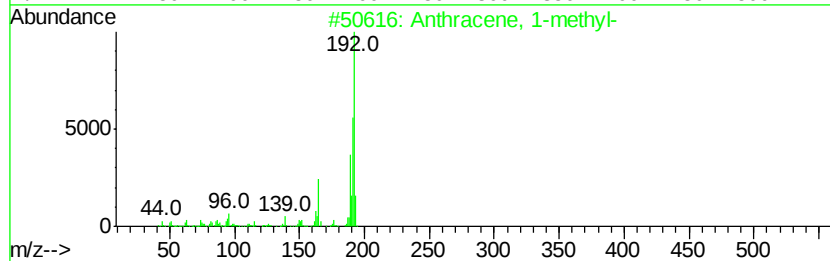
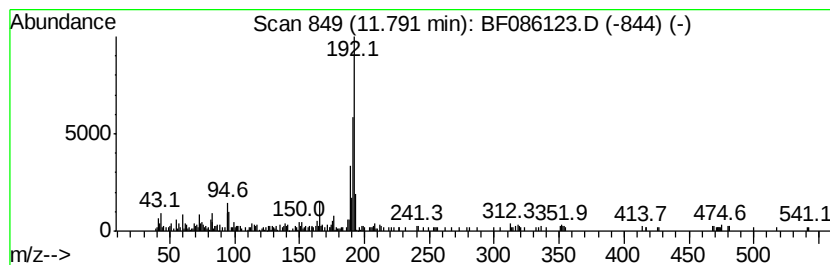
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Anthracene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	9.03 ng	270806	Phenanthrene-d10	11.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	86
3		Anthracene, 9-methyl-	192	C15H12	000779-02-2	86
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	86
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086123.D
Acq On : 7 Apr 2016 3:13
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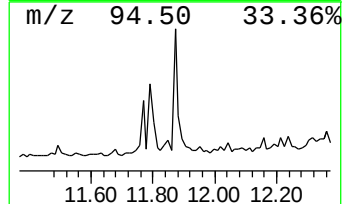
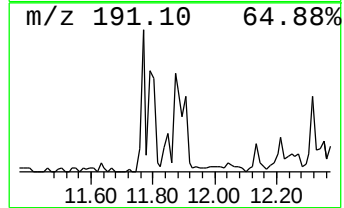
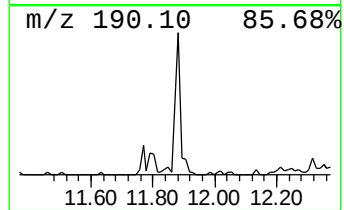
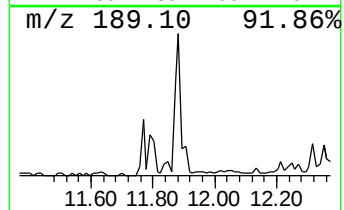
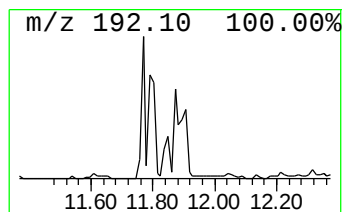
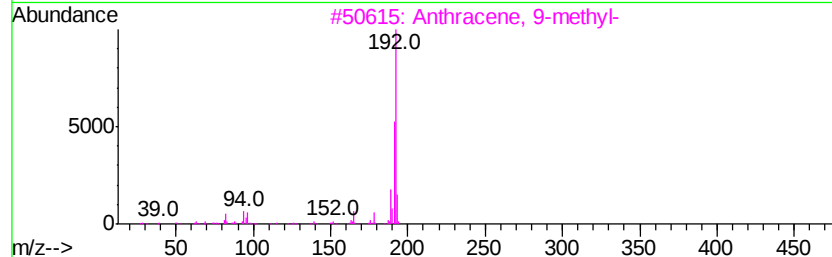
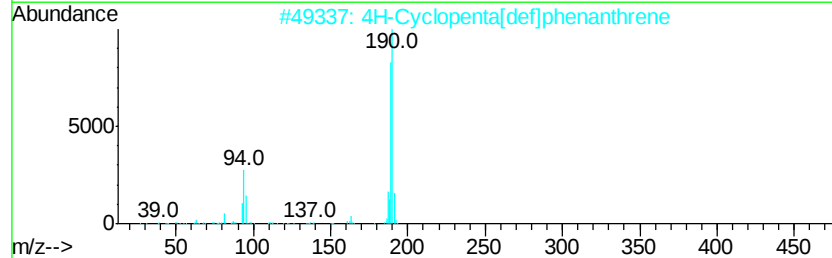
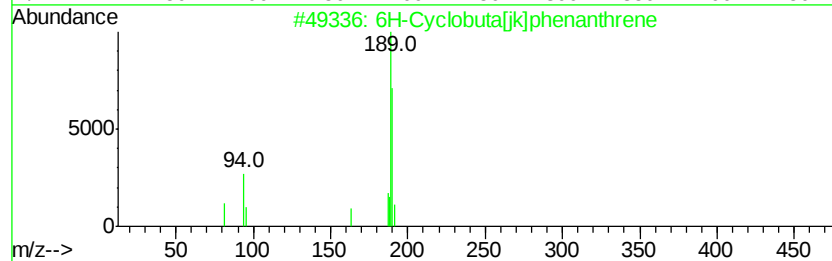
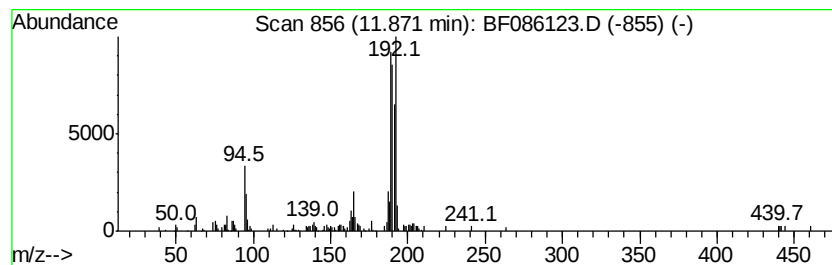
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown11.87 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.87	6.15 ng	184562	Phenanthrene-d10	11.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	49
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	46
4			Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42
5			Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086123.D
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Operator : UM/SJ
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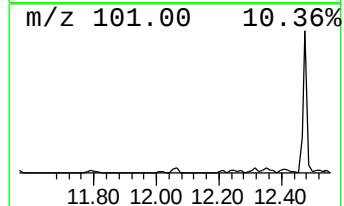
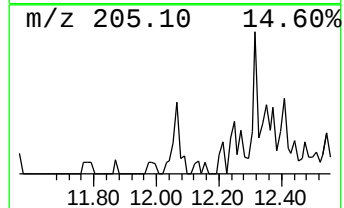
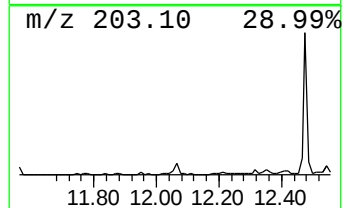
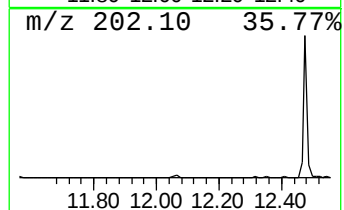
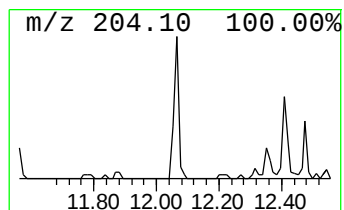
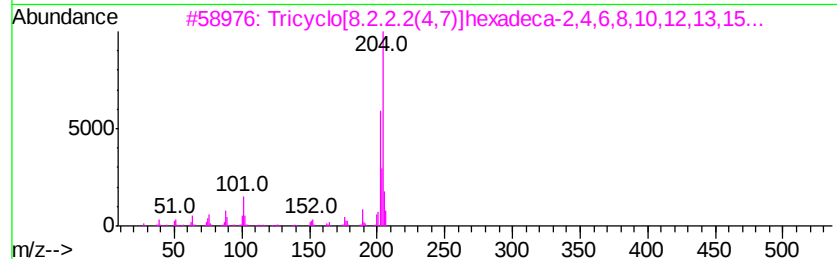
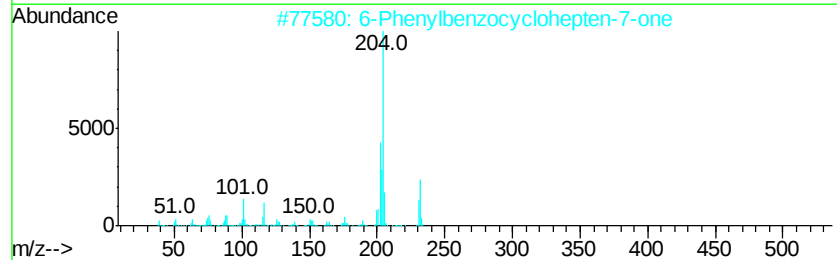
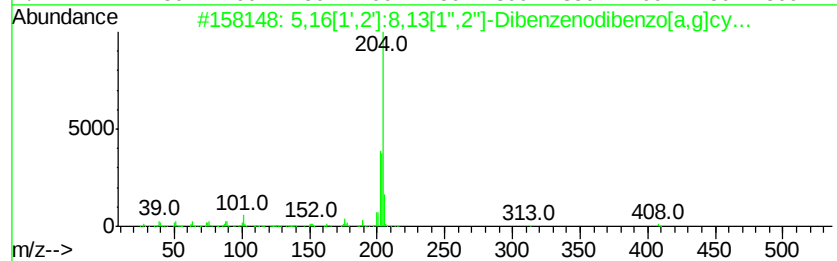
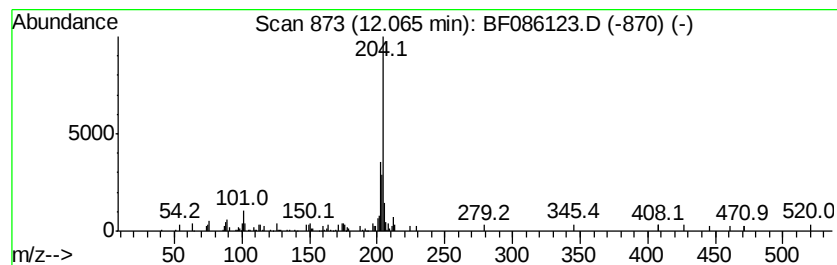
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 5,16[1',2']:8,13[1'',2'']-D... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.06	5.43 ng	162972	Phenanthrene-d10	11.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	91
2		6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	72
3		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	64
4		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	64
5		2-Phenyl-naphthalene	204	C16H12	035465-71-5	52



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040616\
Data File : BF086123.D
Acq On : 7 Apr 2016 3:13
Operator : UM/SJ
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Misc :
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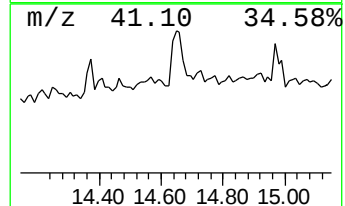
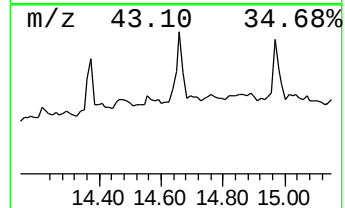
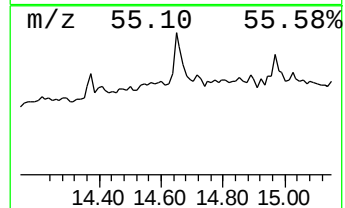
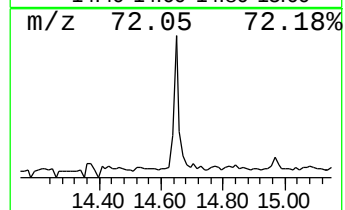
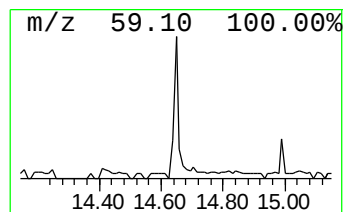
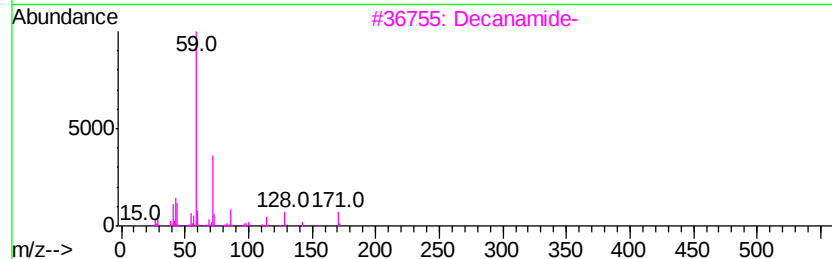
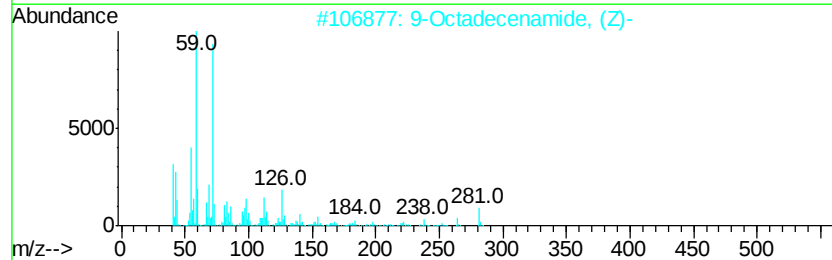
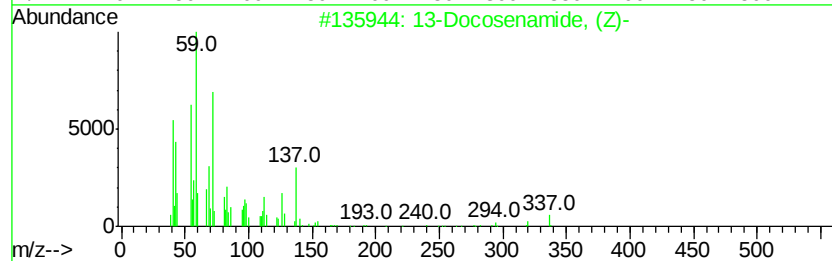
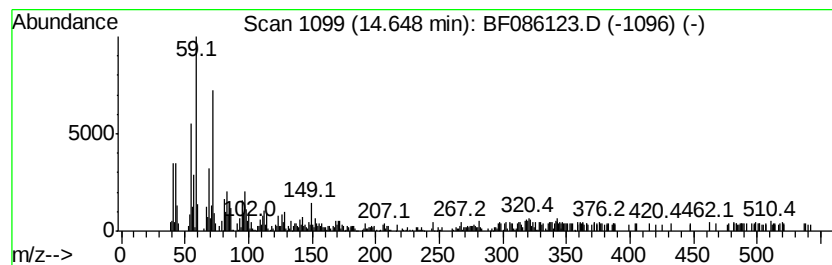
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 13-Docosenamide, (Z)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.65	9.98 ng	306377	Perylene-d12	15.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	78
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	60
3			Decanamide-	171	C10H21NO	002319-29-1	49
4			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	49
5			Cyclohexane, (ethoxymethoxy)-	158	C9H18O2	054699-29-5	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
Data File : BF086123.D
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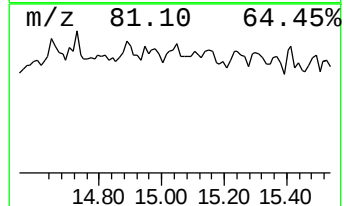
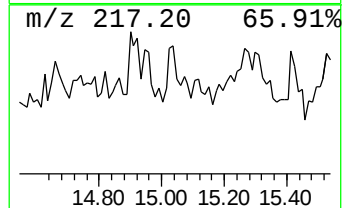
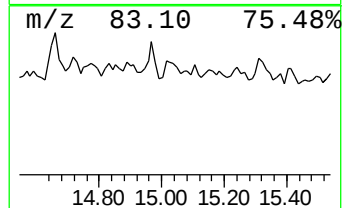
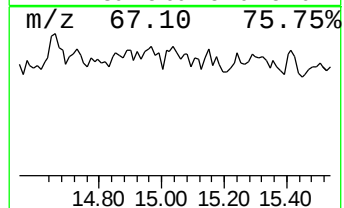
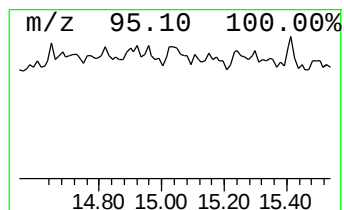
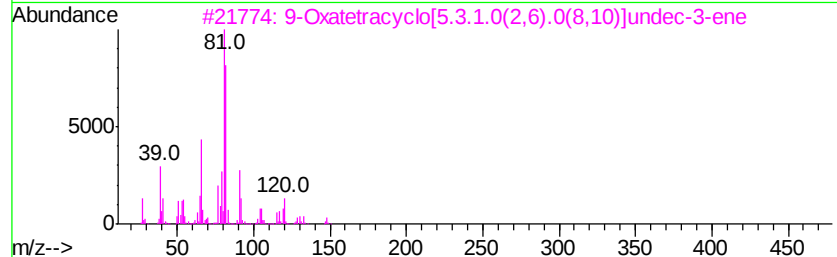
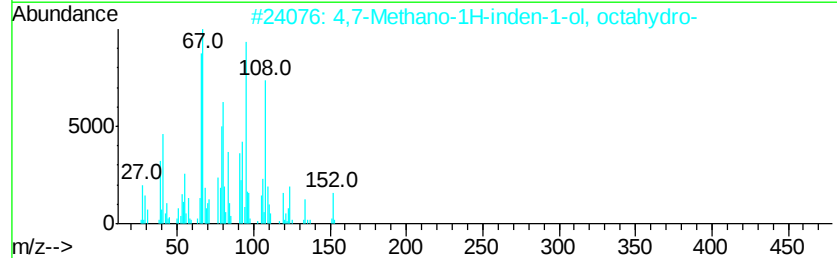
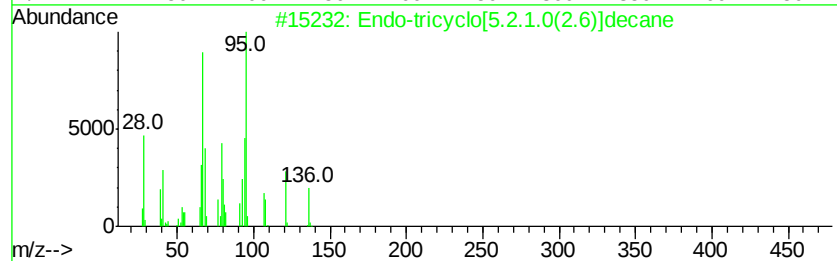
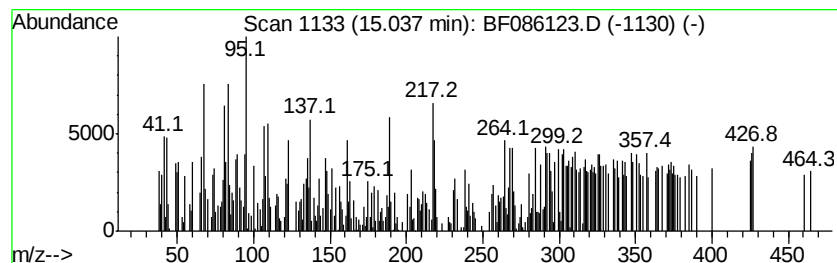
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 unknown15.04 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.04	5.67 ng	174024	Perylene-d12	15.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Endo-tricyclo[5.2.1.0(2.6)]decane	136	C10H16	1000215-28-8	25
2		4,7-Methano-1H-inden-1-ol, octah...	152	C10H16O	055255-97-5	22
3		9-Oxatetracyclo[5.3.1.0(2.6).0(8...	148	C10H12O	1000191-87-2	22
4		1,3-Bis(cyclopentyl)-1-cyclopentene	204	C15H24	1000215-55-1	22
5		Bicyclo[2.2.1]heptane, 2-(2-meth...	150	C11H18	061142-27-6	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040616\
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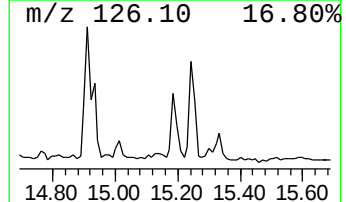
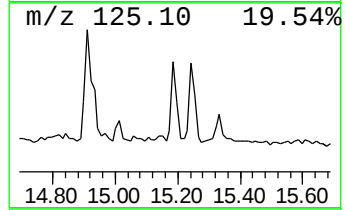
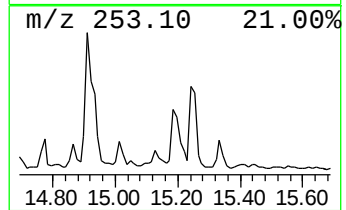
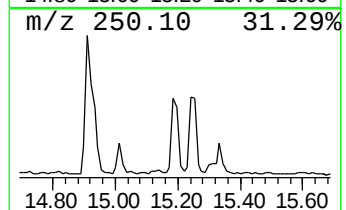
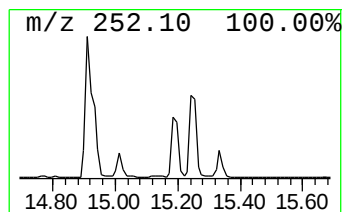
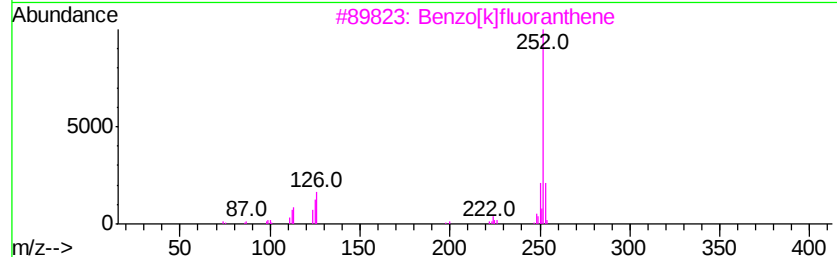
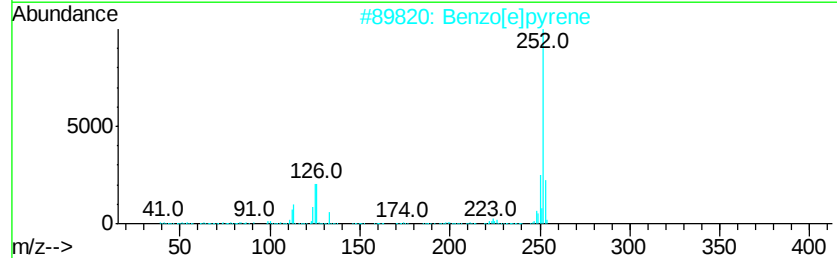
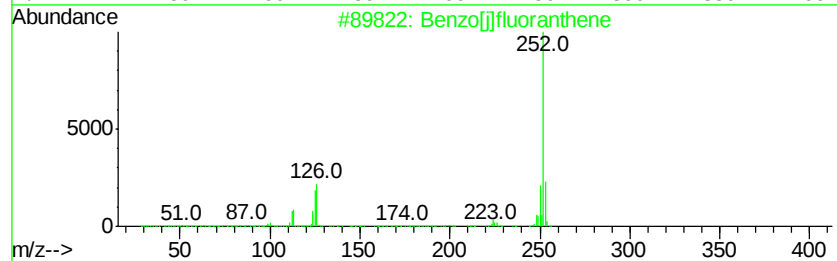
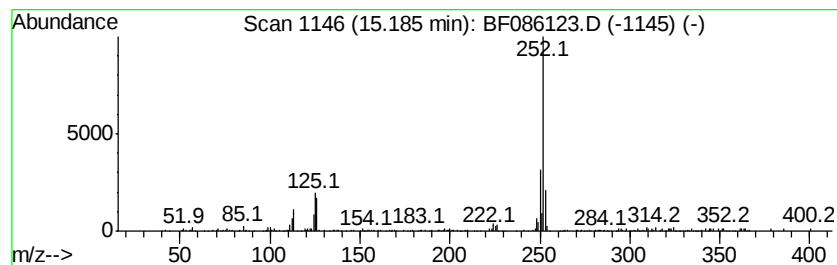
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 Benzo[j]fluoranthene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.19	5.30 ng	162862	Perylene-d12	15.30

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[i]fluoranthene	252	C20H12	000205-82-3	98
2		Benzo[e]pyrene	252	C20H12	000192-97-2	98
3		Benzo[k]fluoranthene	252	C20H12	000207-08-9	98
4		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	98
5		Perylene	252	C20H12	000198-55-0	95



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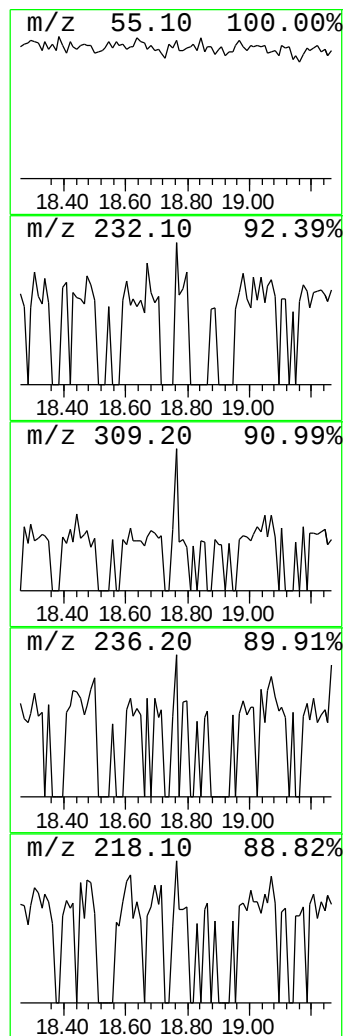
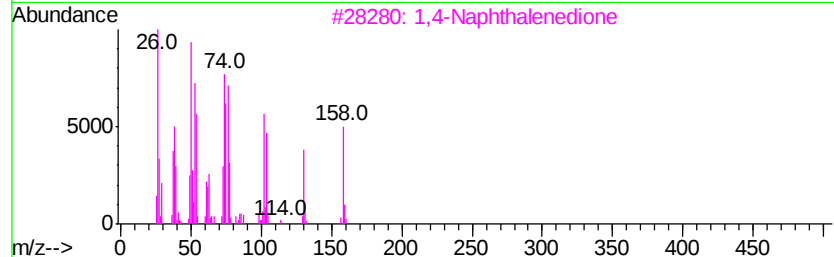
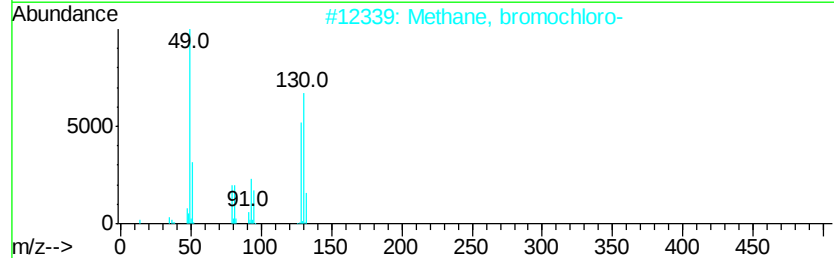
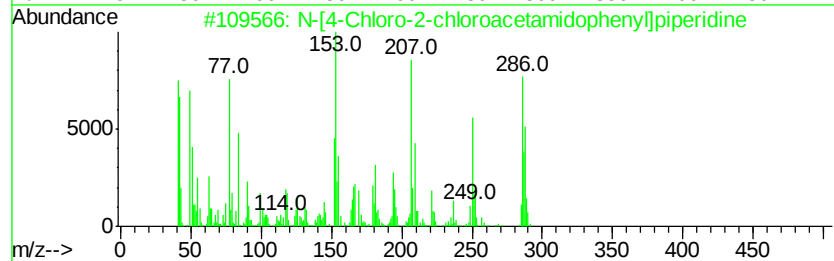
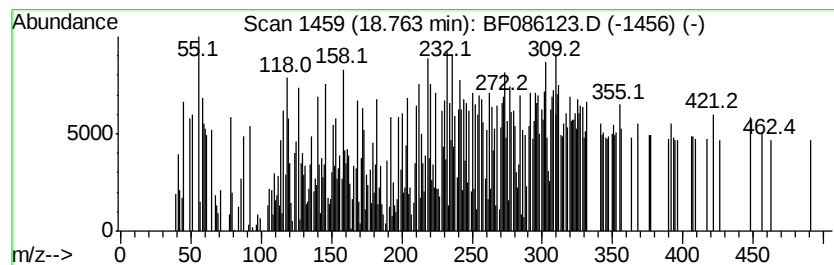
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 unknown18.76 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.76	5.54 ng	170126	Perylene-d12	15.30

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-[4-Chloro-2-chloroacetamidophe...	286	C13H16Cl12N2O	1000254-96-6	35
2			Methane, bromochloro-	128	CH2BrCl	000074-97-5	18
3			1,4-Naphthalenedione	158	C10H6O2	000130-15-4	14
4			Ethyl 2-(2-chloroacetamido)-2-(3...	390	C13H12Cl12F4N2O3	339352-71-5	10
5			1H-Quinoline-2(1H)-one, 3-bromom...	237	C10H8BrNO	035740-85-3	10



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040616\
 Data File : BF086123.D
 Acq On : 7 Apr 2016 3:13
 Operator : UM/SJ
 Sample : H2191-02 5X
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-B1COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.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
unknown4.09	4.09	13.4	ng	420688	1	6.74	629576	20.0
Decachlorodipheny...	4.17	7.8	ng	244926	1	6.74	629576	20.0
N,5-Bis[3,4-dichl...	5.68	45.5	ng	1433320	1	6.74	629576	20.0
unknown6.51	6.51	18.5	ng	581888	1	6.74	629576	20.0
Anthracene, 1-met...	11.79	9.0	ng	270806	4	11.26	599941	20.0
unknown11.87	11.87	6.2	ng	184562	4	11.26	599941	20.0
5,16[1',2']:8,13[...	12.06	5.4	ng	162972	4	11.26	599941	20.0
13-Docosenamide, ...	14.65	10.0	ng	306377	6	15.30	614157	20.0
unknown15.04	15.04	5.7	ng	174024	6	15.30	614157	20.0
Benzo[j]fluoranthene	15.19	5.3	ng	162862	6	15.30	614157	20.0
unknown18.76	18.76	5.5	ng	170126	6	15.30	614157	20.0