

Data Path : V:\HPCHEM1\BNA F\DATA\BF121515\
 Data File : BF083821.D
 Acq On : 15 Dec 2015 16:31
 Operator : UM/IZ
 Sample : G4739-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 K1206-0-2

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

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

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.093	260	263	266	rBV	509269	686263	11.98%	0.751%
2	5.516	297	300	302	rVB	2073745	2265473	39.56%	2.478%
3	5.733	316	319	324	rBV2	104216	193244	3.37%	0.211%
4	6.533	387	389	392	rVB	1629972	2255201	39.38%	2.467%
5	6.670	398	401	404	rBV	2192244	2105600	36.77%	2.303%
6	6.739	404	407	409	rVV3	166854	225696	3.94%	0.247%
7	6.899	419	421	424	rVV	643446	604806	10.56%	0.661%
8	7.059	433	435	438	rVV	1990678	1748479	30.54%	1.912%
9	7.459	467	470	473	rBV	1108173	1346007	23.51%	1.472%
10	8.190	531	534	538	rVV2	733607	1158276	20.23%	1.267%
11	8.899	594	596	598	rVB	465976	397255	6.94%	0.434%
12	9.002	603	605	606	rVV	229798	212178	3.71%	0.232%
13	9.265	625	628	630	rVB	2625362	2279648	39.81%	2.493%
14	9.528	648	651	653	rBV	117021	166446	2.91%	0.182%
15	9.608	656	658	661	rBV2	282045	453229	7.92%	0.496%
16	9.665	661	663	666	rVV2	455041	632485	11.05%	0.692%
17	9.802	672	675	678	rBV	1475167	1292979	22.58%	1.414%
18	9.939	684	687	689	rBV	884879	943228	16.47%	1.032%
19	10.248	712	714	716	rBV2	120494	203759	3.56%	0.223%
20	10.396	724	727	728	rVB2	141333	213234	3.72%	0.233%
21	10.488	733	735	738	rVB3	189056	300165	5.24%	0.328%
22	10.602	742	745	747	rBV2	168293	213805	3.73%	0.234%
23	10.648	747	749	753	rVB2	97522	198975	3.47%	0.218%
24	10.728	753	756	758	rVB	1299155	1259464	22.00%	1.377%
25	10.854	765	767	770	rBV	331111	422618	7.38%	0.462%
26	11.036	779	783	784	rBV2	117697	198624	3.47%	0.217%
27	11.185	792	796	798	rBV2	160837	258975	4.52%	0.283%
28	11.231	798	800	802	rVB	368271	368920	6.44%	0.403%
29	11.322	805	808	811	rVB	380990	496469	8.67%	0.543%
30	11.425	815	817	818	rBV	601135	708488	12.37%	0.775%
31	11.448	818	819	822	rVB	2043321	2721420	47.53%	2.976%
32	11.505	822	824	825	rBV	495701	616176	10.76%	0.674%
33	11.642	834	836	838	rBV2	117033	162090	2.83%	0.177%
34	11.722	838	843	844	rBV	475587	527398	9.21%	0.577%

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

35	11.756	844	846	848	rVB	547193	430216	7.51%	0.471%
36	11.836	848	853	855	rVB2	283062	469566	8.20%	0.514%
37	11.882	855	857	858	rBV	240166	237899	4.15%	0.260%
38	11.928	859	861	862	rVV	1179091	1085034	18.95%	1.187%
39	11.962	862	864	866	rVB	921855	1179746	20.60%	1.290%
40	12.042	869	871	875	rVB2	1278317	2223260	38.83%	2.432%
41	12.111	875	877	880	rBV4	133645	245966	4.30%	0.269%
42	12.225	883	887	888	rBV	832781	1072283	18.73%	1.173%
43	12.317	892	895	897	rBV3	108340	211003	3.68%	0.231%
44	12.374	897	900	901	rBV2	236552	406823	7.10%	0.445%
45	12.408	901	903	906	rVB2	228757	517926	9.05%	0.566%
46	12.477	906	909	911	rBV2	721848	1264473	22.08%	1.383%
47	12.511	911	912	915	rVV2	332171	559145	9.76%	0.612%
48	12.568	915	917	919	rVV	1102919	1210626	21.14%	1.324%
49	12.648	919	924	926	rVV	3328814	4108158	71.74%	4.493%
50	12.682	926	927	930	rVV2	185297	399510	6.98%	0.437%
51	12.728	930	931	933	rVB	493245	674826	11.79%	0.738%
52	12.774	933	935	941	rBV3	428624	993423	17.35%	1.087%
53	12.877	941	944	946	rBV	3498253	5726099	100.00%	6.263%
54	12.922	946	948	950	rVB2	238543	351612	6.14%	0.385%
55	13.014	954	956	960	rVB	1379981	1716283	29.97%	1.877%
56	13.082	960	962	965	rBV	686758	1172842	20.48%	1.283%
57	13.139	966	967	968	rVV	264434	220272	3.85%	0.241%
58	13.174	968	970	973	rVB2	712695	1521491	26.57%	1.664%
59	13.242	973	976	979	rBV3	194937	531434	9.28%	0.581%
60	13.299	979	981	983	rVV	1083211	1058515	18.49%	1.158%
61	13.391	987	989	990	rVV	827056	752977	13.15%	0.824%
62	13.425	990	992	994	rVV	844323	816333	14.26%	0.893%
63	13.505	997	999	1001	rVB3	118360	156889	2.74%	0.172%
64	13.597	1003	1007	1010	rBV6	128952	356673	6.23%	0.390%
65	13.699	1014	1016	1018	rVB	971039	971235	16.96%	1.062%
66	13.814	1023	1026	1027	rBV	855990	1200455	20.96%	1.313%
67	13.837	1027	1028	1029	rVV	668100	702352	12.27%	0.768%
68	13.905	1032	1034	1037	rVV	874333	1148570	20.06%	1.256%
69	13.974	1038	1040	1044	rVB2	194151	392090	6.85%	0.429%
70	14.054	1044	1047	1049	rBV2	2073905	3526814	61.59%	3.857%
71	14.100	1049	1051	1053	rVV	2233377	2850997	49.79%	3.118%

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72	14.145	1054	1055	1058	rVV2	155110	299593	5.23%	0.328%
73	14.202	1058	1060	1065	rVV3	795132	1619388	28.28%	1.771%
74	14.317	1068	1070	1073	rVV	360605	603599	10.54%	0.660%
75	14.385	1073	1076	1077	rVV2	173400	352128	6.15%	0.385%
76	14.408	1077	1078	1079	rVV	277282	367941	6.43%	0.402%
77	14.454	1079	1082	1083	rVV	709699	1207397	21.09%	1.321%
78	14.488	1083	1085	1086	rVV2	507083	719855	12.57%	0.787%
79	14.534	1086	1089	1091	rVV3	451018	865803	15.12%	0.947%
80	14.568	1091	1092	1096	rVV2	393314	871589	15.22%	0.953%
81	14.637	1096	1098	1100	rVB	209546	292972	5.12%	0.320%
82	14.854	1115	1117	1120	rVB2	296258	410006	7.16%	0.448%
83	14.922	1120	1123	1127	rVV4	191928	477249	8.33%	0.522%
84	15.002	1127	1130	1135	rVB2	187966	410881	7.18%	0.449%
85	15.094	1135	1138	1143	rBV	1291075	3168278	55.33%	3.465%
86	15.197	1143	1147	1150	rVB2	462095	850017	14.84%	0.930%
87	15.391	1161	1164	1167	rVV	972696	1207416	21.09%	1.321%
88	15.448	1167	1169	1172	rVV	1002007	1272934	22.23%	1.392%
89	15.505	1172	1174	1176	rVV2	340260	533908	9.32%	0.584%
90	15.540	1176	1177	1182	rVB2	274875	415397	7.25%	0.454%
91	15.974	1212	1215	1217	rVV	545637	807456	14.10%	0.883%
92	16.614	1267	1271	1279	rVB2	236989	645980	11.28%	0.707%
93	16.785	1280	1286	1294	rVB4	182066	680269	11.88%	0.744%
94	16.946	1296	1300	1304	rBV	478629	983161	17.17%	1.075%
95	17.117	1312	1315	1317	rBV	115029	233994	4.09%	0.256%
96	17.163	1317	1319	1322	rVV2	108395	255908	4.47%	0.280%
97	17.220	1322	1324	1328	rVB	84594	184134	3.22%	0.201%
98	17.380	1333	1338	1343	rVB	409853	877084	15.32%	0.959%
99	17.517	1347	1350	1354	rVB5	75334	180820	3.16%	0.198%
100	18.534	1435	1439	1446	rVB5	98626	335515	5.86%	0.367%

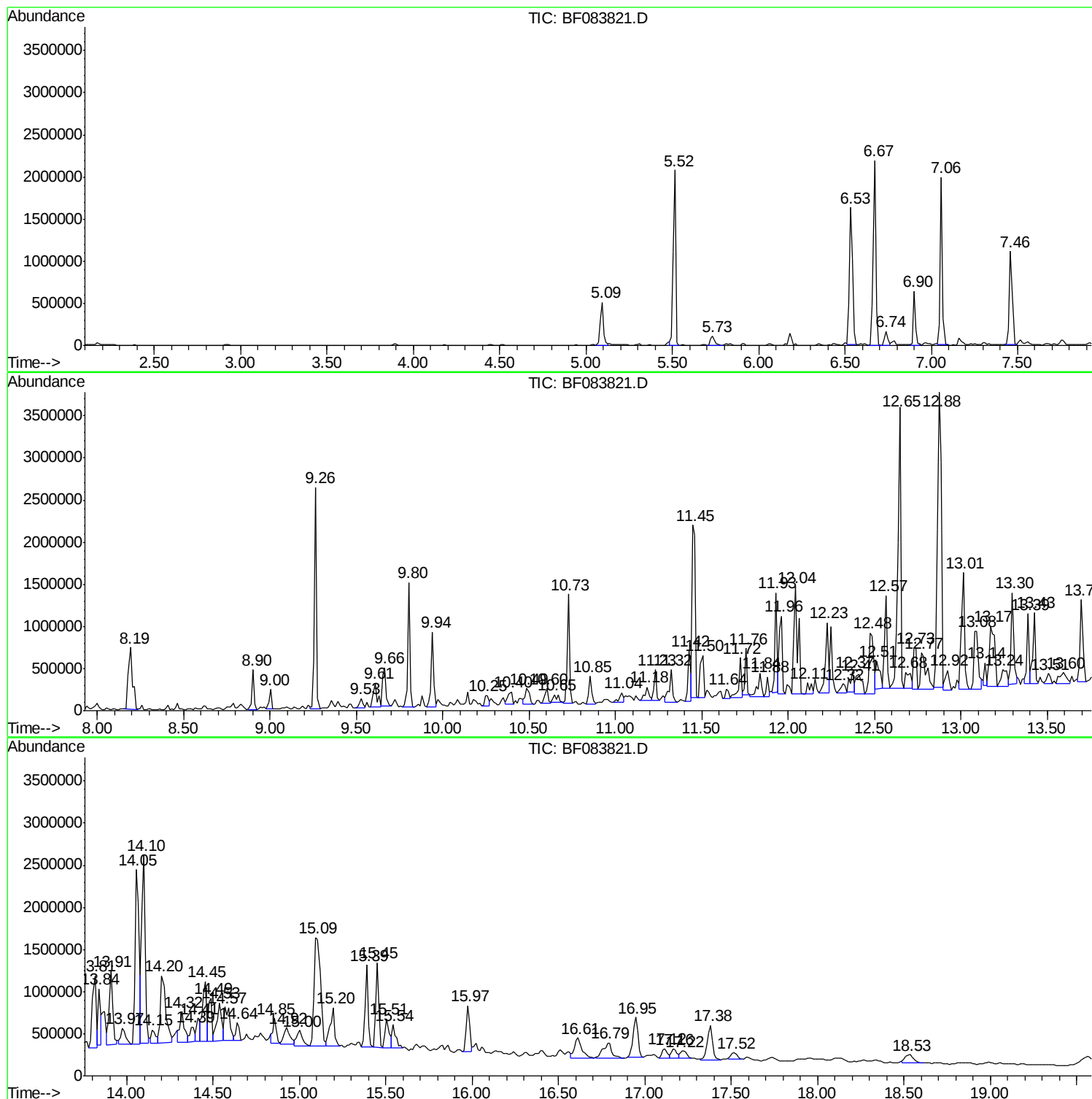
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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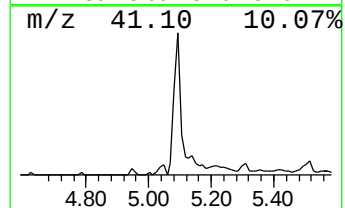
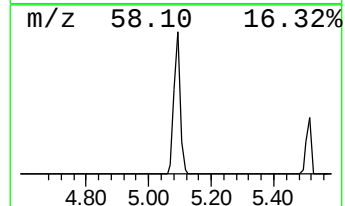
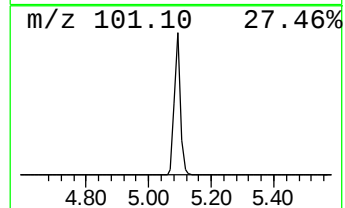
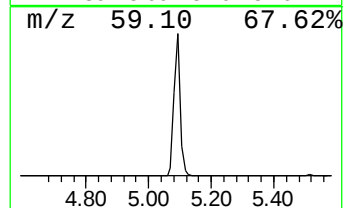
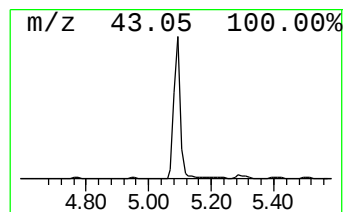
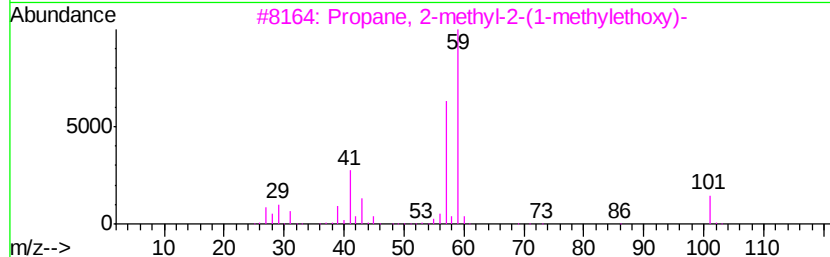
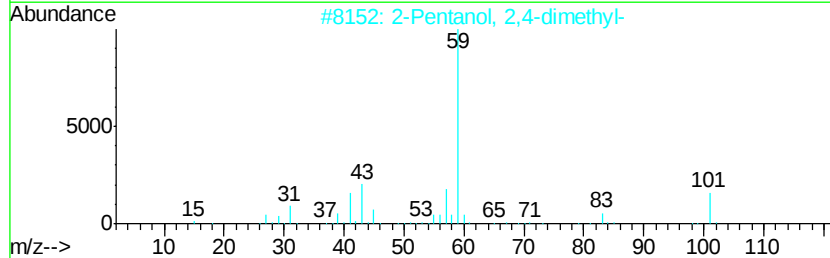
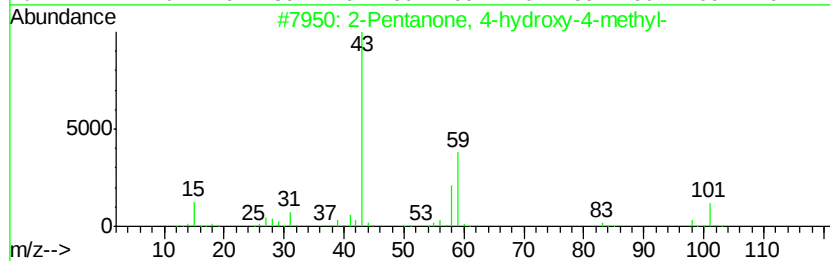
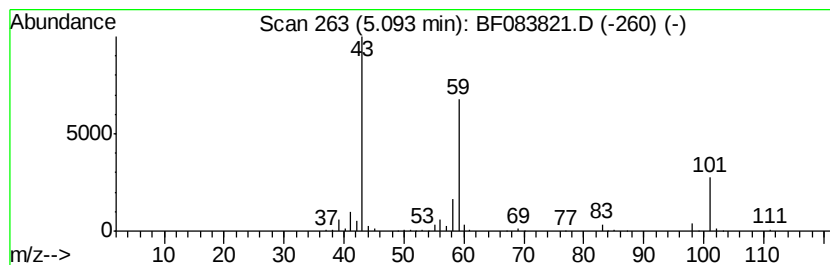
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	22.69 ng	686263	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	40
3			Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	33
4			4-Hydroxy-3-hexanone	116	C6H12O2	004984-85-4	23
5			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	17



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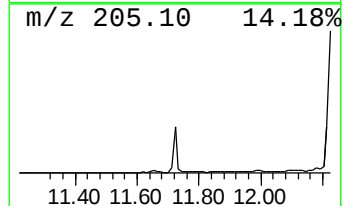
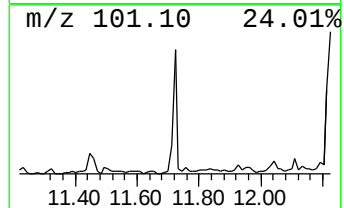
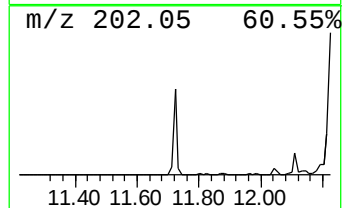
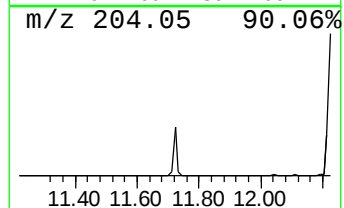
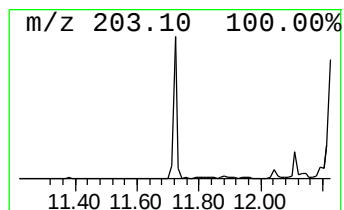
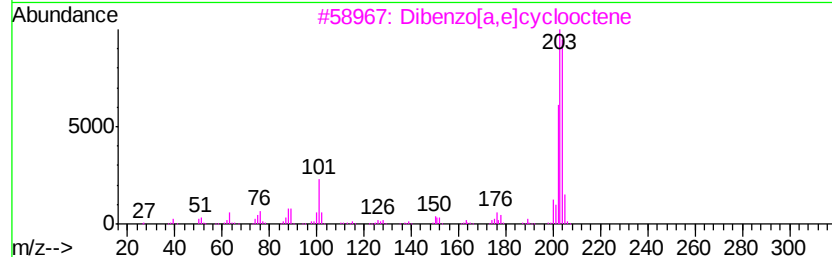
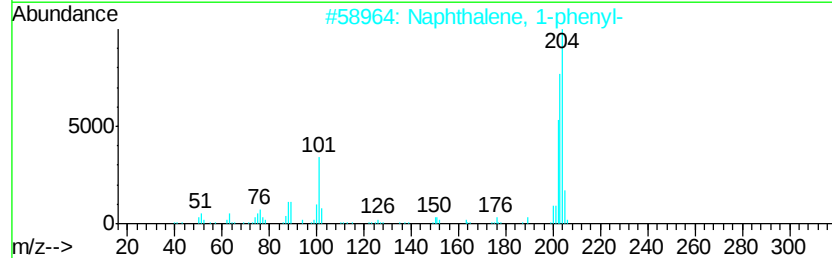
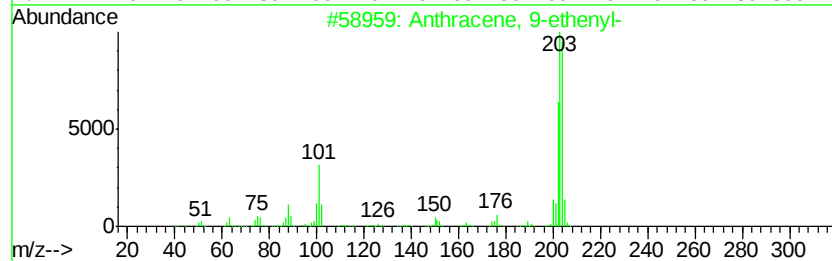
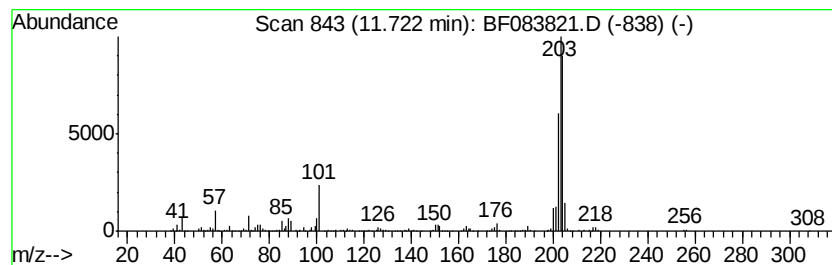
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Anthracene, 9-ethenyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.72	14.89 ng	527398	Phenanthrene-d10	11.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-ethenyl-	204	C16H12	002444-68-0	95
2		Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	91
3		Dibenzo[a,e]cyclooctene	204	C16H12	000262-89-5	90
4		1H-Indene, 1-(phenylmethylene)-	204	C16H12	005394-86-5	81
5		Cyclobuta[1'',2'':3,4;3'',4'':3'...	204	C16H12	006574-36-3	76



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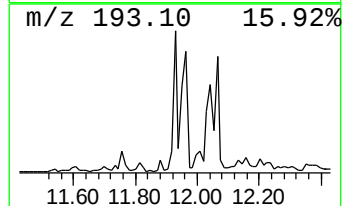
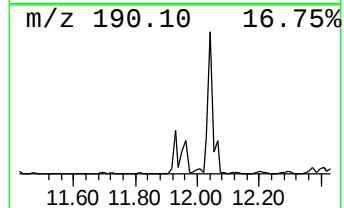
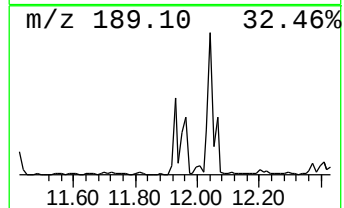
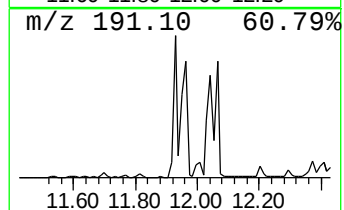
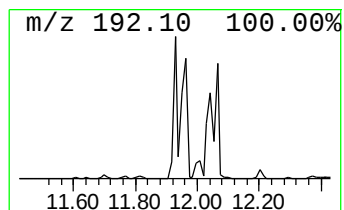
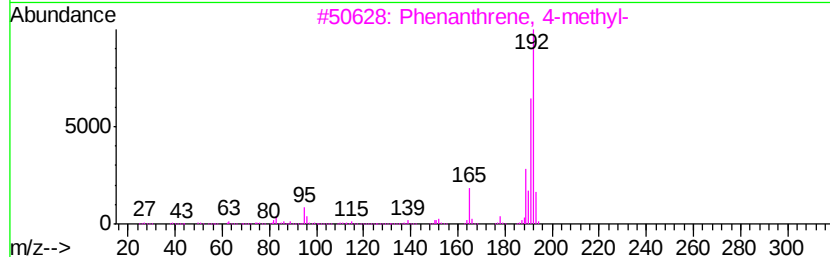
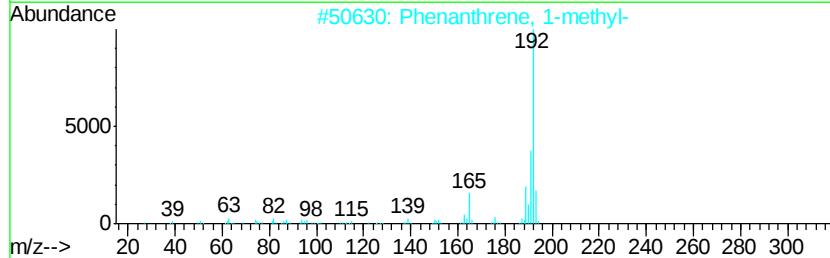
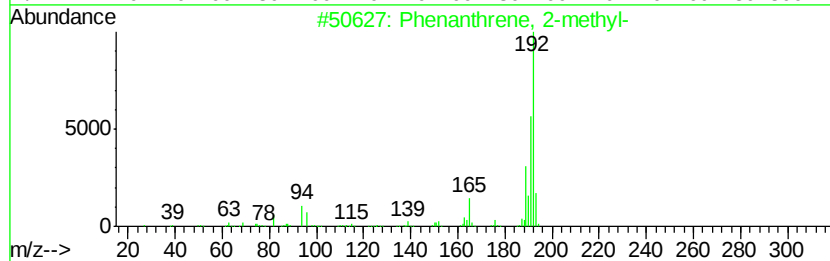
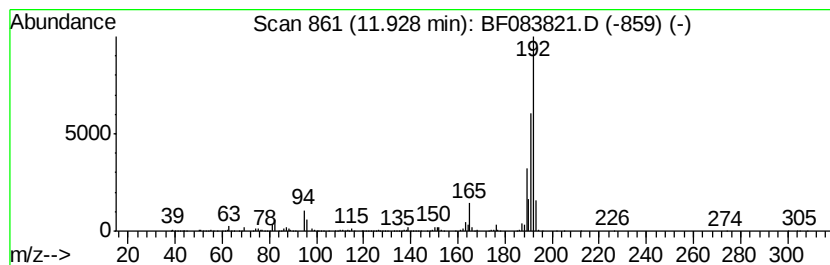
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	30.63 ng	1085030	Phenanthrene-d10	11.43

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	95
3		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	94
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	91



Data Path : V:\HPCHEM1\BNA_F\DATA\BF121515\
Data File : BF083821.D
Acq On : 15 Dec 2015 16:31
Operator : UM/IZ
Sample : G4739-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

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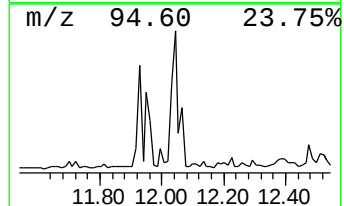
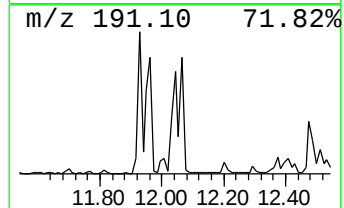
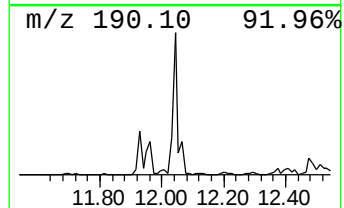
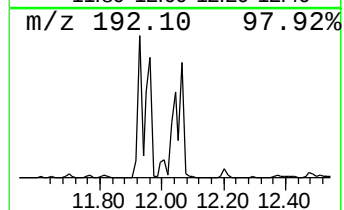
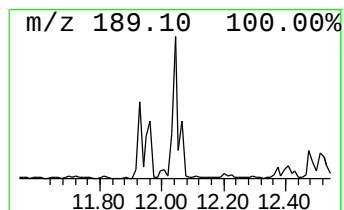
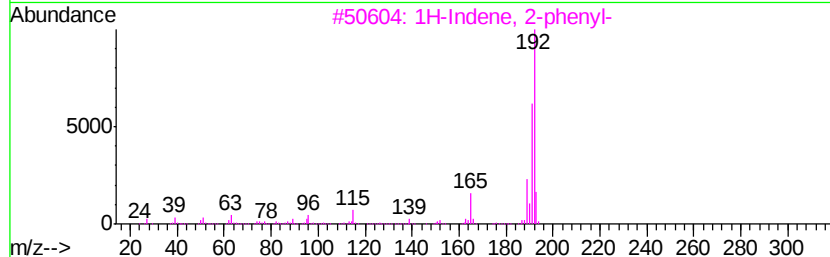
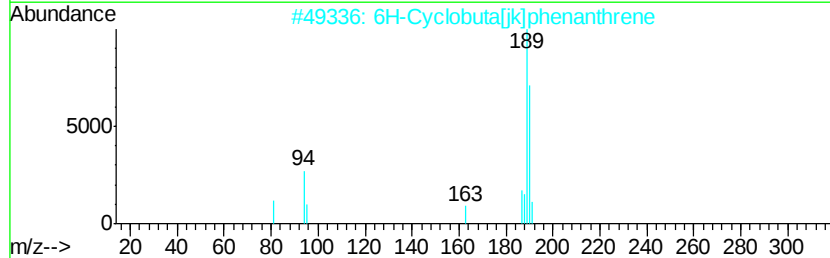
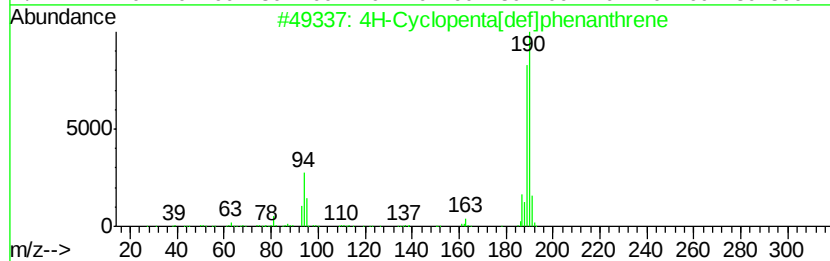
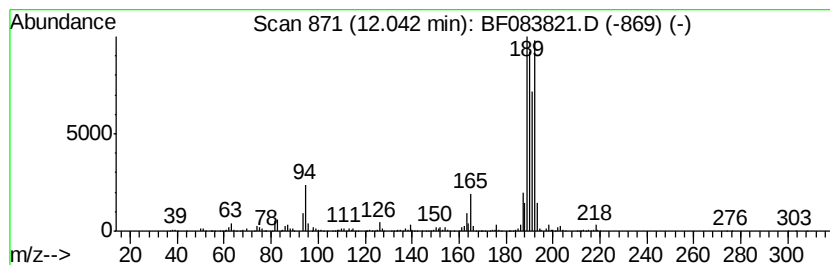
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.04	62.76 ng	2223260	Phenanthrene-d10	11.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
2		6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	49
3		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	46
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	46
5		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	46



Data Path : V:\HPCHEM1\BNA F\DATA\BF121515\
Data File : BF083821.D
Acq On : 15 Dec 2015 16:31
Operator : UM/IZ
Sample : G4739-07
Misc :
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ClientSampleId :
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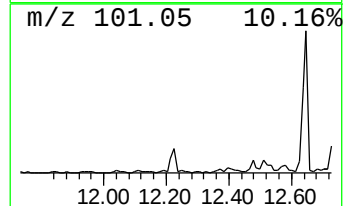
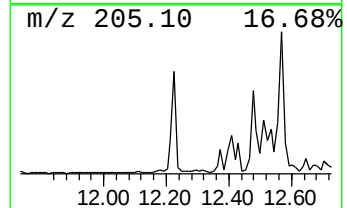
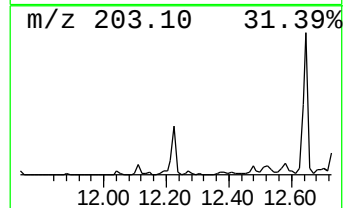
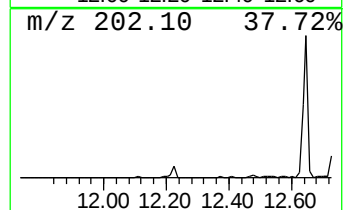
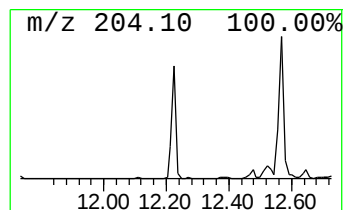
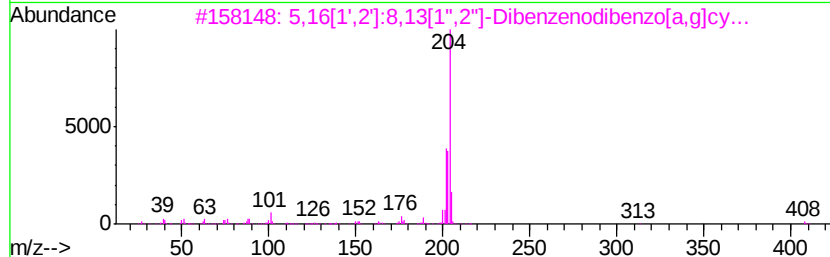
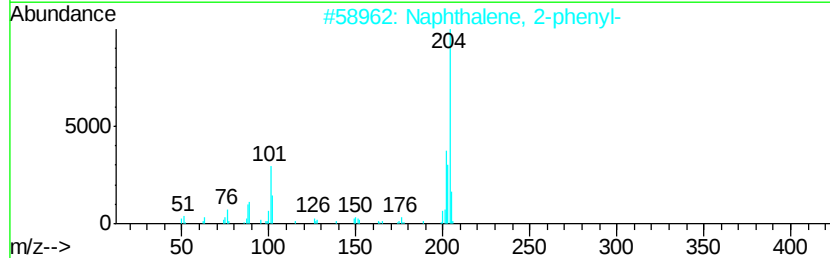
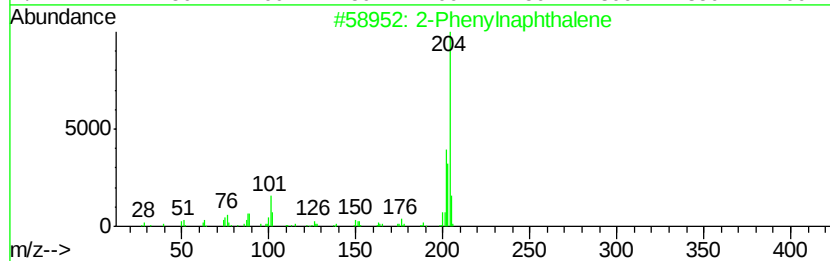
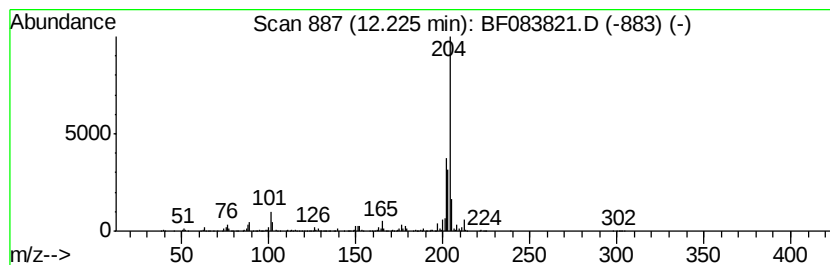
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 2-Phenylnaphthalene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.23	30.27 ng	1072280	Phenanthrene-d10	11.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Phenylnaphthalene	204	C16H12	035465-71-5	95
2		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	93
3		5,16[1',2']8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	87
4		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	74
5		Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	53



Data Path : V:\HPCHEM1\BNA F\DATA\BF121515\
Data File : BF083821.D
Acq On : 15 Dec 2015 16:31
Operator : UM/IZ
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Misc :
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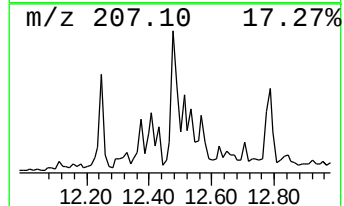
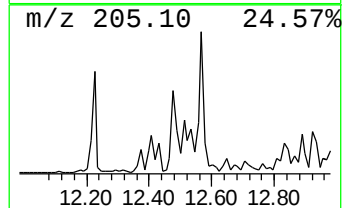
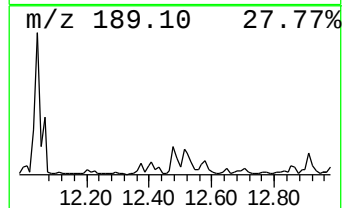
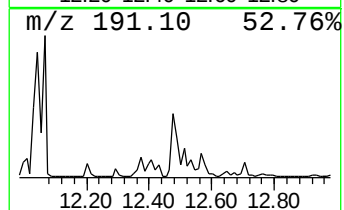
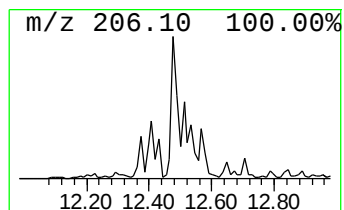
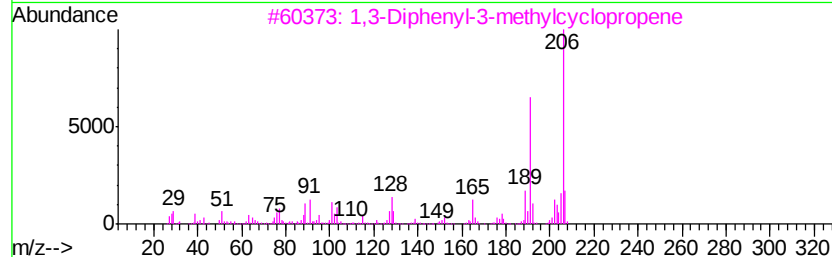
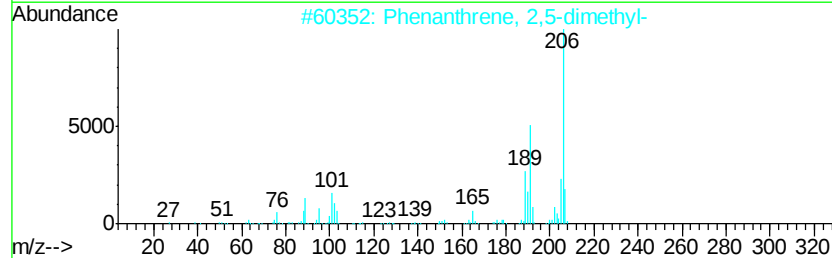
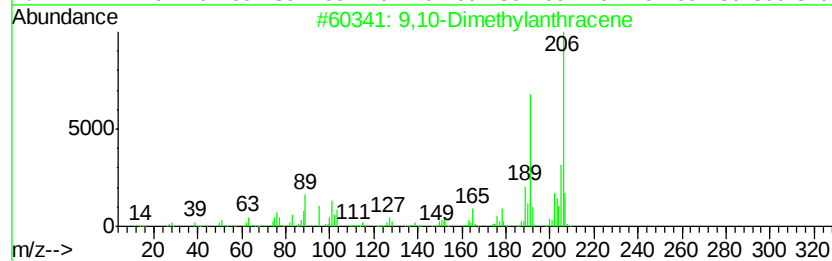
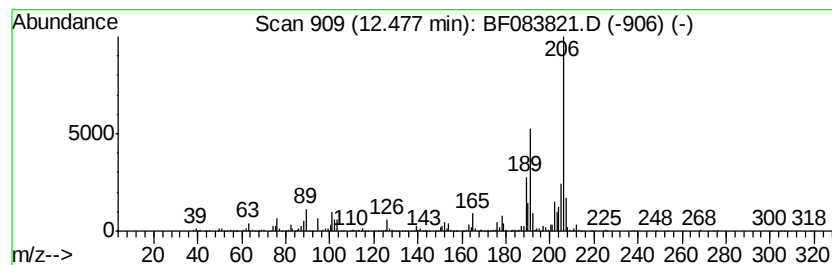
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9,10-Dimethylantracene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.48	35.70 ng	1264470	Phenanthrene-d10	11.43

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylantracene	206	C16H14	000781-43-1	97	
2	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	96	
3	1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	94	
4	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	91	
5	Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	90	



Data Path : V:\HPCHEM1\BNA F\DATA\BF121515\
Data File : BF083821.D
Acq On : 15 Dec 2015 16:31
Operator : UM/IZ
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Misc :
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ClientSampleId :
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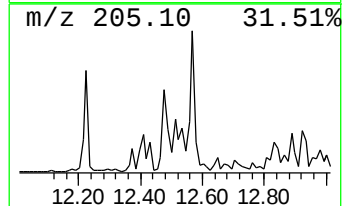
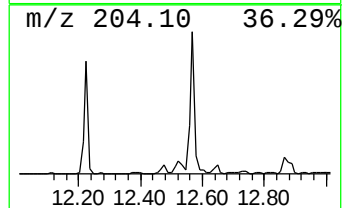
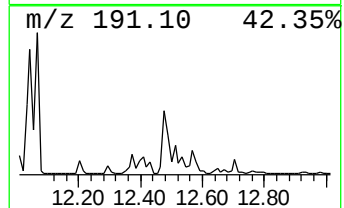
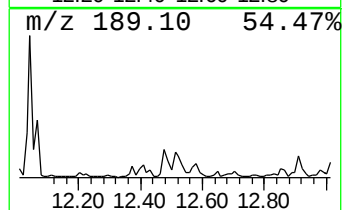
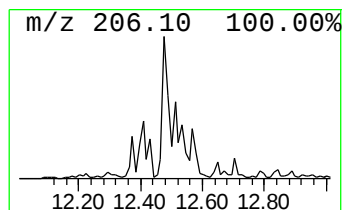
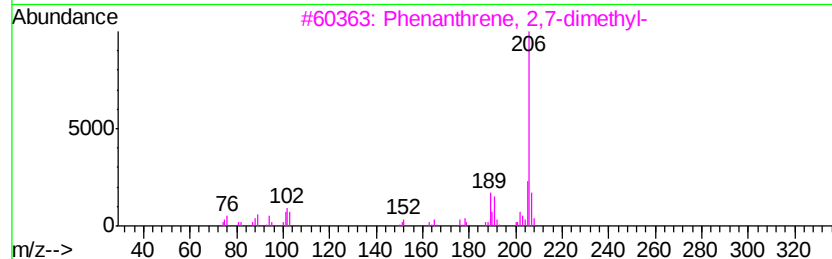
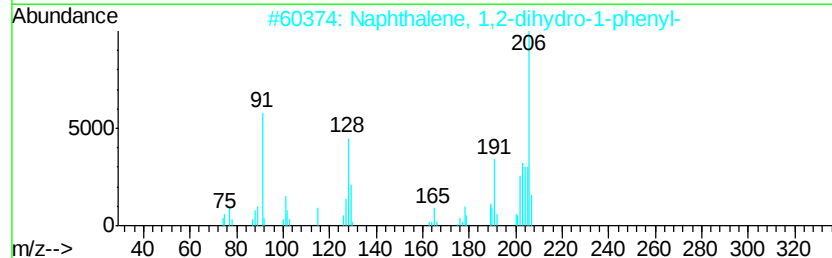
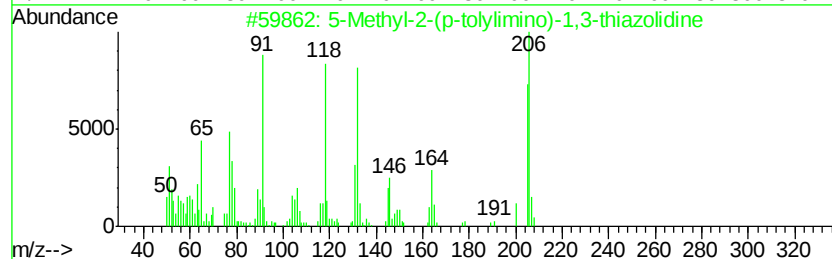
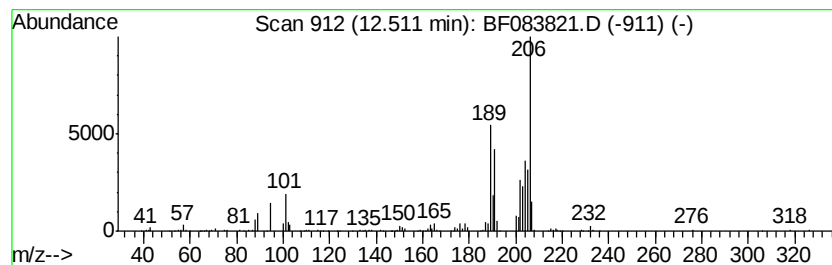
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 5-Methyl-2-(p-tolylimino)-1... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.51	15.78 ng	559145	Phenanthrene-d10	11.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Methyl-2-(p-tolylimino)-1,3-th...	206	C11H14N2S	070446-87-6	80
2		Naphthalene, 1,2-dihydro-1-phenyl-	206	C16H14	016606-46-5	64
3		Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	64
4		Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	59
5		Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	58



Data Path : V:\HPCHEM1\BNA_F\DATA\BF121515\
Data File : BF083821.D
Acq On : 15 Dec 2015 16:31
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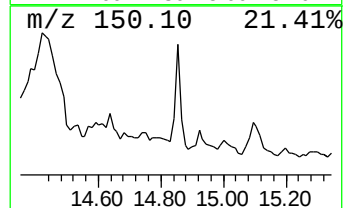
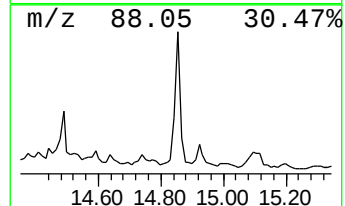
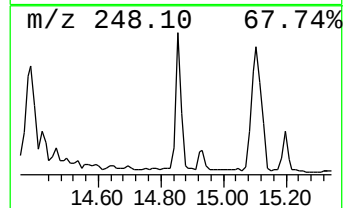
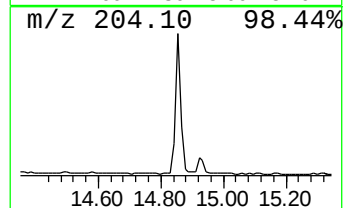
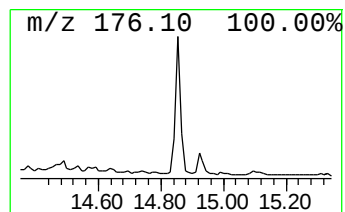
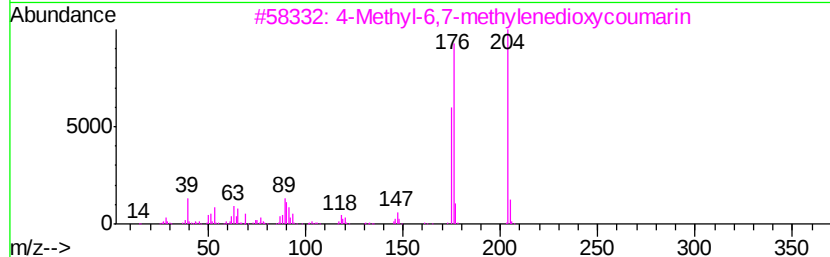
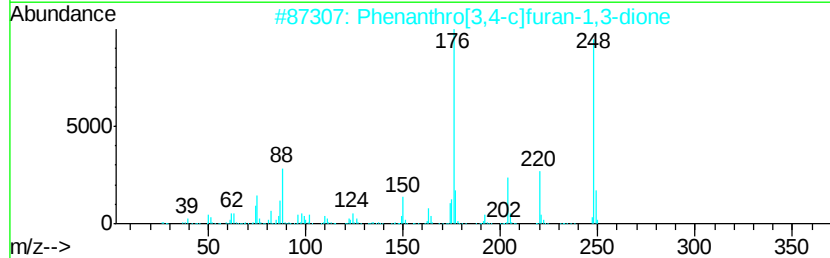
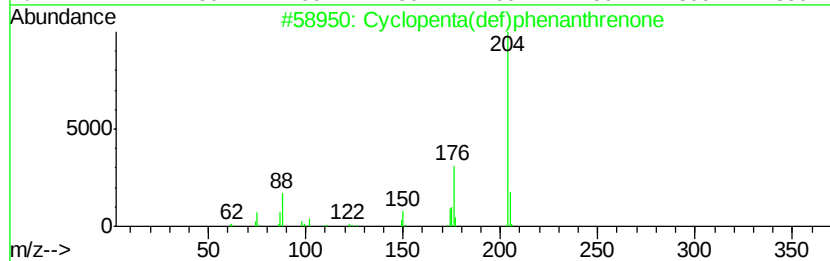
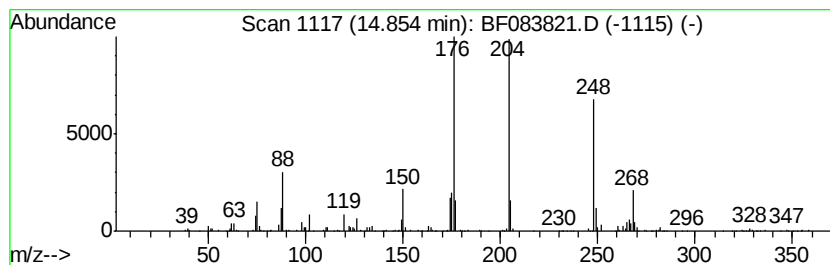
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Cyclopenta(def)phenanthrenone Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	15.36 ng	410006	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	50
2		Phenanthro[3,4-c]furan-1,3-dione	248	C16H8O3	005723-54-6	46
3		4-Methyl-6,7-methylenedioxy coumarin	204	C11H8O4	015071-04-2	43
4		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	38
5		4-Methoxy-6-(3-fluorophenyl)pyri...	204	C11H9FN2O	076128-74-0	22



Data Path : V:\HPCHEM1\BNA F\DATA\BF121515\
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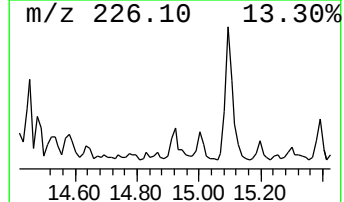
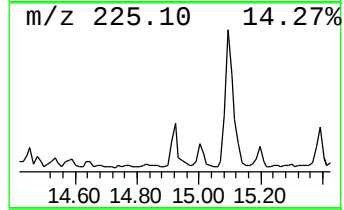
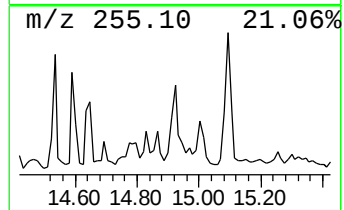
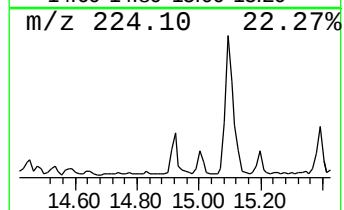
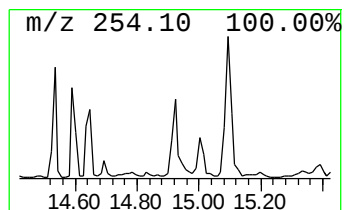
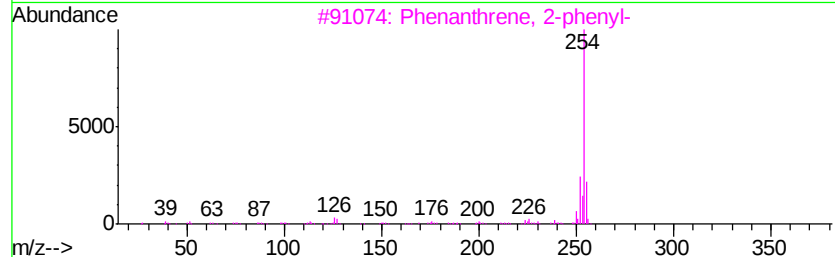
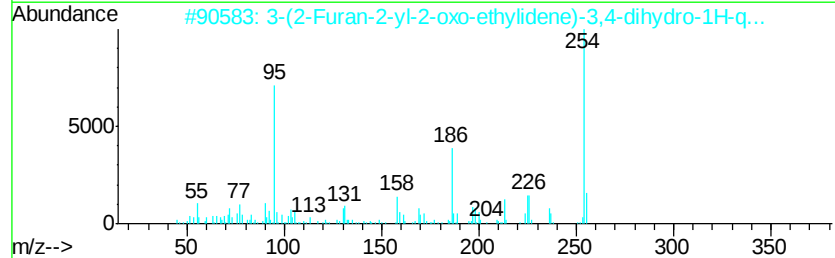
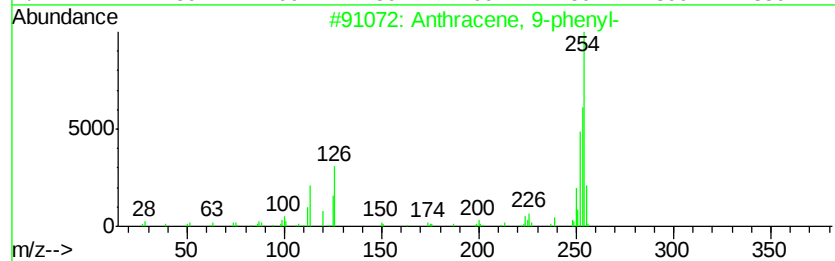
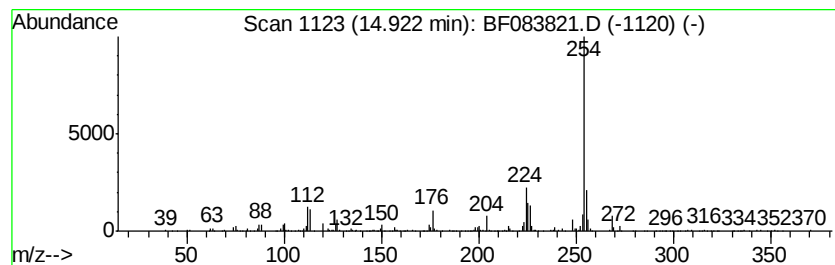
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Anthracene, 9-phenyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.92	17.88 ng	477249	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-phenyl-	254	C20H14	000602-55-1	59
2		3-(2-Furan-2-yl-2-oxo-ethylidene...	254	C14H10N2O3	1000297-32-1	50
3		Phenanthrene, 2-phenyl-	254	C20H14	004325-77-3	47
4		9,10-Anthracenedione, 1,8-dihydr...	254	C15H10O4	000481-74-3	43
5		9,10-Ethenoanthracene-11,12-dica...	254	C18H10N2	019067-49-3	42



Data Path : V:\HPCHEM1\BNA_F\DATA\BF121515\
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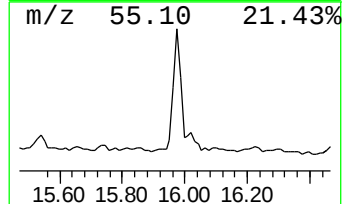
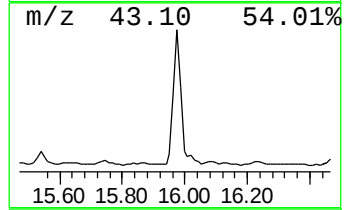
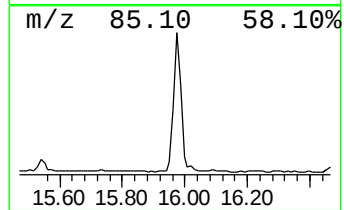
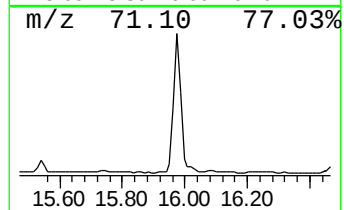
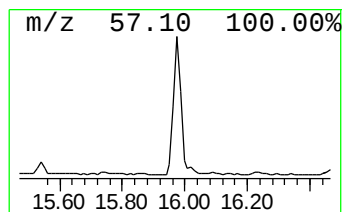
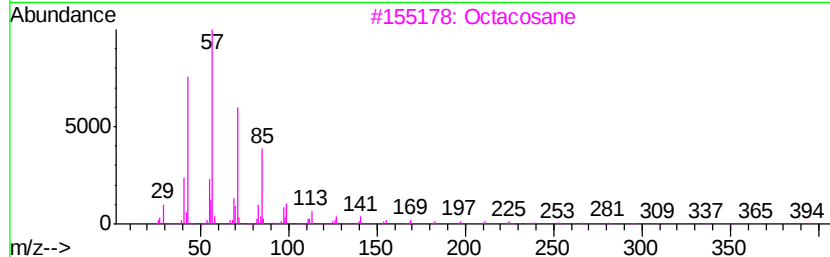
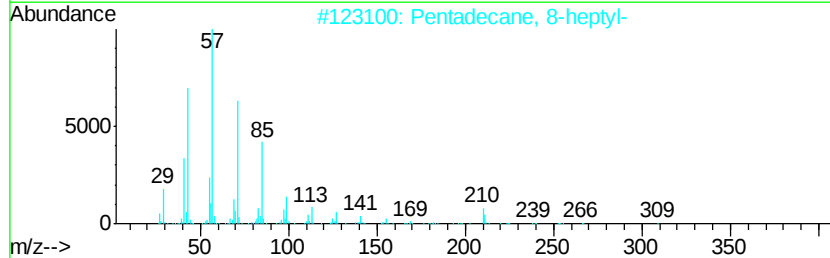
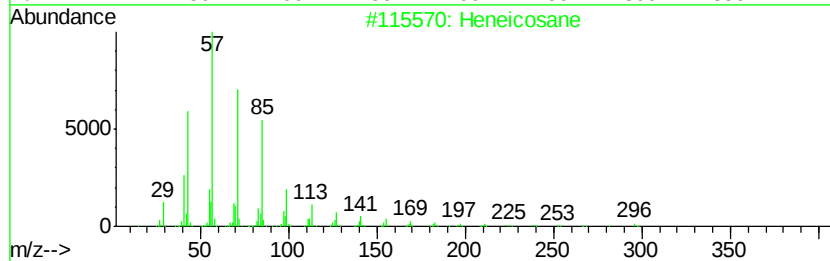
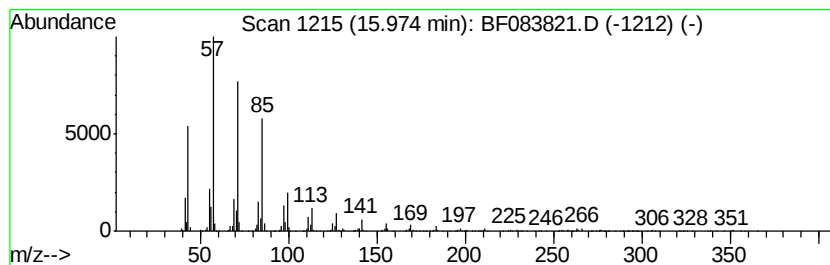
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Heneicosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	30.25 ng	807456	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	91
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
3		Octacosane	394	C28H58	000630-02-4	90
4		Octadecane	254	C18H38	000593-45-3	87
5		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	87



Data Path : V:\HPCHEM1\BNA F\DATA\BF121515\
Data File : BF083821.D
Acq On : 15 Dec 2015 16:31
Operator : UM/IZ
Sample : G4739-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
K1206-0-2

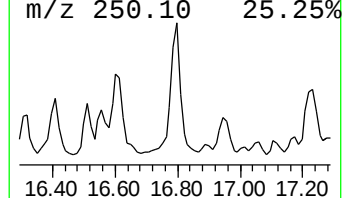
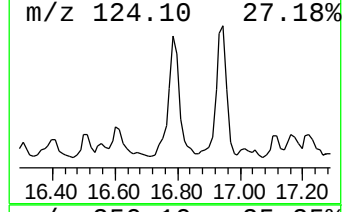
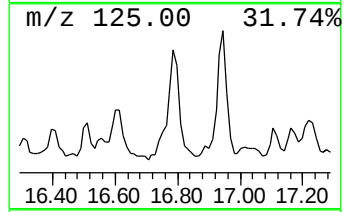
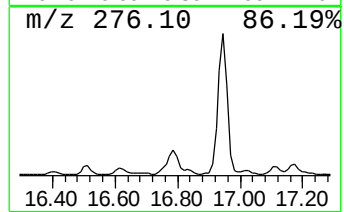
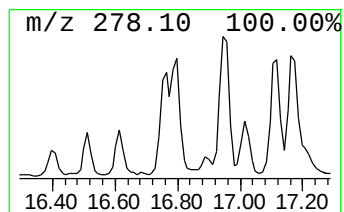
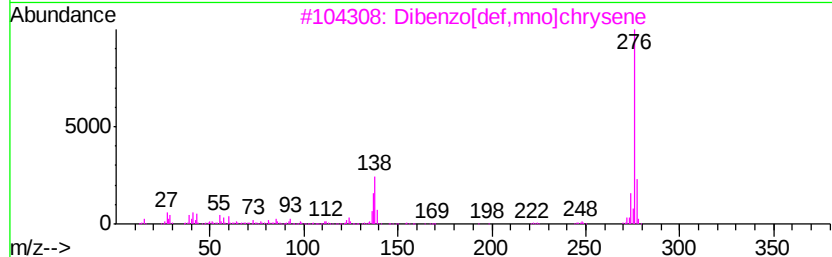
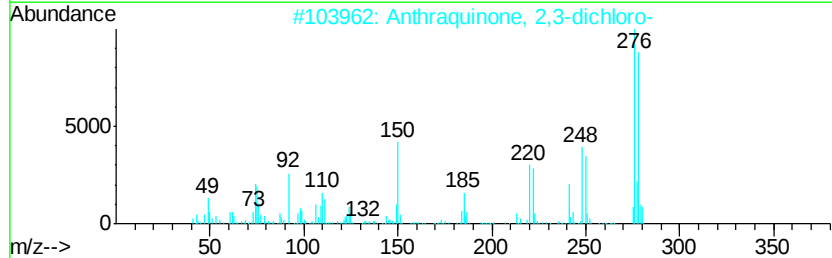
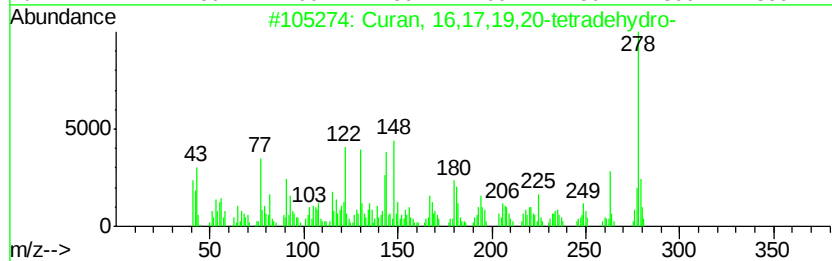
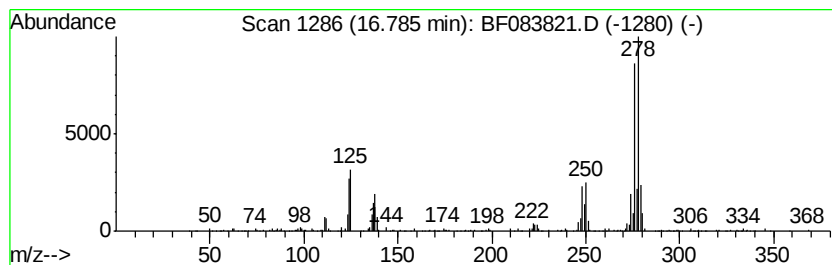
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Curan, 16,17,19,20-tetradeh... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.79	25.48 ng	680269	Perylene-d12	15.51

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Curan, 16,17,19,20-tetradehydro-	278	C19H22N2	056053-17-9	89
2		Anthraquinone, 2,3-dichloro-	276	C14H6Cl2O2	000084-45-7	83
3		Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	80
4		Anthraquinone, 2,6-dichloro-	276	C14H6Cl2O2	000605-40-3	53
5		Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	42



Data Path : V:\HPCHEM1\BNA_F\DATA\BF121515\
Data File : BF083821.D
Acq On : 15 Dec 2015 16:31
Operator : UM/IZ
Sample : G4739-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
K1206-0-2

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.09	22.7	ng	686263	1	6.90	604806	20.0
Anthracene, 9-eth...	11.72	14.9	ng	527398	4	11.43	708488	20.0
Phenanthrene, 2-m...	11.93	30.6	ng	1085030	4	11.43	708488	20.0
4H-Cyclopenta[def...	12.04	62.8	ng	2223260	4	11.43	708488	20.0
2-Phenylnaphthalene	12.23	30.3	ng	1072280	4	11.43	708488	20.0
9,10-Dimethylanth...	12.48	35.7	ng	1264470	4	11.43	708488	20.0
5-Methyl-2-(p-tol...	12.51	15.8	ng	559145	4	11.43	708488	20.0
Cyclopenta(def)ph...	14.85	15.4	ng	410006	6	15.51	533908	20.0
Anthracene, 9-phe...	14.92	17.9	ng	477249	6	15.51	533908	20.0
Heneicosane	15.97	30.3	ng	807456	6	15.51	533908	20.0
Curan, 16,17,19,2...	16.79	25.5	ng	680269	6	15.51	533908	20.0