

Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
 Data File : BF079426.D
 Acq On : 22 May 2015 7:51
 Operator : TP/IZ
 Sample : G2289-08
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-29D

Integration Parameters: rteint.p

Integrator: RTE

Soothing : 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-BF043015.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF079426.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.336	11	13	25	rVB	146550	401760	4.55%	0.780%
2	1.496	25	27	34	rVV	245074	333735	3.78%	0.648%
3	1.610	36	37	46	rVV	740799	1281201	14.50%	2.488%
4	1.724	46	47	86	rVB	4136338	8838539	100.00%	17.162%
5	4.810	315	317	323	rBV	96063	131827	1.49%	0.256%
6	5.347	362	364	371	rBV	1158873	1282639	14.51%	2.491%
7	6.650	475	478	480	rBV	976600	1337749	15.14%	2.598%
8	6.776	487	489	492	rBV	1218314	1368024	15.48%	2.656%
9	7.073	513	515	517	rBV	293193	350872	3.97%	0.681%
10	7.119	517	519	520	rBV	94776	112547	1.27%	0.219%
11	7.256	528	531	533	rBV	1018376	997696	11.29%	1.937%
12	7.759	574	575	578	rBV	515880	717626	8.12%	1.393%
13	8.651	650	653	654	rBV	373359	570422	6.45%	1.108%
14	8.673	654	655	657	rVB	415782	310271	3.51%	0.602%
15	9.531	728	730	734	rBV	451382	424659	4.80%	0.825%
16	9.645	738	740	743	rBV	263686	378332	4.28%	0.735%
17	9.988	768	770	773	rBV	1106351	1441461	16.31%	2.799%
18	10.228	787	791	796	rVB	239156	279000	3.16%	0.542%
19	10.308	796	798	801	rBV	124464	214177	2.42%	0.416%
20	10.399	803	806	808	rBV	194440	224971	2.55%	0.437%
21	10.434	808	809	812	rVB	218996	172190	1.95%	0.334%
22	10.502	812	815	817	rBV	115293	121931	1.38%	0.237%
23	10.548	817	819	824	rVB	79756	142804	1.62%	0.277%
24	10.628	824	826	830	rBV2	180506	285359	3.23%	0.554%
25	10.811	840	842	844	rVB	503302	548588	6.21%	1.065%
26	10.856	844	846	848	rBV	405443	385308	4.36%	0.748%
27	10.959	853	855	858	rBV	104951	189245	2.14%	0.367%
28	11.074	862	865	867	rBV	237365	307496	3.48%	0.597%
29	11.325	885	887	890	rBV3	125084	212008	2.40%	0.412%
30	11.428	895	896	900	rVB2	162498	215312	2.44%	0.418%
31	11.497	900	902	904	rVB	388763	389258	4.40%	0.756%
32	11.565	904	908	910	rBV2	101376	224663	2.54%	0.436%
33	11.691	918	919	923	rVB2	66406	118662	1.34%	0.230%
34	11.794	925	928	931	rBV	1065806	1109690	12.56%	2.155%
35	11.919	935	939	941	rBV2	67415	132986	1.50%	0.258%
36	12.182	958	962	963	rVV2	81707	166506	1.88%	0.323%
37	12.274	966	970	972	rVV4	58179	136901	1.55%	0.266%

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Integrator: RTE

Soothing : ON

Sampling : 1

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Max Peaks: 100

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M

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38	12.434	982	984	986	rVB2	111130	119266	1.35%	0.232%
39	12.525	990	992	993	rVB	114665	119277	1.35%	0.232%
40	12.685	1001	1006	1008	rBV	2384396	2935858	33.22%	5.701%
41	12.742	1008	1011	1014	rVB	592933	670256	7.58%	1.301%
42	12.960	1028	1030	1031	rBV	98802	104719	1.18%	0.203%
43	13.280	1055	1058	1060	rBV	384706	459568	5.20%	0.892%
44	13.314	1060	1061	1064	rVB	624163	559193	6.33%	1.086%
45	13.371	1064	1066	1068	rBV	175941	210025	2.38%	0.408%
46	13.417	1068	1070	1071	rVV	444717	608204	6.88%	1.181%
47	13.440	1071	1072	1075	rVB	514175	634795	7.18%	1.233%
48	13.657	1089	1091	1095	rVB2	189776	252290	2.85%	0.490%
49	13.840	1105	1107	1109	rBV2	445122	711192	8.05%	1.381%
50	13.965	1114	1118	1120	rBV2	165878	290974	3.29%	0.565%
51	14.011	1120	1122	1123	rVV	131690	219788	2.49%	0.427%
52	14.034	1123	1124	1126	rVB	274636	257338	2.91%	0.500%
53	14.080	1126	1128	1130	rVB	86368	127510	1.44%	0.248%
54	14.160	1130	1135	1137	rBV	1615251	1913178	21.65%	3.715%
55	14.228	1139	1141	1143	rVB2	409613	474723	5.37%	0.922%
56	14.331	1148	1150	1152	rBV3	56551	96027	1.09%	0.186%
57	14.434	1157	1159	1161	rVB	1699112	1869974	21.16%	3.631%
58	14.525	1166	1167	1170	rVB	264890	270123	3.06%	0.525%
59	14.594	1171	1173	1175	rBV	126315	151470	1.71%	0.294%
60	14.651	1175	1178	1180	rBV	1495611	1731741	19.59%	3.363%
61	14.720	1181	1184	1186	rBV2	101130	159342	1.80%	0.309%
62	14.811	1189	1192	1194	rBV	691132	986034	11.16%	1.915%
63	14.857	1194	1196	1198	rVB2	252868	294936	3.34%	0.573%
64	14.937	1198	1203	1205	rBV	271392	306411	3.47%	0.595%
65	14.983	1205	1207	1209	rBV	188777	259006	2.93%	0.503%
66	15.085	1215	1216	1218	rVB	85823	106740	1.21%	0.207%
67	15.131	1218	1220	1223	rBV	108262	133942	1.52%	0.260%
68	15.394	1238	1243	1245	rBV2	41011	95398	1.08%	0.185%
69	15.623	1261	1263	1265	rBV	86702	122171	1.38%	0.237%
70	15.657	1265	1266	1270	rVV2	86085	187716	2.12%	0.365%
71	15.840	1277	1282	1286	rBV2	190908	363323	4.11%	0.705%
72	15.943	1286	1291	1293	rVV2	843558	1746887	19.76%	3.392%
73	15.977	1293	1294	1296	rVV	659578	618834	7.00%	1.202%
74	16.068	1299	1302	1305	rVV2	97903	153751	1.74%	0.299%
75	16.251	1317	1318	1323	rVB3	56717	108023	1.22%	0.210%
76	16.457	1333	1336	1338	rBV	151696	196912	2.23%	0.382%
77	16.583	1344	1347	1348	rVV2	76945	125527	1.42%	0.244%
78	16.617	1348	1350	1351	rVV2	71464	107500	1.22%	0.209%

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

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79	16.674	1353	1355	1358	rVV2	47924	110696	1.25%	0.215%
80	17.269	1405	1407	1413	rBV	456662	1157446	13.10%	2.247%
81	17.394	1416	1418	1421	rVB	126227	158032	1.79%	0.307%
82	17.600	1434	1436	1440	rVB	288774	413598	4.68%	0.803%
83	17.669	1440	1442	1446	rBV	584762	685807	7.76%	1.332%
84	17.737	1446	1448	1450	rBV	498289	679052	7.68%	1.319%
85	18.949	1551	1554	1557	rVB3	50662	112636	1.27%	0.219%
86	19.109	1565	1568	1572	rBV2	216025	460898	5.21%	0.895%
87	19.177	1572	1574	1580	rVB2	47660	105536	1.19%	0.205%
88	19.269	1580	1582	1585	rBV	55648	126584	1.43%	0.246%
89	19.326	1585	1587	1596	rVB3	75132	245470	2.78%	0.477%
90	19.497	1599	1602	1608	rVB	219069	393145	4.45%	0.763%
91	19.714	1618	1621	1626	rVB	61803	134047	1.52%	0.260%

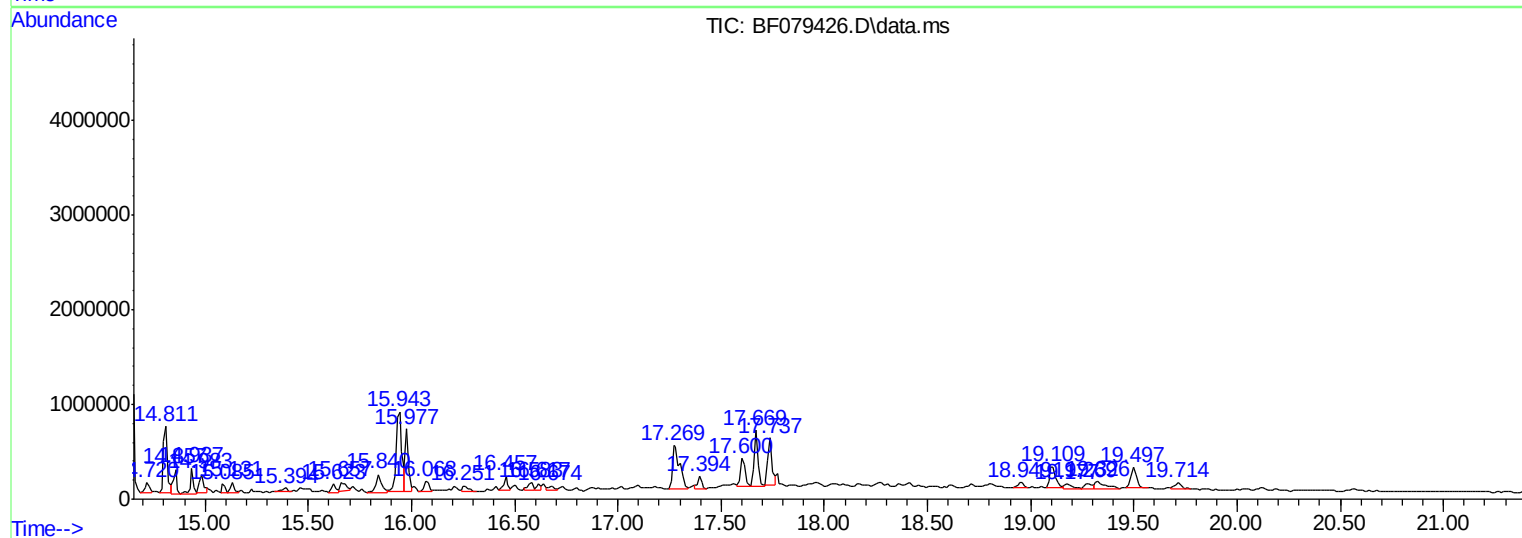
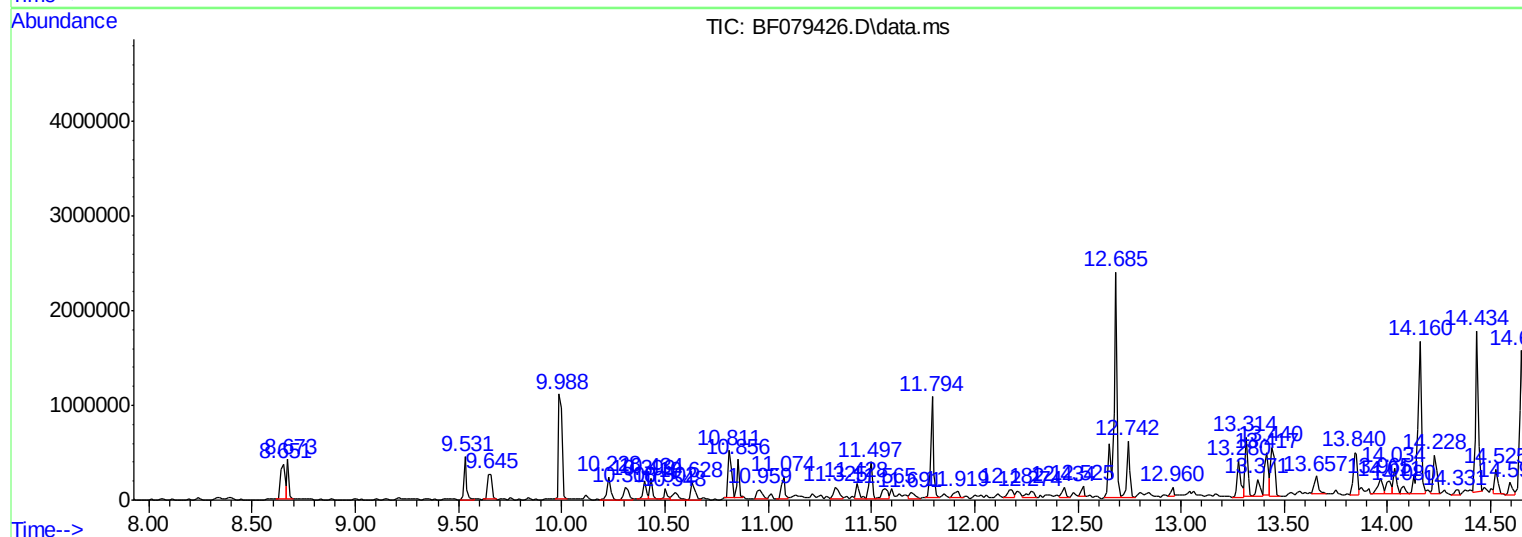
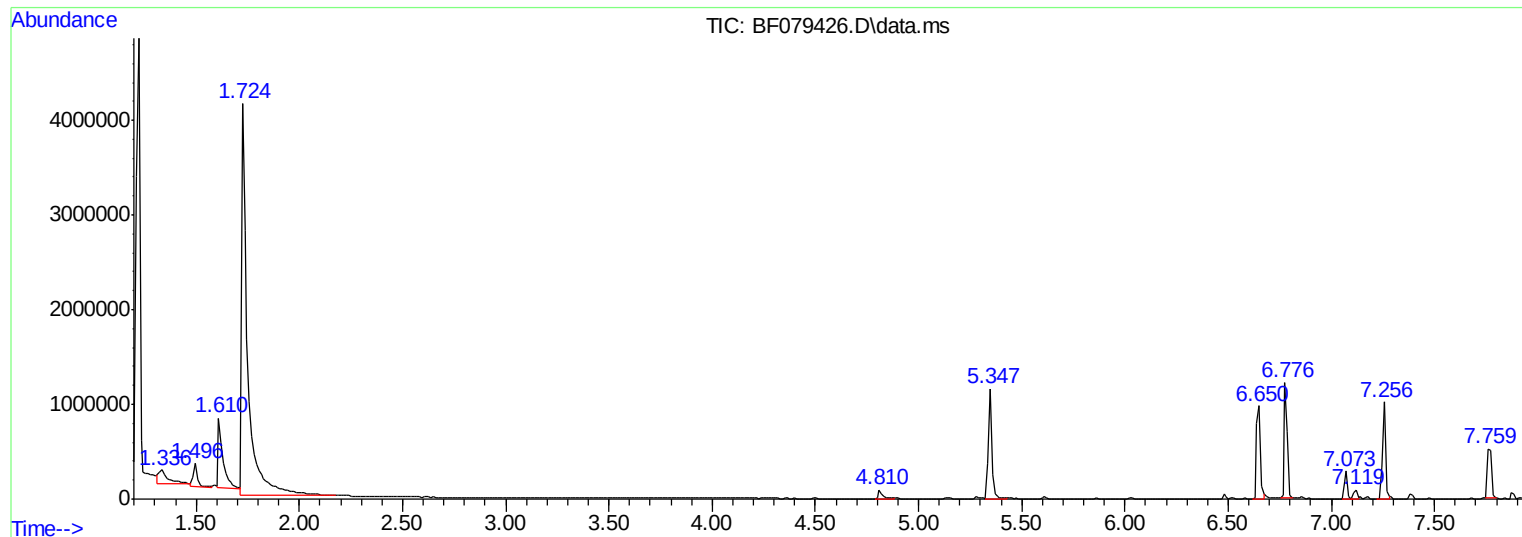
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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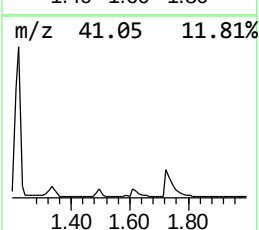
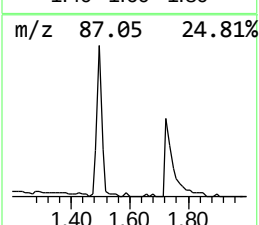
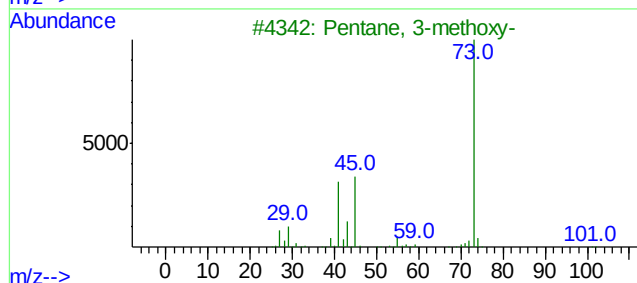
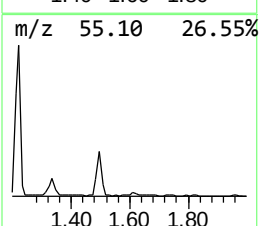
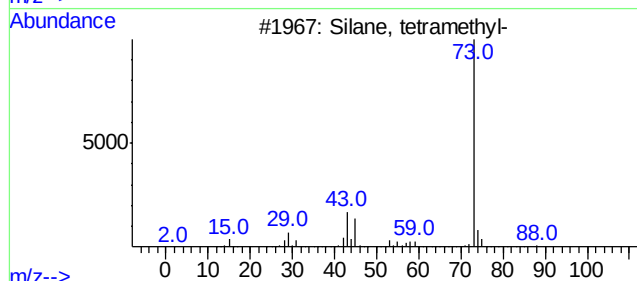
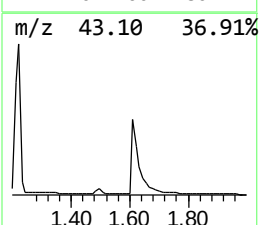
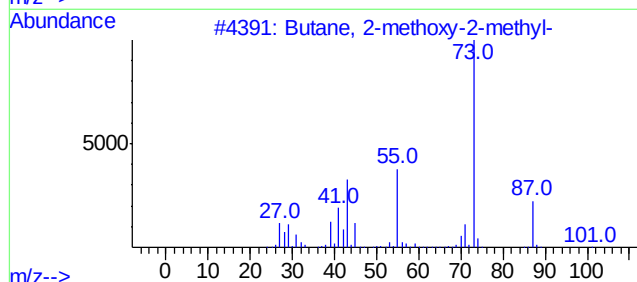
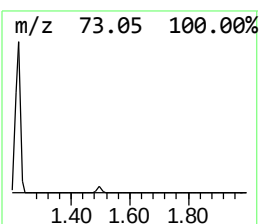
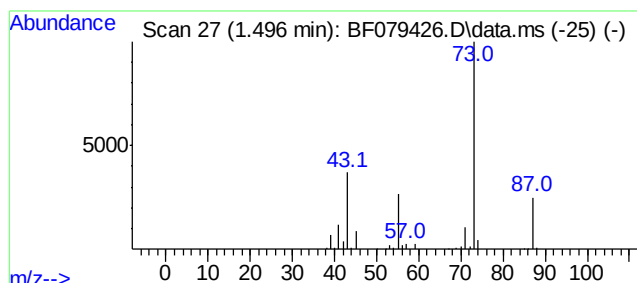
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.496	19.02 ng	333735	1,4-Dichlorobenzene-d4	7.073

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	Silane, tetramethyl-	88	C4H12Si	000075-76-3	25
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4	N-Ethylformamide	73	C3H7NO	000627-45-2	9
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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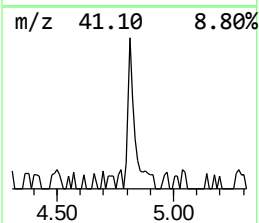
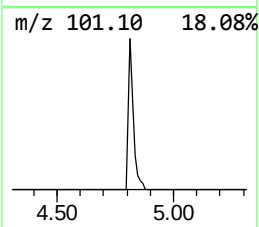
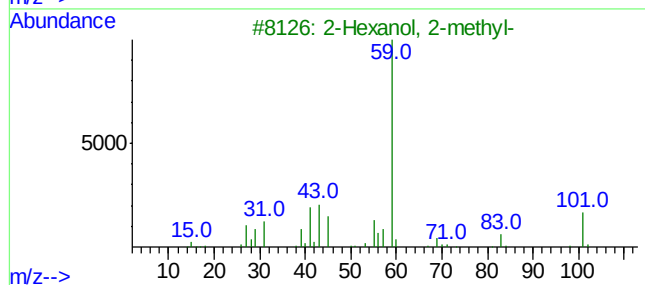
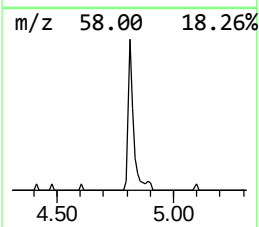
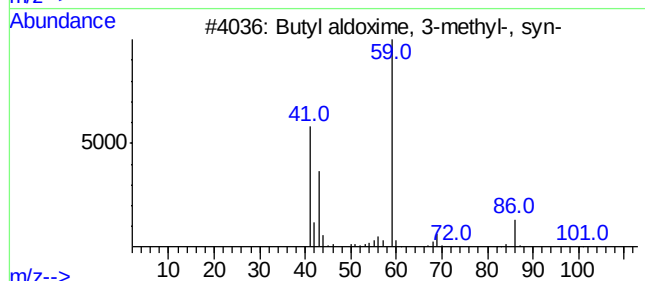
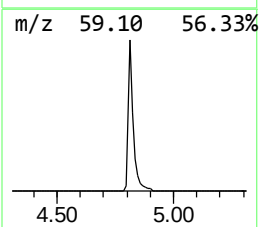
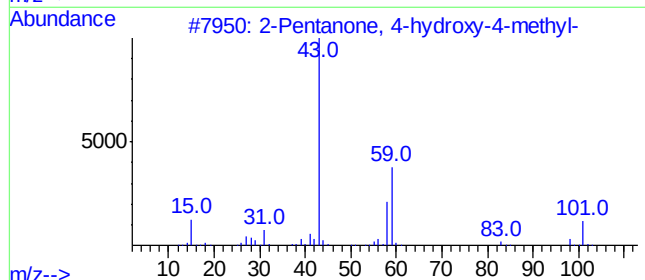
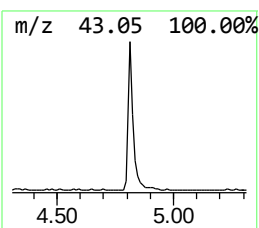
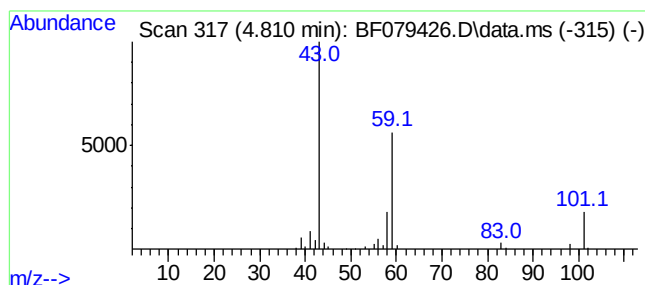
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.810	7.51 ng	131827	1,4-Dichlorobenzene-d4	7.073

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	32
3			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16



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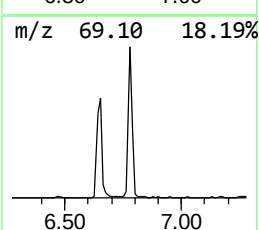
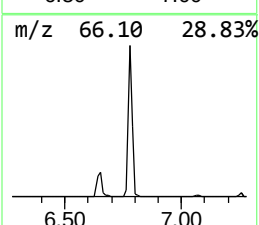
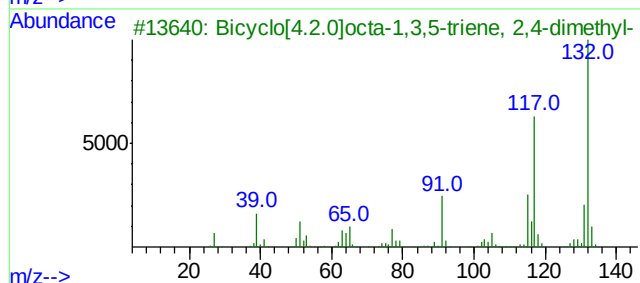
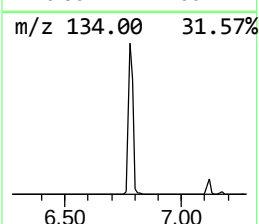
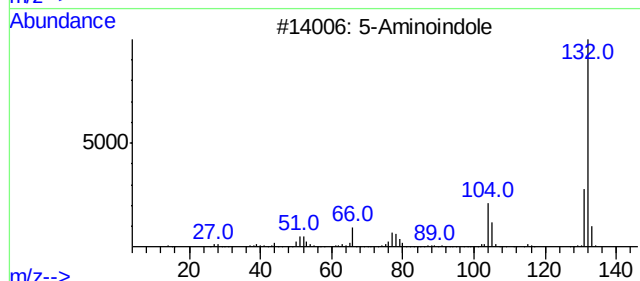
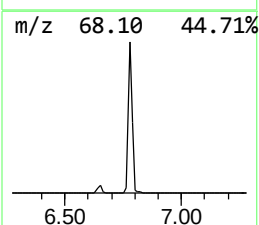
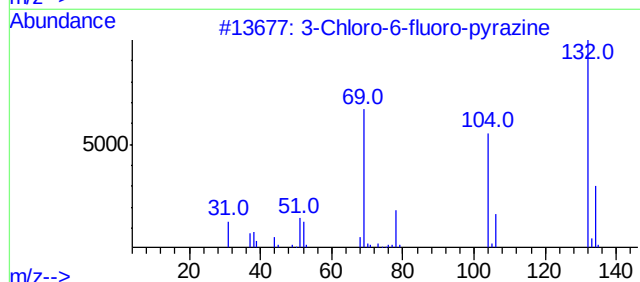
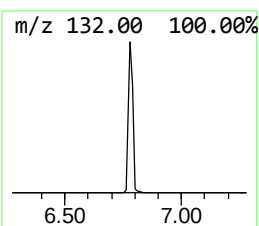
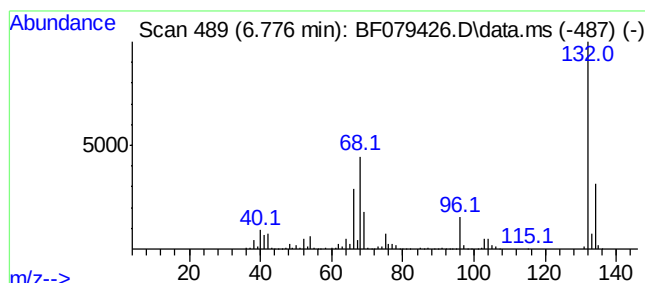
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown6.776 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.776	77.98 ng	1368020	1,4-Dichlorobenzene-d4	7.073

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole	132	C8H8N2	005192-03-0	12
3	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4	Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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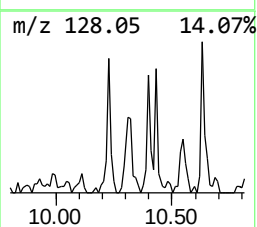
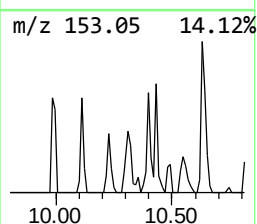
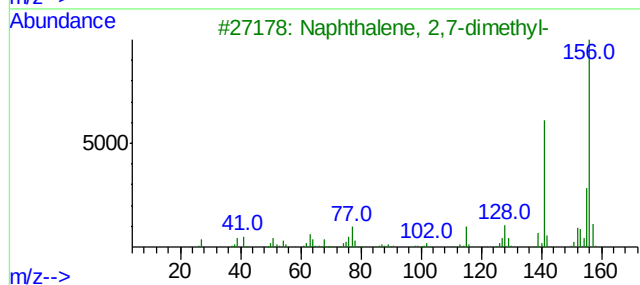
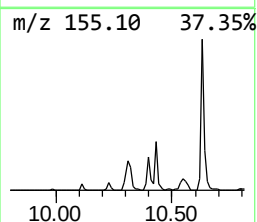
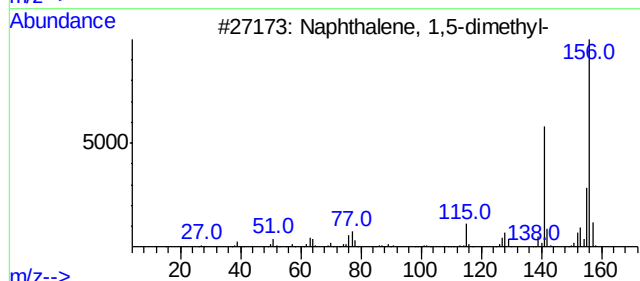
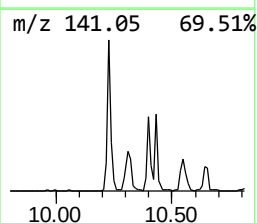
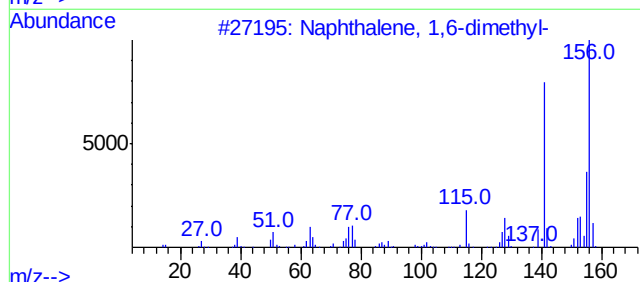
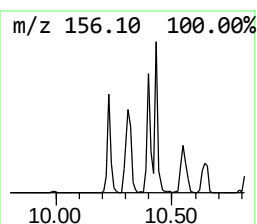
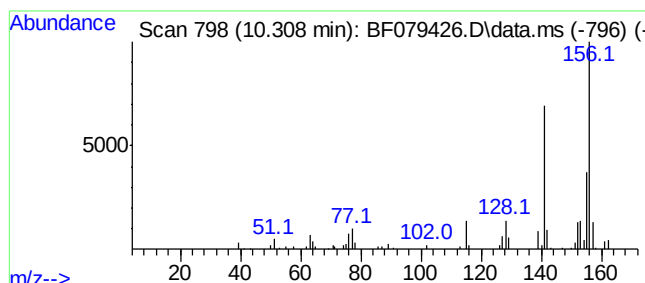
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Naphthalene, 1,6-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.308	7.81 ng	214177	Acenaphthene-d10	10.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079426.D
Acq On : 22 May 2015 7:51
Operator : TP/IZ
Sample : G2289-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-29D

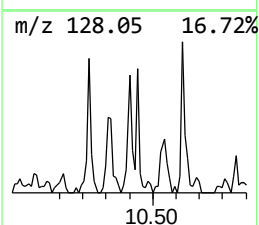
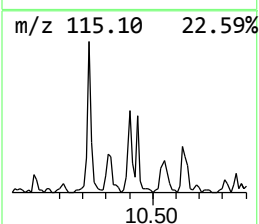
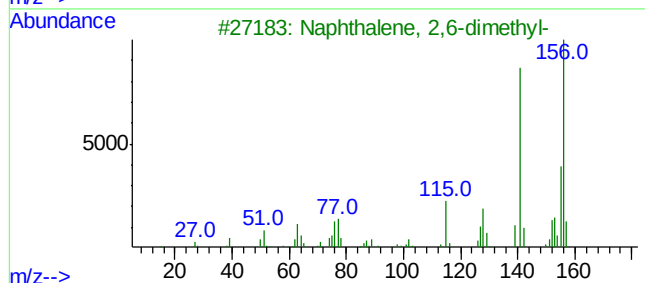
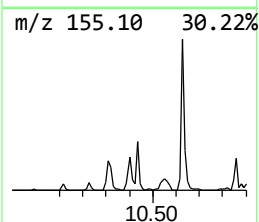
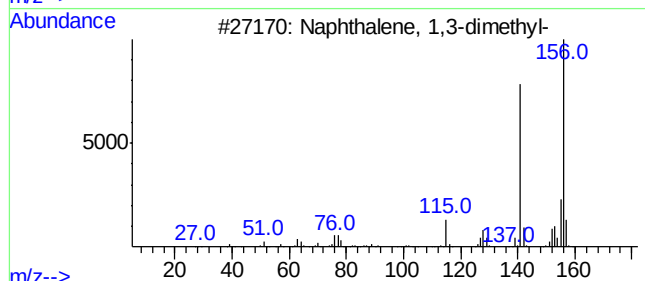
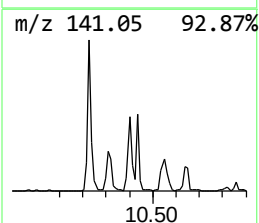
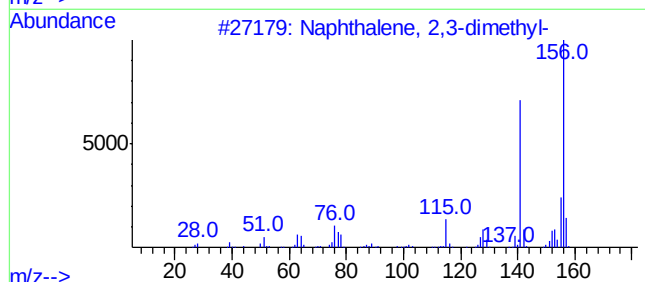
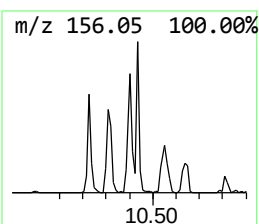
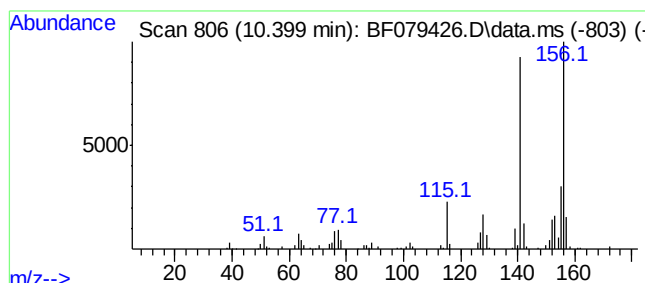
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 2,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.399	8.20 ng	224971	Acenaphthene-d10	10.811

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
3	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079426.D
Acq On : 22 May 2015 7:51
Operator : TP/IZ
Sample : G2289-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-29D

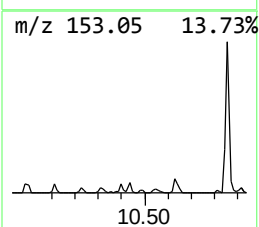
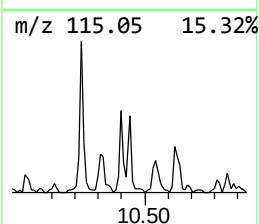
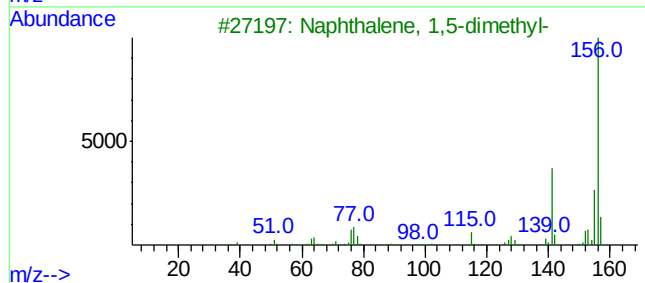
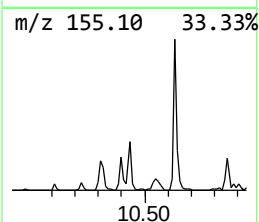
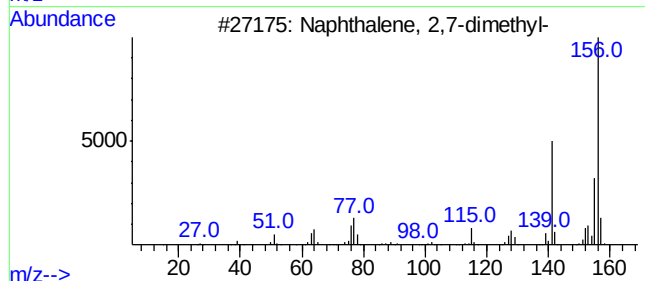
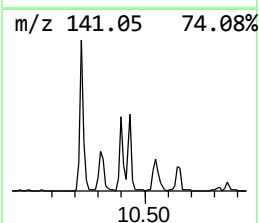
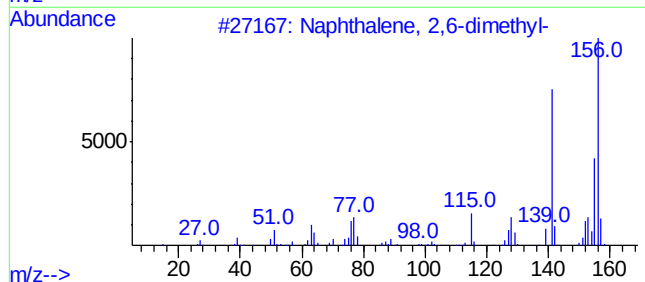
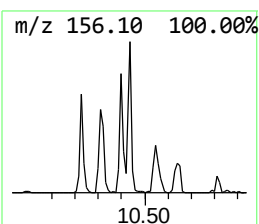
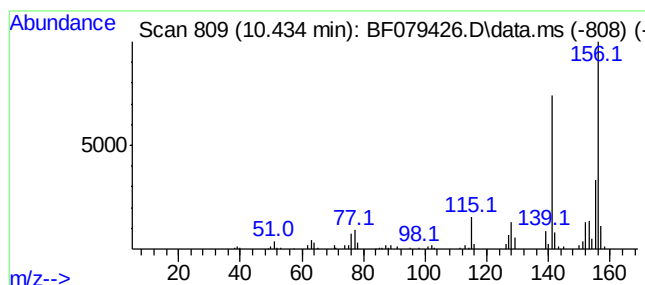
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, 2,6-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.434	6.28 ng	172190	Acenaphthene-d10	10.811

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
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Acq On : 22 May 2015 7:51
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Sample : G2289-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

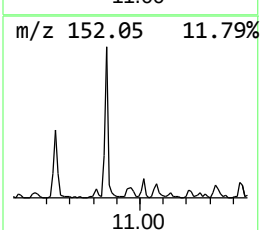
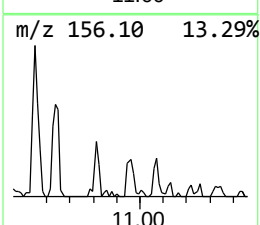
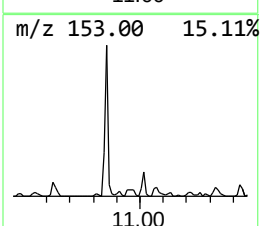
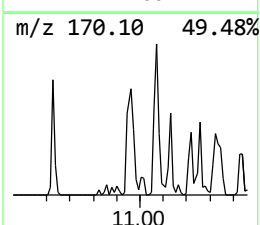
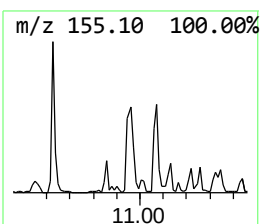
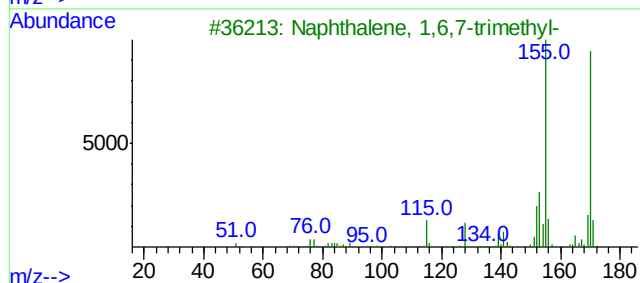
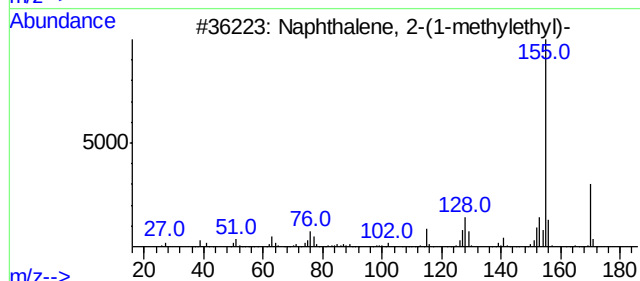
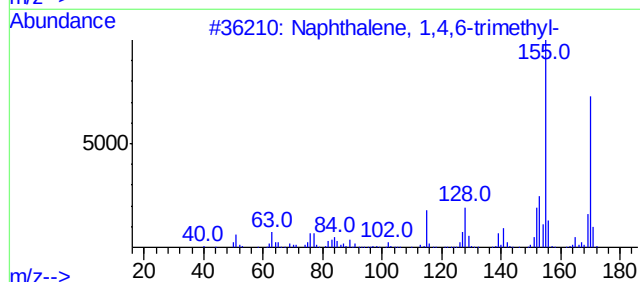
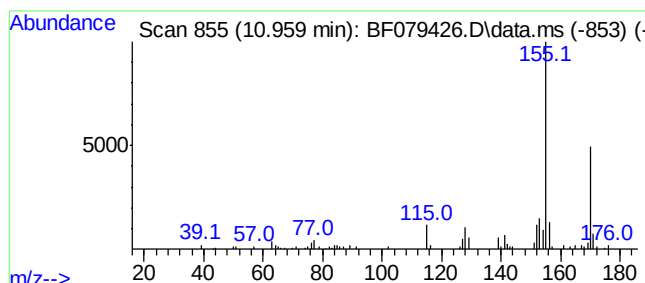
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Naphthalene, 1,4,6-trimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD			R.T.
10.959	6.90 ng	189245	Acenaphthene-d10			10.811
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene, 1,4,6-trimethyl-		170	C13H14	002131-42-2	93
2	Naphthalene, 2-(1-methylethyl)-		170	C13H14	002027-17-0	91
3	Naphthalene, 1,6,7-trimethyl-		170	C13H14	002245-38-7	89
4	3-(2-Methyl-propenyl)-1H-indene		170	C13H14	1000187-78-5	80
5	1'-Acetonaphthone		170	C12H10O	000941-98-0	72



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
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Misc :
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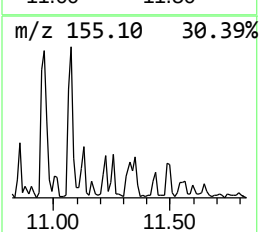
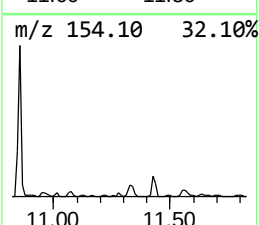
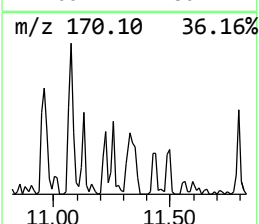
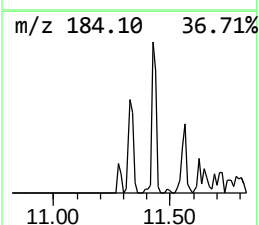
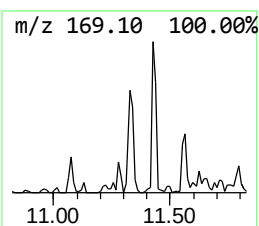
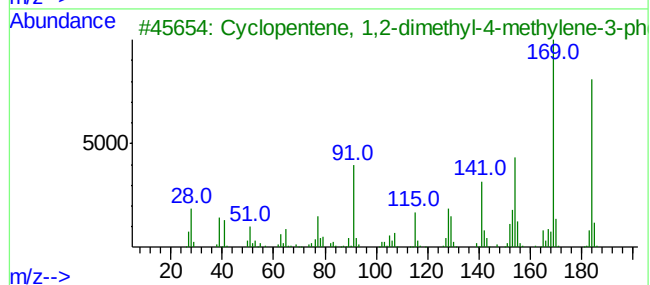
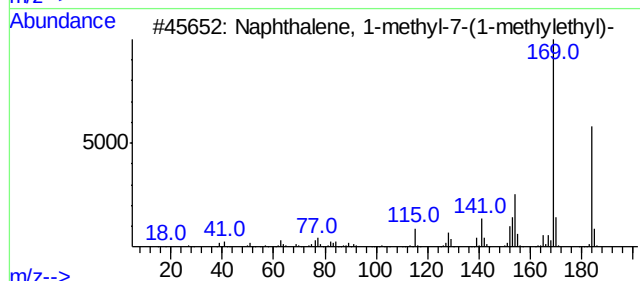
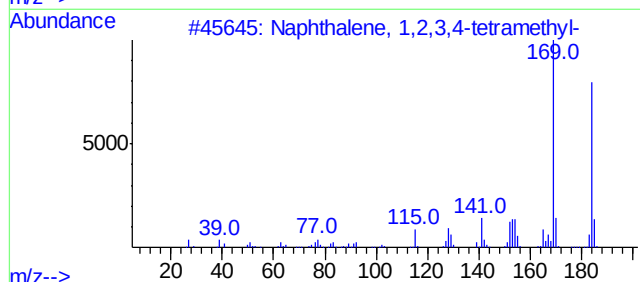
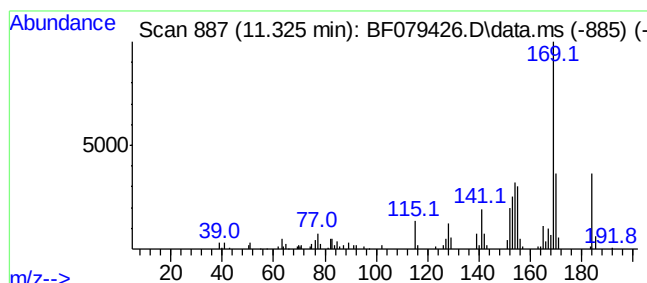
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Naphthalene, 1,2,3,4-tetram... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.325	7.73 ng	212008	Acenaphthene-d10	10.811	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetramethyl-	184	C14H16	003031-15-0	87
2	Naphthalene, 1-methyl-7-(1-methy...	184	C14H16	000490-65-3	81
3	Cyclopentene, 1,2-dimethyl-4-met...	184	C14H16	1000152-22-4	53
4	Naphthalene, 2-(1,1-dimethylethyl)-	184	C14H16	002876-35-9	47
5	[1,1'-Biphenyl]-4-amine	169	C12H11N	000092-67-1	45



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079426.D
Acq On : 22 May 2015 7:51
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Misc :
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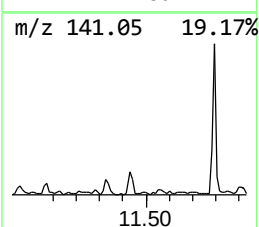
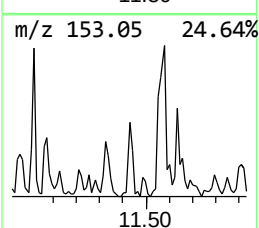
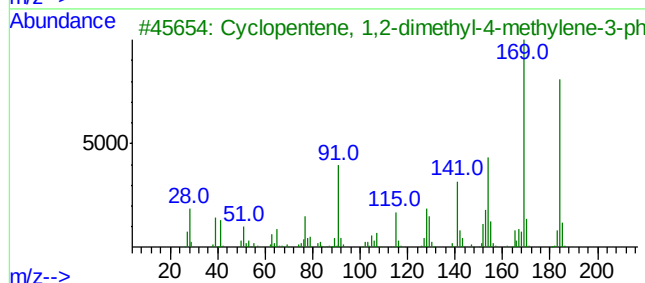
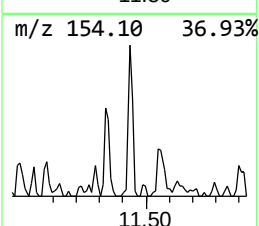
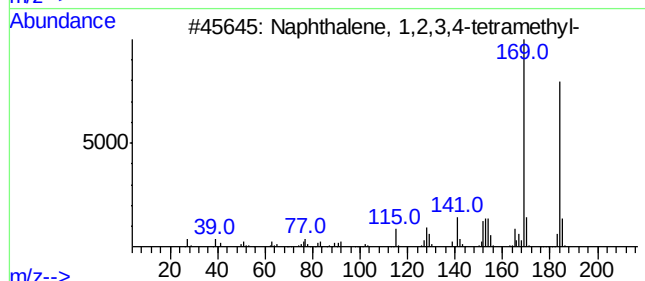
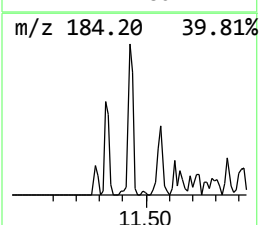
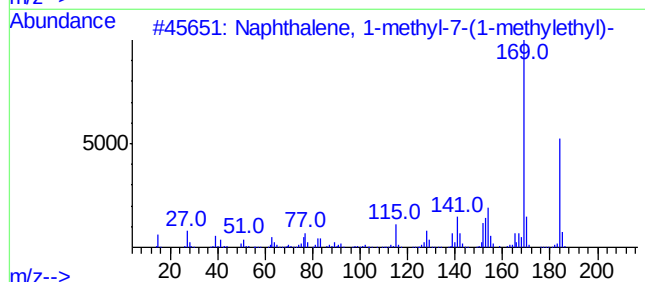
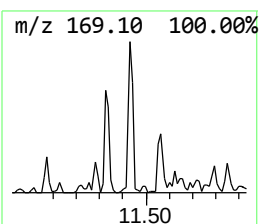
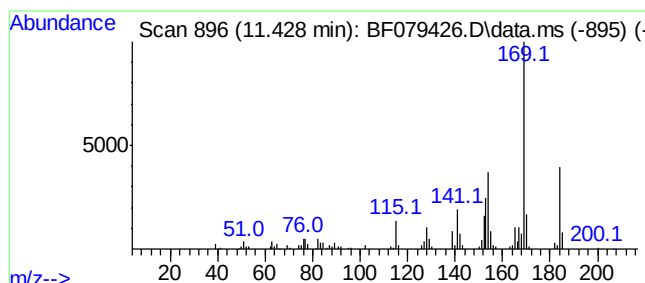
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Naphthalene, 1-methyl-7-(1-... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.428	7.85 ng	215312	Acenaphthene-d10	10.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-7-(1-methy...	184	C14H16	000490-65-3	96
2			Naphthalene, 1,2,3,4-tetramethyl-	184	C14H16	003031-15-0	94
3			Cyclopentene, 1,2-dimethyl-4-met...	184	C14H16	1000152-22-4	89
4			Naphthalene, 1-(1,1-dimethylethyl)-	184	C14H16	017085-91-5	74
5			Naphthalene, 2-(1,1-dimethylethyl)-	184	C14H16	002876-35-9	59



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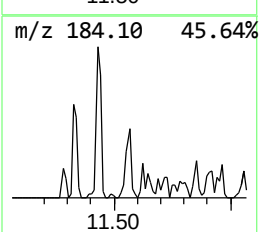
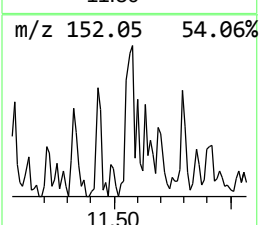
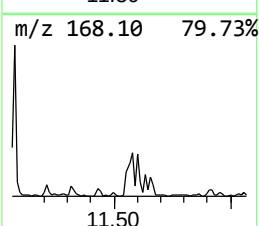
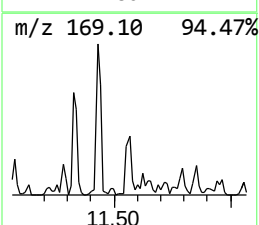
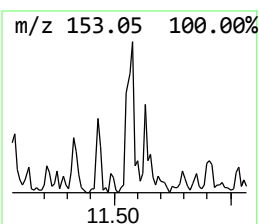
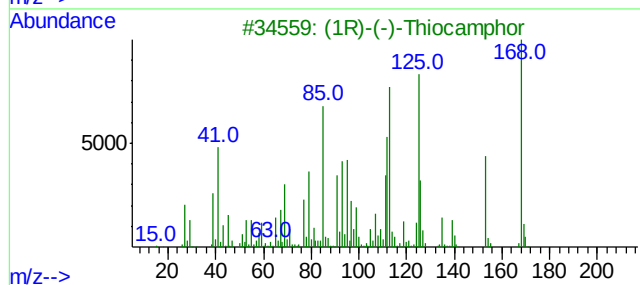
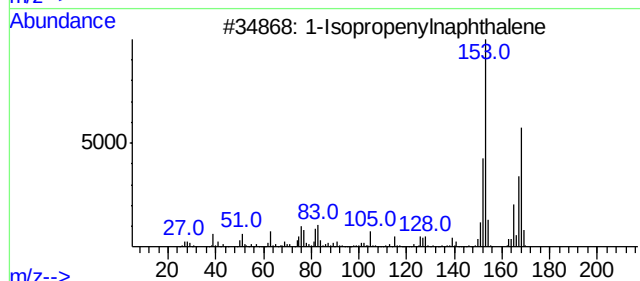
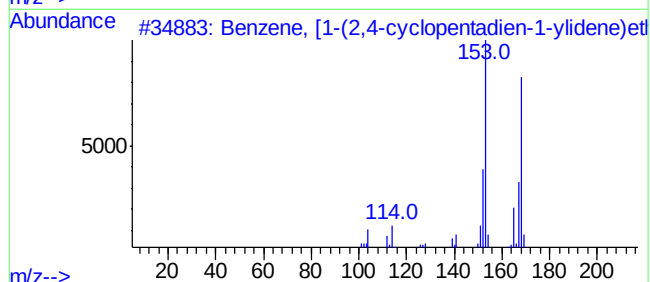
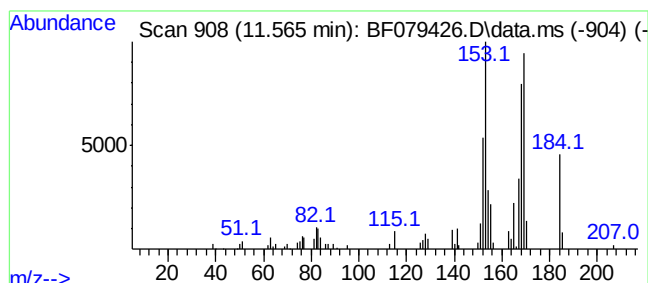
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown11.565 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.565	8.19 ng	224663	Acenaphthene-d10	10.811

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, [1-(2,4-cyclopentadien-...	168	C13H12	002320-32-3	45
2			1-Isopropenyl naphthalene	168	C13H12	001855-47-6	43
3			(1R)-(-)-Thiocamphor	168	C10H16S	053402-10-1	35
4			1,2-Dihydro-5-acenaphthylenamine	169	C12H11N	004657-93-6	35
5			Cyclopentene, 1,2-dimethyl-4-met...	184	C14H16	1000152-22-4	35



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079426.D
Acq On : 22 May 2015 7:51
Operator : TP/IZ
Sample : G2289-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-29D

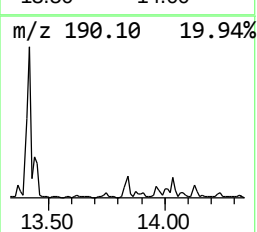
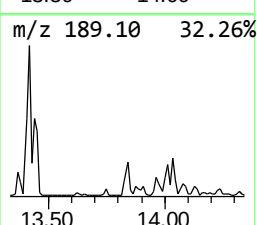
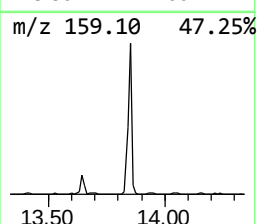
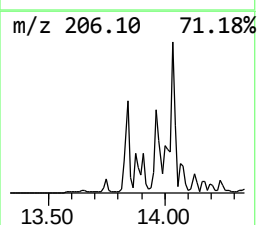
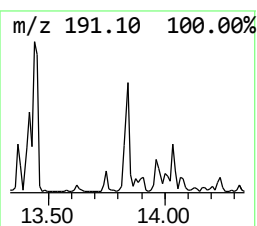
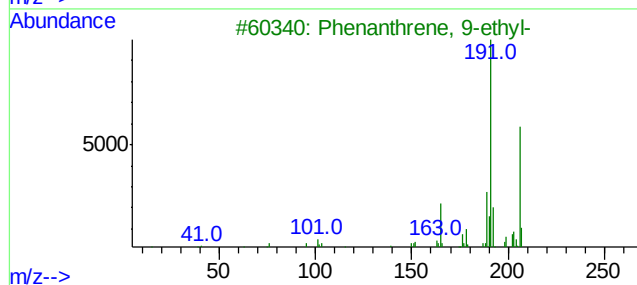
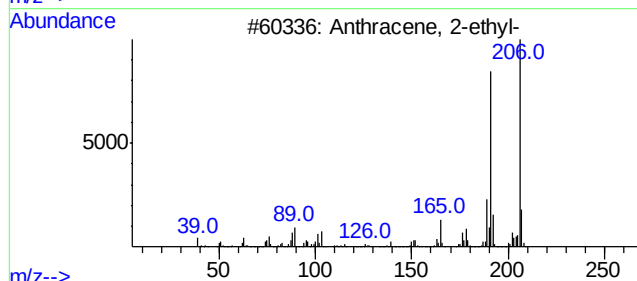
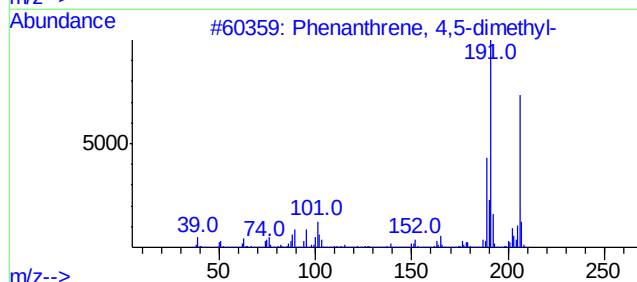
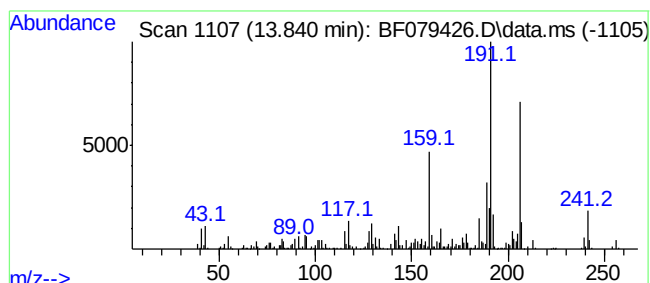
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Phenanthrene, 4,5-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.840	4.84 ng	711192	Phenanthrene-d10	12.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	93
2	Anthracene, 2-ethyl-	206	C16H14	052251-71-5	83
3	Phenanthrene, 9-ethyl-	206	C16H14	003674-75-7	62
4	Anthracene, 9-ethyl-	206	C16H14	000605-83-4	60
5	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	49



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079426.D
Acq On : 22 May 2015 7:51
Operator : TP/IZ
Sample : G2289-08
Misc :
ALS Vial : 23 Sample Multiplier: 1

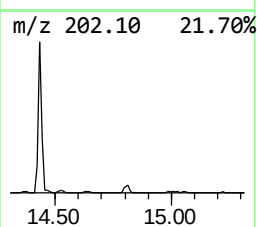
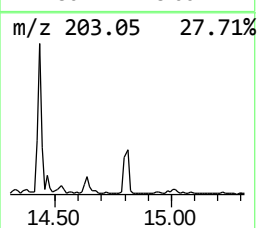
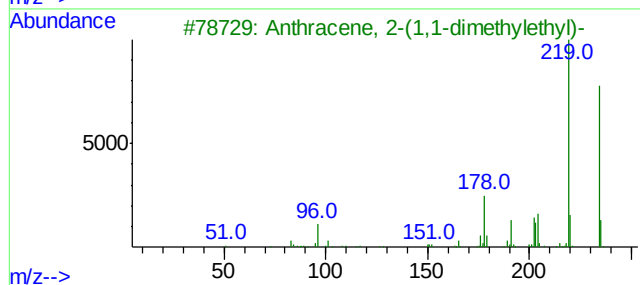
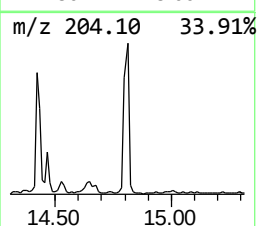
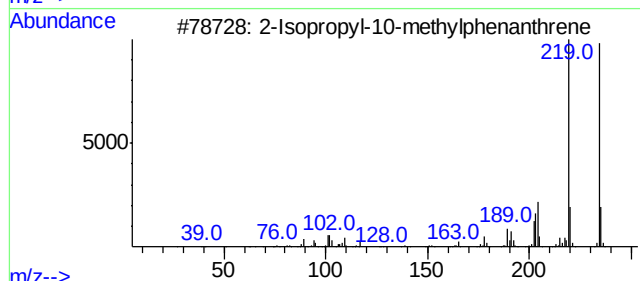
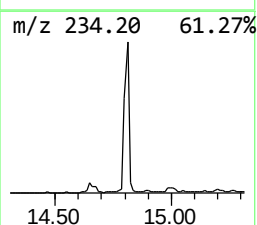
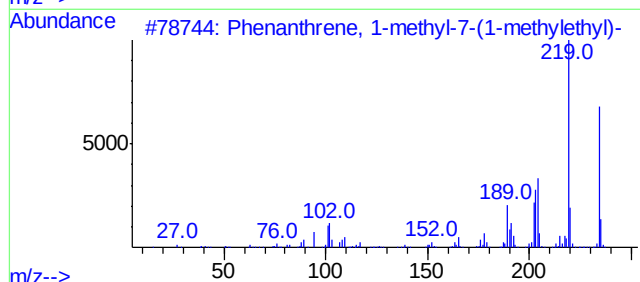
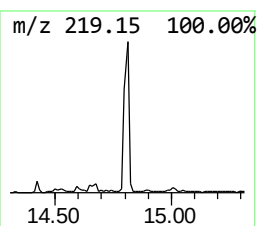
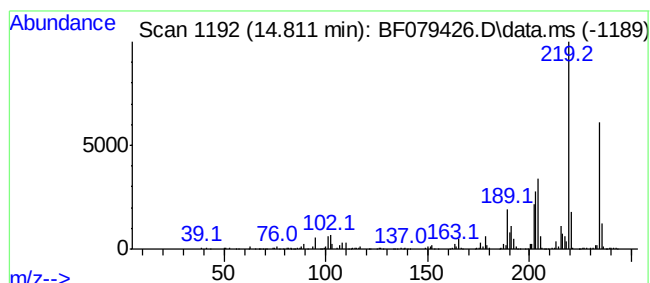
Instrument :
BNA_F
ClientSampleId :
SB-29D

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Phenanthrene, 1-methyl-7-(1... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.811	11.29 ng	986034	Chrysene-d12			15.943
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-7-(1-meth...		234	C18H18	000483-65-8	98
2	2-Isopropyl-10-methylphenanthrene		234	C18H18	066552-97-4	93
3	Anthracene, 2-(1,1-dimethylethyl)-		234	C18H18	018801-00-8	64
4	1-(10-Methylanthracen-9-yl)ethanone		234	C17H14O	036778-18-4	58
5	Phenanthrene, 2,4,5,7-tetramethyl-		234	C18H18	007396-38-5	58



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
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 Operator : TP/IZ
 Sample : G2289-08
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-29D

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.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
Butane, 2-metho...	1.496	19.0	ng	333735	1	7.073	350872	20.0
2-Pentanone, 4-...	4.810	7.5	ng	131827	1	7.073	350872	20.0
unknown6.776	6.776	78.0	ng	1368020	1	7.073	350872	20.0
Naphthalene, 1,...	10.308	7.8	ng	214177	3	10.811	548588	20.0
Naphthalene, 2,...	10.399	8.2	ng	224971	3	10.811	548588	20.0
Naphthalene, 2,...	10.434	6.3	ng	172190	3	10.811	548588	20.0
Naphthalene, 1,...	10.959	6.9	ng	189245	3	10.811	548588	20.0
Naphthalene, 1,...	11.325	7.7	ng	212008	3	10.811	548588	20.0
Naphthalene, 1-...	11.428	7.8	ng	215312	3	10.811	548588	20.0
unknown11.565	11.565	8.2	ng	224663	3	10.811	548588	20.0
Phenanthrene, 4...	13.840	4.8	ng	711192	4	12.651	2935860	20.0
Phenanthrene, 1...	14.811	11.3	ng	986034	5	15.943	1746890	20.0