

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
 Data File : BF079491.D
 Acq On : 26 May 2015 22:50
 Operator : TP/IZ
 Sample : G2289-10 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-30D

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-BF052615.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.278	6	8	19	rVB2	130551	321615	5.48%	0.981%
2	1.427	19	21	34	rVB	107083	296860	5.06%	0.905%
3	1.587	34	35	43	rVB	66035	149564	2.55%	0.456%
4	1.701	43	45	88	rVB	2565292	5870717	100.00%	17.902%
5	5.302	358	360	372	rBV	286770	380350	6.48%	1.160%
6	6.616	472	475	483	rBV	296411	405716	6.91%	1.237%
7	6.730	483	485	488	rBV	394406	422637	7.20%	1.289%
8	7.016	508	510	513	rBV	621639	569668	9.70%	1.737%
9	7.199	524	526	529	rBV	276904	274926	4.68%	0.838%
10	7.713	569	571	576	rBV	239486	234223	3.99%	0.714%
11	8.593	645	648	650	rBV	907892	856957	14.60%	2.613%
12	9.599	734	736	741	rVB	92937	86127	1.47%	0.263%
13	9.942	764	766	768	rBV	460171	397696	6.77%	1.213%
14	10.262	791	794	796	rVB	35565	61575	1.05%	0.188%
15	10.353	799	802	803	rBV	65256	94847	1.62%	0.289%
16	10.491	812	814	817	rVB	49287	67606	1.15%	0.206%
17	10.765	836	838	840	rVV	811955	951940	16.22%	2.903%
18	10.799	840	841	845	rVB	371439	310837	5.29%	0.948%
19	11.016	857	860	863	rVV	167169	175509	2.99%	0.535%
20	11.074	863	865	871	rVV2	33658	78074	1.33%	0.238%
21	11.279	880	883	888	rBV2	23734	61478	1.05%	0.187%
22	11.439	895	897	900	rVB	267813	237643	4.05%	0.725%
23	11.519	900	904	905	rBV	39735	68965	1.17%	0.210%
24	11.645	913	915	918	rVB	68287	84196	1.43%	0.257%
25	11.736	921	923	926	rVV2	287540	412785	7.03%	1.259%
26	12.125	952	957	959	rVV	67821	108051	1.84%	0.329%
27	12.228	961	966	967	rVB3	36738	74007	1.26%	0.226%
28	12.308	970	973	975	rBV3	27208	65272	1.11%	0.199%
29	12.422	981	983	985	rBV	42875	60455	1.03%	0.184%
30	12.468	985	987	989	rVV	110112	101285	1.73%	0.309%
31	12.594	996	998	1000	rBV2	770216	1236754	21.07%	3.771%
32	12.628	1000	1001	1004	rVV	1886984	1510933	25.74%	4.607%
33	12.697	1004	1007	1009	rVB	403732	572621	9.75%	1.746%
34	12.902	1023	1025	1028	rVV	94637	117805	2.01%	0.359%

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
 Data File : BF079491.D
 Acq On : 26 May 2015 22:50
 Operator : TP/IZ
 Sample : G2289-10 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-30D

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-BF052615.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	13.222	1051	1053	1055	rBV	188574	226763	3.86%	0.691%
36	13.257	1055	1056	1059	rVB	211253	233237	3.97%	0.711%
37	13.314	1059	1061	1063	rBV	98571	134480	2.29%	0.410%
38	13.359	1063	1065	1067	rVV	345901	475651	8.10%	1.450%
39	13.394	1067	1068	1071	rVB	106086	85079	1.45%	0.259%
40	13.599	1084	1086	1091	rVB	129996	171275	2.92%	0.522%
41	13.782	1100	1102	1104	rBV2	47500	59378	1.01%	0.181%
42	13.908	1110	1113	1115	rBV	106915	164020	2.79%	0.500%
43	13.954	1115	1117	1121	rVB2	62129	104466	1.78%	0.319%
44	14.022	1121	1123	1126	rVB3	46584	75784	1.29%	0.231%
45	14.102	1127	1130	1132	rBV	1869677	1830395	31.18%	5.582%
46	14.171	1134	1136	1139	rBV3	154585	191157	3.26%	0.583%
47	14.285	1143	1146	1147	rBV	80489	118279	2.01%	0.361%
48	14.377	1151	1154	1156	rBV	1626213	1848474	31.49%	5.637%
49	14.445	1159	1160	1166	rVB2	67882	114128	1.94%	0.348%
50	14.537	1166	1168	1170	rBV	86582	104299	1.78%	0.318%
51	14.594	1170	1173	1175	rVB2	381528	488973	8.33%	1.491%
52	14.662	1176	1179	1181	rBV2	69978	135313	2.30%	0.413%
53	14.800	1185	1191	1193	rVB	226057	403848	6.88%	1.231%
54	14.880	1193	1198	1200	rBV	271628	375101	6.39%	1.144%
55	14.925	1200	1202	1204	rBV	141029	200787	3.42%	0.612%
56	15.028	1210	1211	1213	rBV	72476	91784	1.56%	0.280%
57	15.074	1213	1215	1218	rVV	65121	91549	1.56%	0.279%
58	15.337	1233	1238	1240	rVV2	40459	124400	2.12%	0.379%
59	15.405	1242	1244	1245	rBV	45480	65385	1.11%	0.199%
60	15.565	1255	1258	1259	rBV	78983	107896	1.84%	0.329%
61	15.600	1259	1261	1264	rVV2	110671	194415	3.31%	0.593%
62	15.645	1264	1265	1268	rVB2	48854	67451	1.15%	0.206%
63	15.771	1274	1276	1278	rBV2	56133	91701	1.56%	0.280%
64	15.874	1281	1285	1287	rVV2	1231239	2134328	36.36%	6.508%
65	15.908	1287	1288	1290	rVV	597660	502466	8.56%	1.532%
66	16.000	1293	1296	1298	rBV2	160993	205337	3.50%	0.626%
67	16.137	1306	1308	1310	rVB	58238	64720	1.10%	0.197%
68	16.171	1310	1311	1316	rVB2	65537	119066	2.03%	0.363%
69	16.320	1322	1324	1326	rBV3	43006	64415	1.10%	0.196%
70	16.365	1326	1328	1330	rBV	93015	147235	2.51%	0.449%
71	16.411	1330	1332	1333	rBV2	74534	95962	1.63%	0.293%

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-30D

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

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	17.154	1395	1397	1403	rBV	521250	1101302	18.76%	3.358%
73	17.268	1406	1407	1409	rVB	109991	123846	2.11%	0.378%
74	17.474	1423	1425	1428	rBV	329001	438644	7.47%	1.338%
75	17.531	1428	1430	1433	rBV	534980	662677	11.29%	2.021%
76	17.600	1433	1436	1438	rBV	878346	1135783	19.35%	3.463%
77	18.926	1549	1552	1556	rBV2	241149	464848	7.92%	1.418%
78	19.314	1583	1586	1590	rBV	209263	437634	7.45%	1.335%

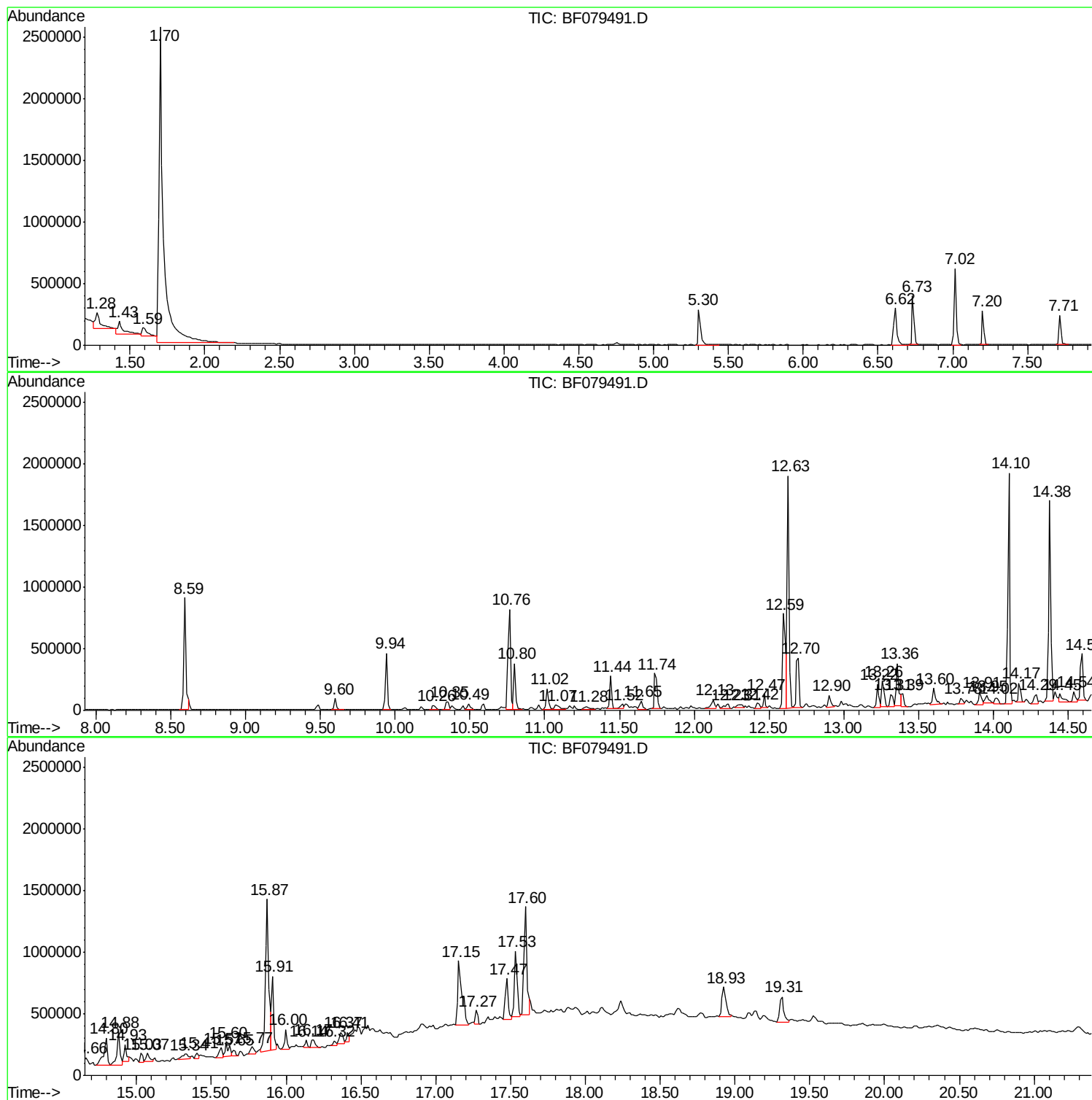
Sum of corrected areas: 32793425

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-30D

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-30D

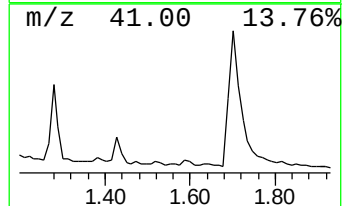
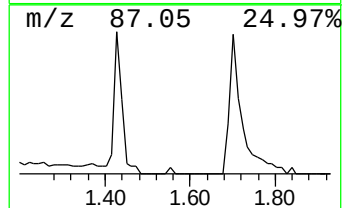
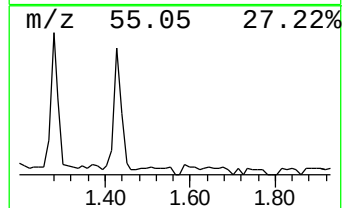
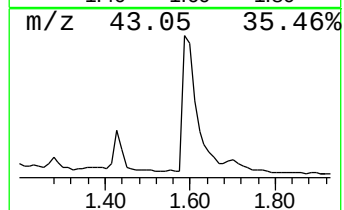
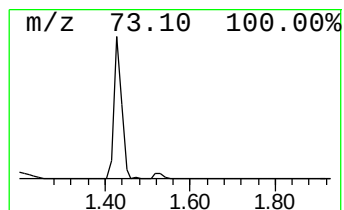
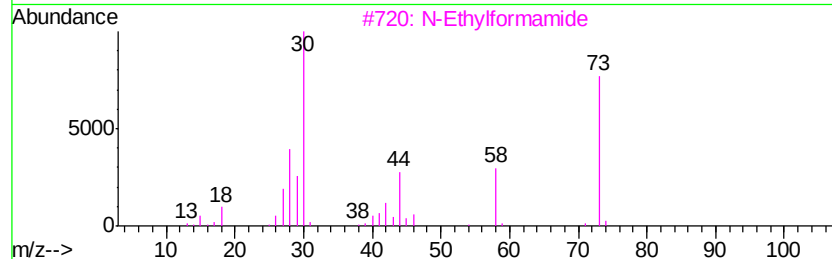
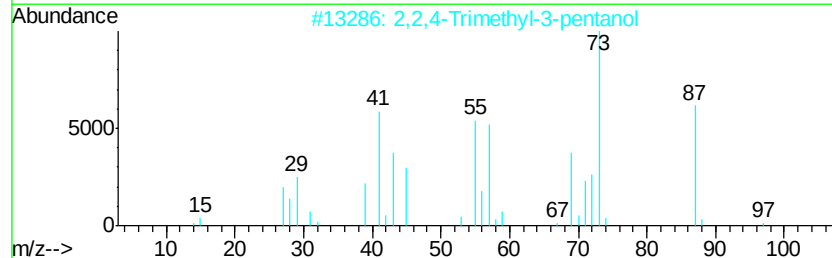
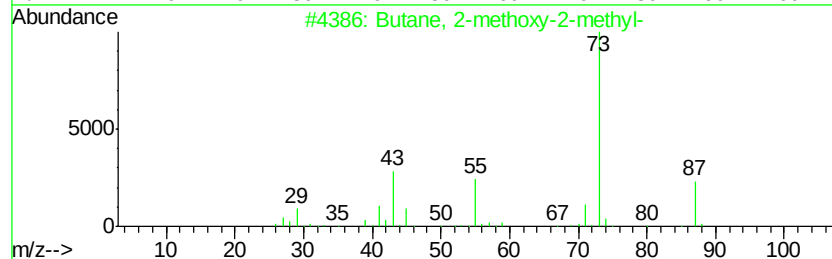
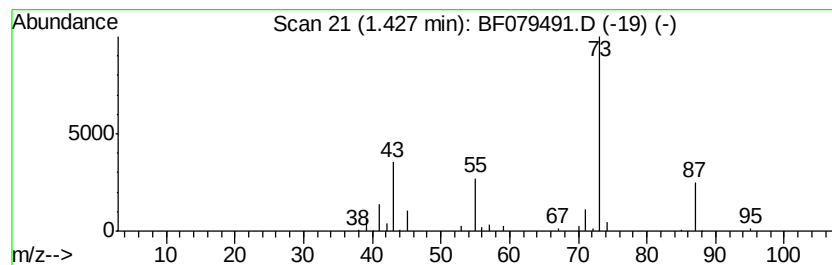
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.43	10.42 ng	296860	1,4-Dichlorobenzene-d4	7.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3		N-Ethylformamide	73	C3H7NO	000627-45-2	9
4		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5		Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-30D

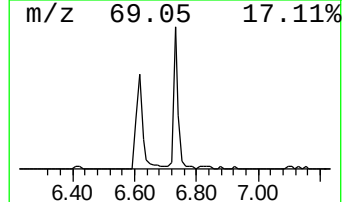
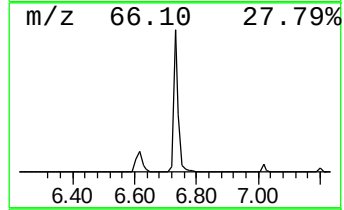
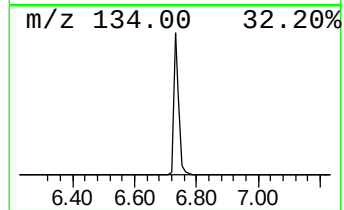
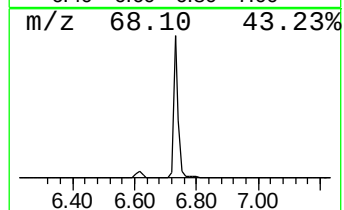
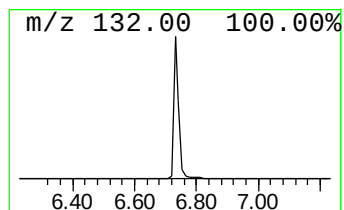
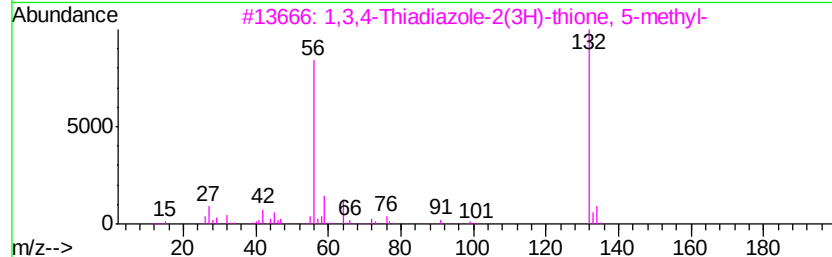
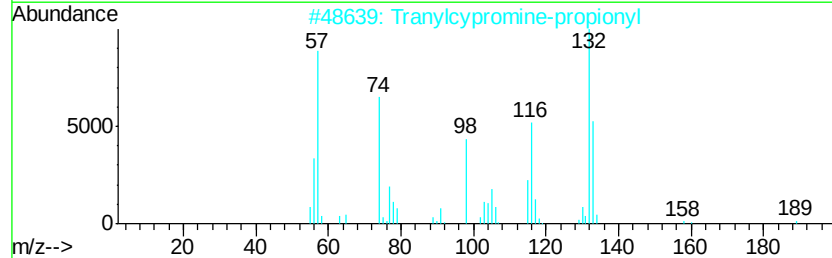
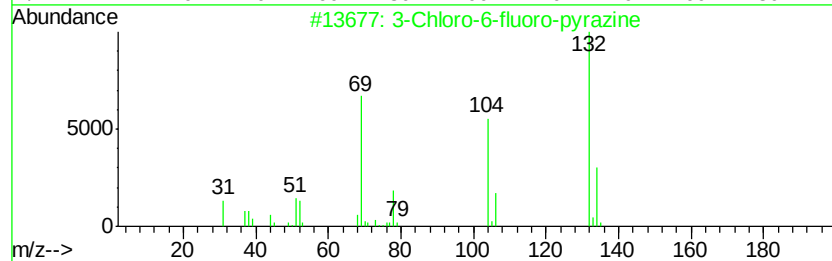
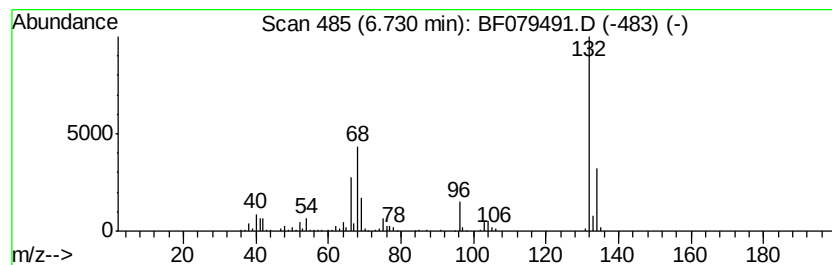
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 unknown6.73 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.73	14.84 ng	422637	1,4-Dichlorobenzene-d4	7.02

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-30D

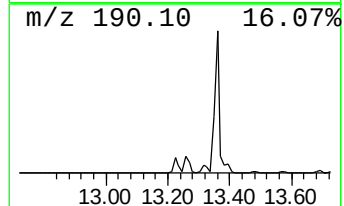
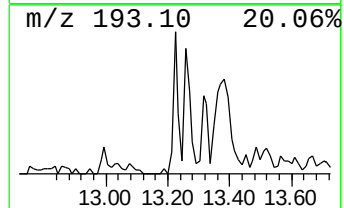
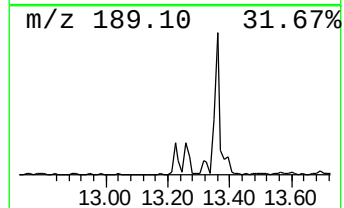
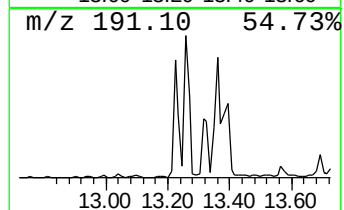
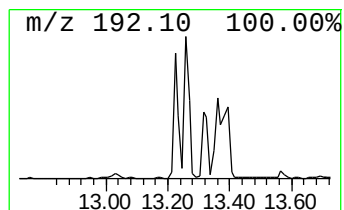
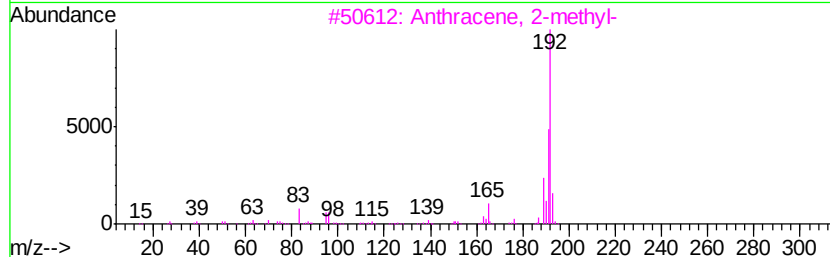
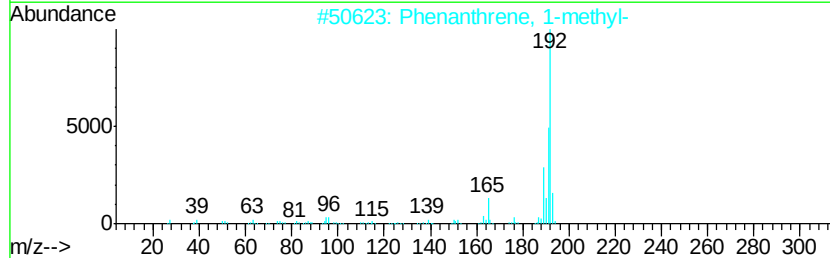
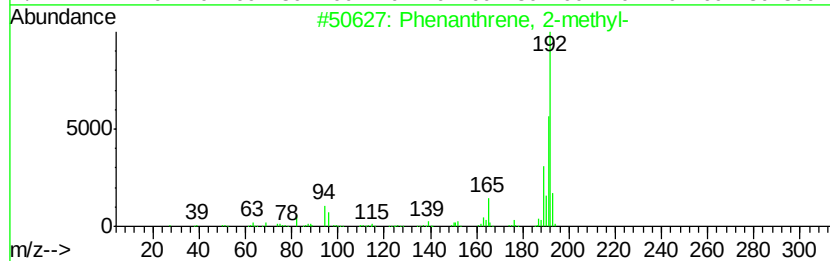
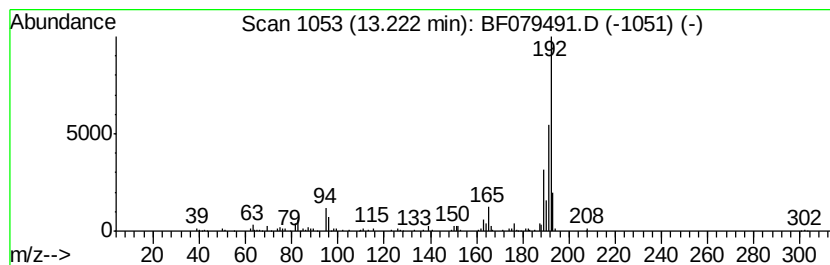
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenanthrene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.22	3.67 ng	226763	Phenanthrene-d10	12.59

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4		Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
5		1H-Cyclopropa[l]phenanthrene, 1a, ...	192	C15H12	000949-41-7	93



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

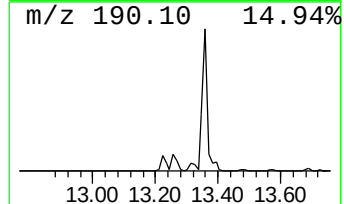
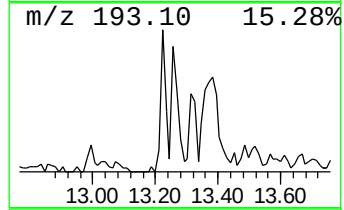
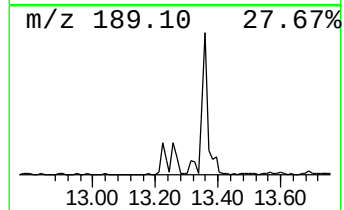
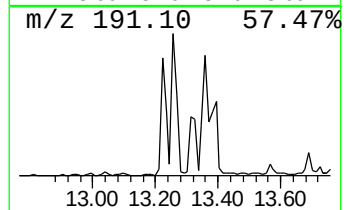
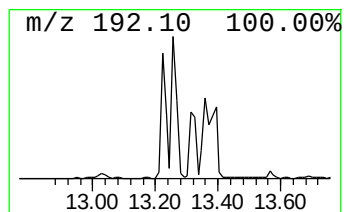
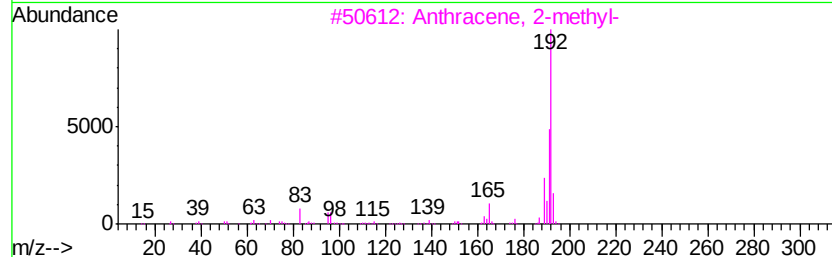
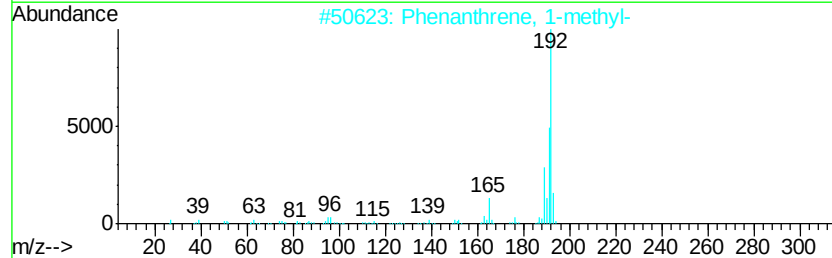
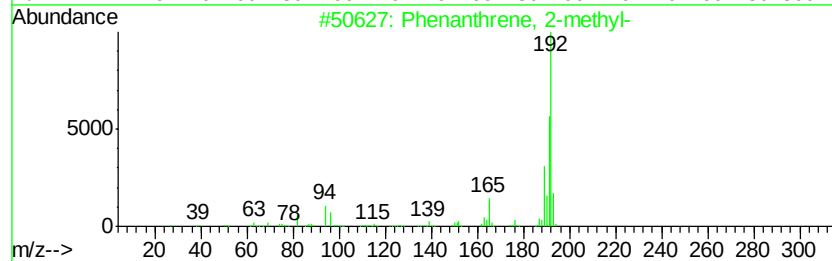
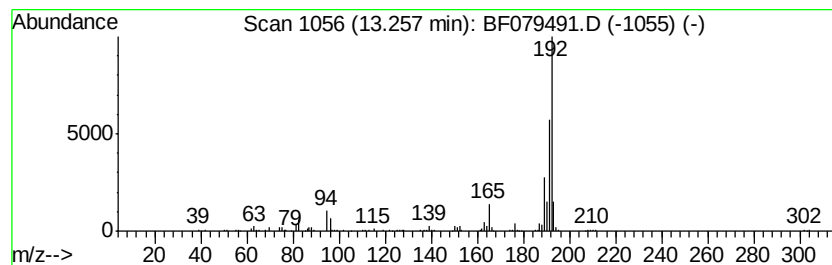
Instrument :
BNA_F
ClientSampleId :
SB-30D

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phenanthrene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.26	3.77 ng	233237	Phenanthrene-d10	12.59	
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	97
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4	1H-Cyclopropa[1,1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-30D

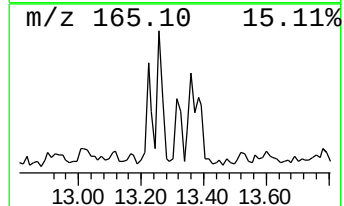
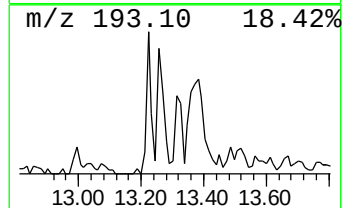
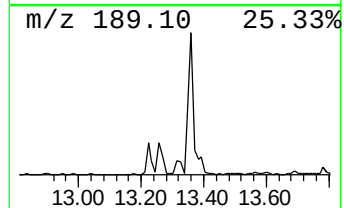
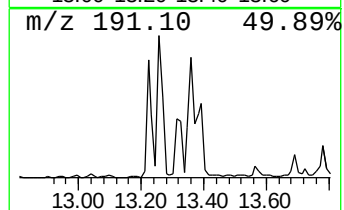
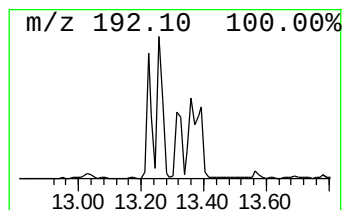
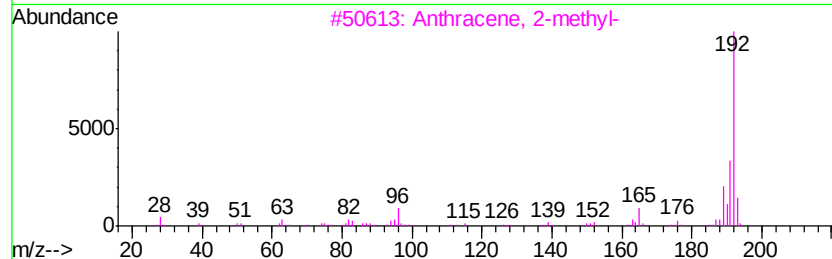
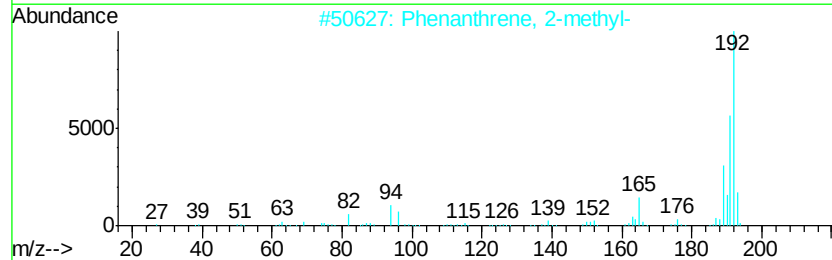
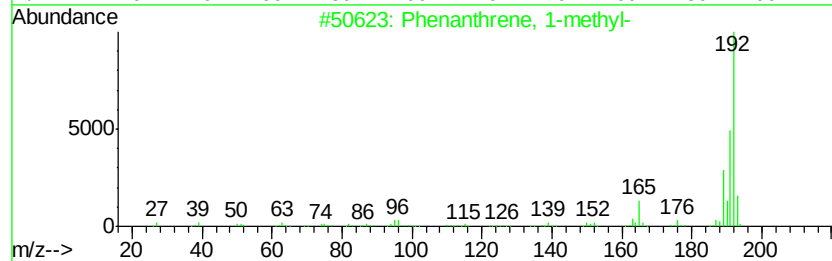
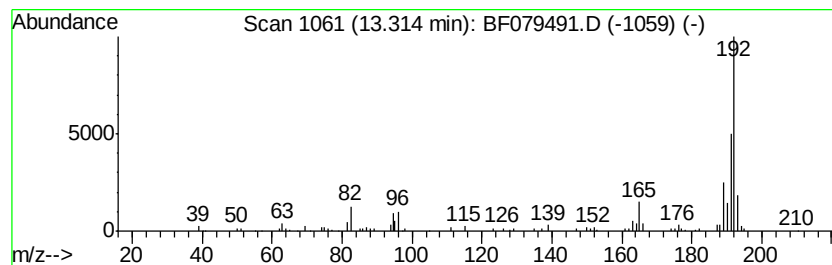
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Anthracene, 2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.31	2.17 ng	134480	Phenanthrene-d10	12.59

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4		1H-Cyclopropa[1,1]phenanthrene, 1a,...	192	C15H12	000949-41-7	96
5		Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	95



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

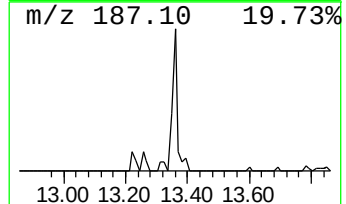
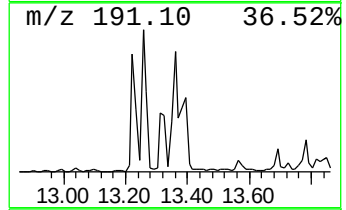
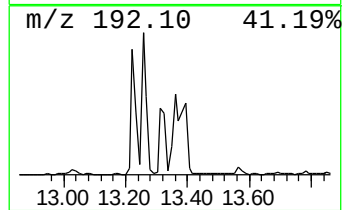
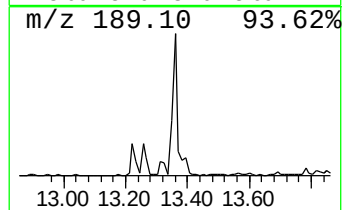
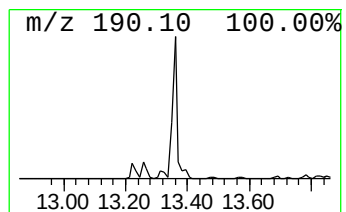
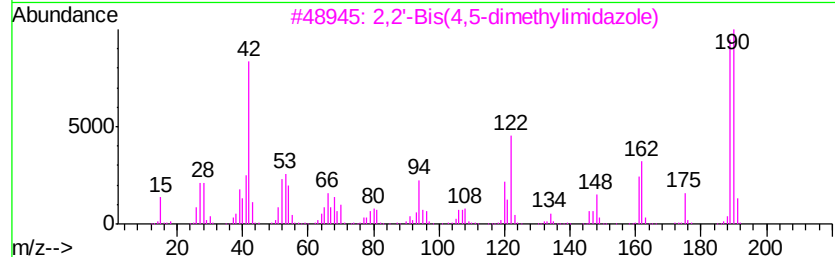
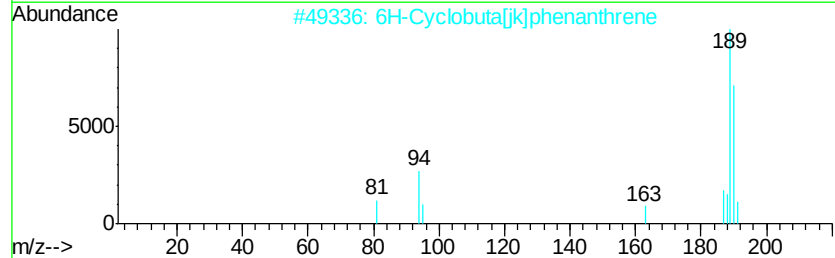
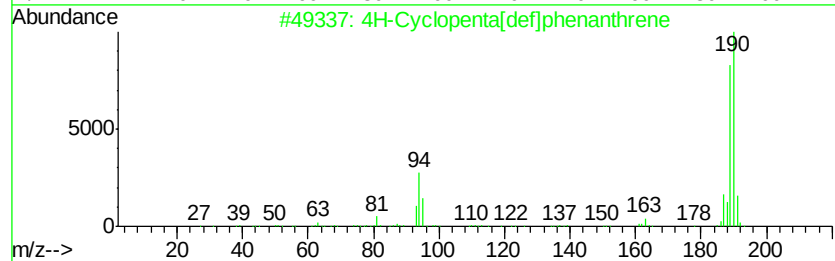
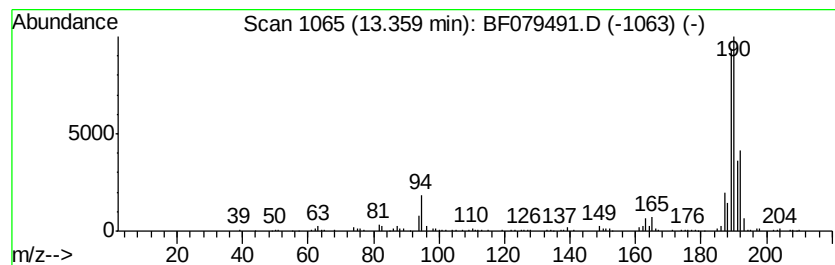
Instrument :
BNA_F
ClientSampleId :
SB-30D

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.36	7.69 ng	475651	Phenanthrene-d10	12.59	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	6H-Cyclobuta[ik]phenanthrene	190	C15H10	083469-43-6	45
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	42
4	4-Isothiazolecarbonitrile, 5-chl...	190	C5H3ClN2S2	054798-93-5	25
5	4,4'-Difluorobiphenyl	190	C12H8F2	000398-23-2	17



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

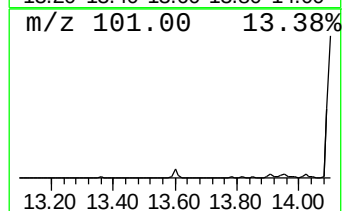
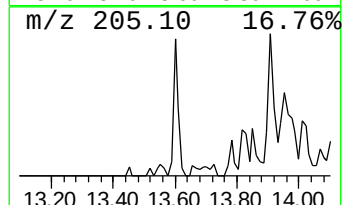
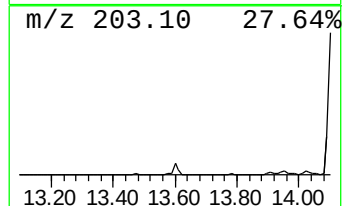
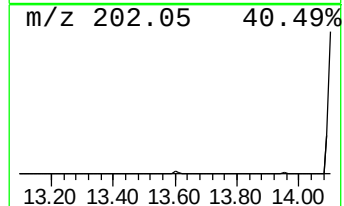
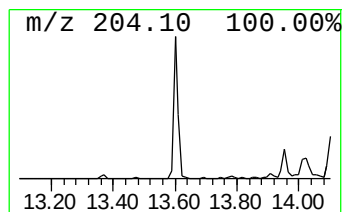
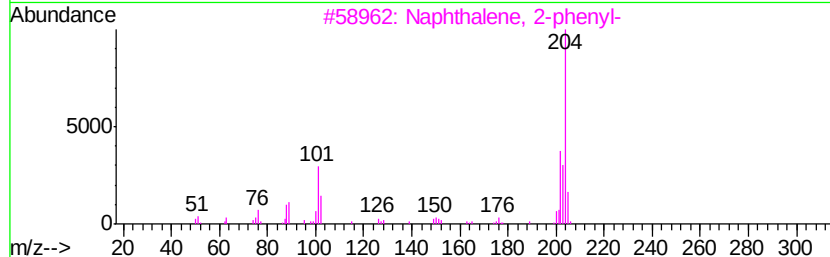
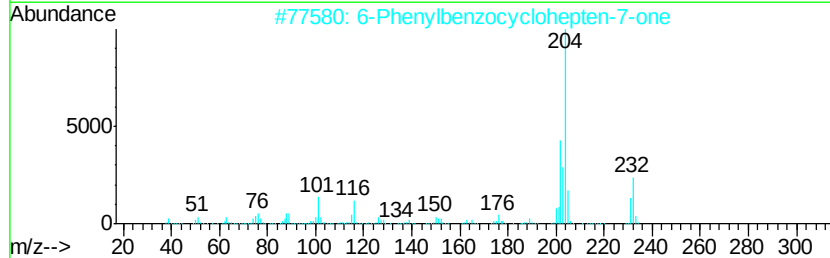
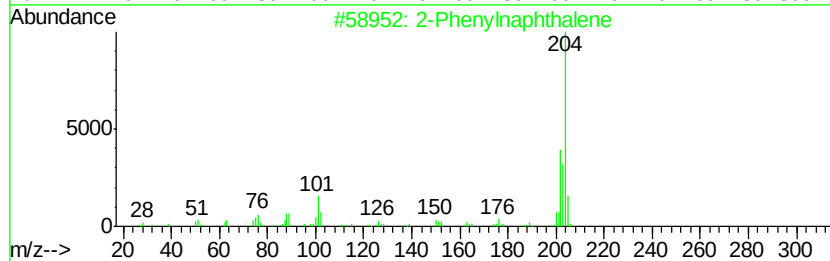
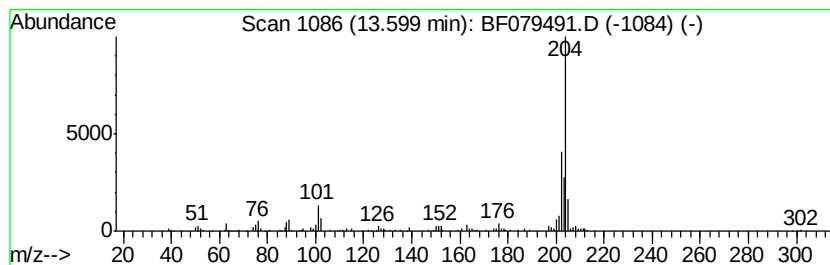
Instrument :
BNA_F
ClientSampleId :
SB-30D

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2-Phenylnaphthalene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.60	2.77 ng	171275	Phenanthrene-d10	12.59	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenylnaphthalene	204	C16H12	035465-71-5	94
2	6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	91
3	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86
4	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	86
5	Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	70



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-30D

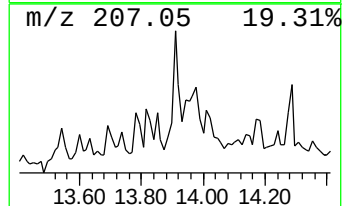
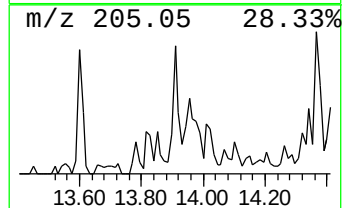
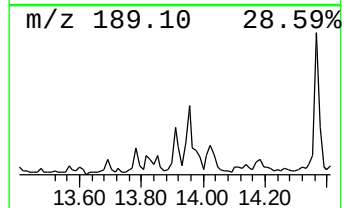
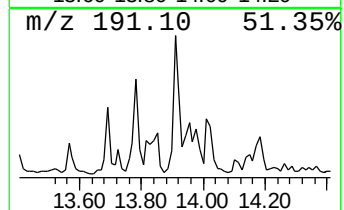
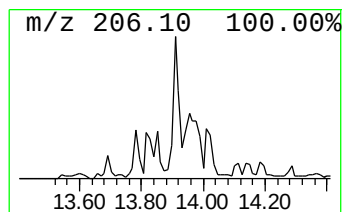
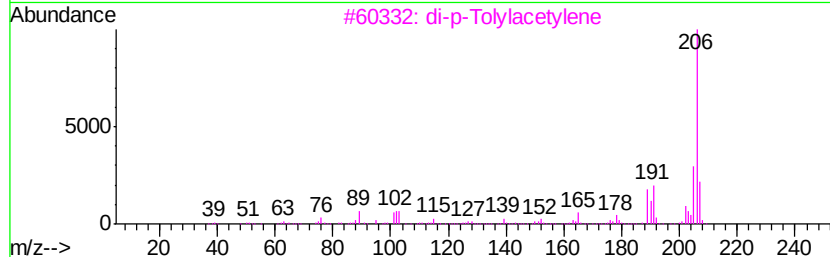
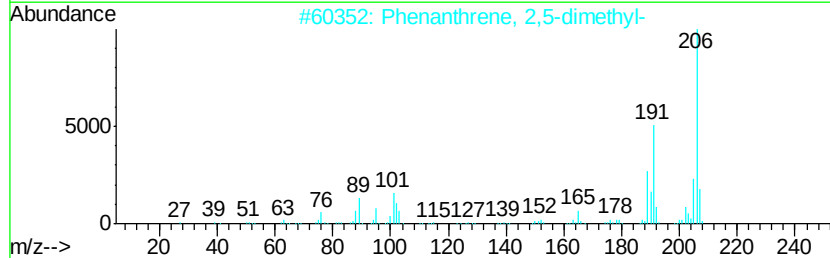
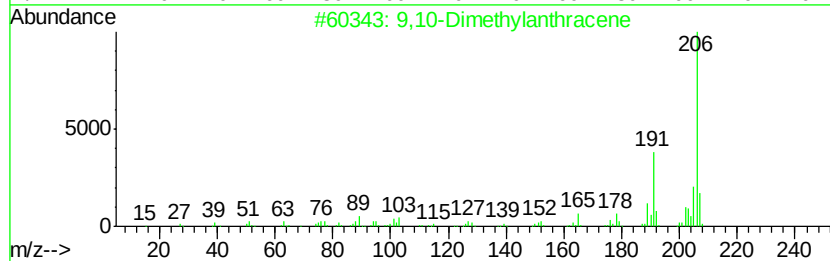
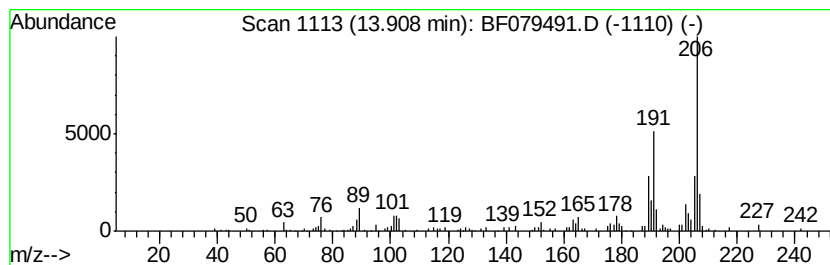
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 9,10-Dimethylantracene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.91	2.65 ng	164020	Phenanthrene-d10	12.59

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	94
3			di-p-Tolylacetylene	206	C16H14	002789-88-0	94
4			Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
5			Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-30D

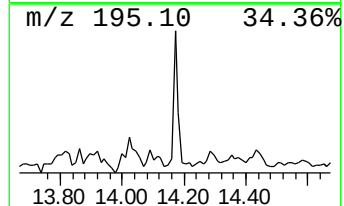
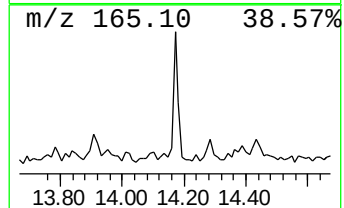
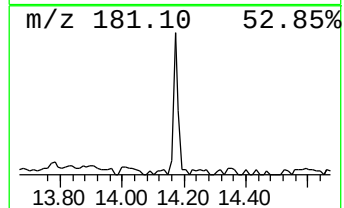
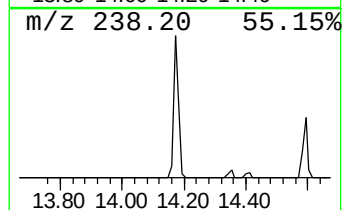
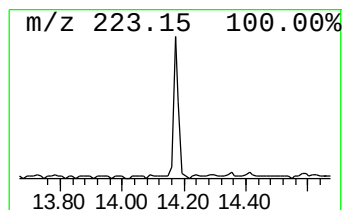
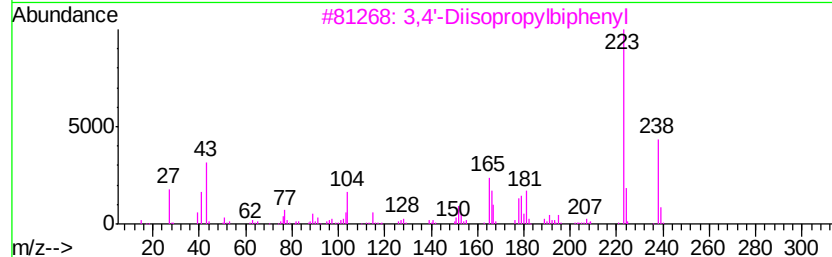
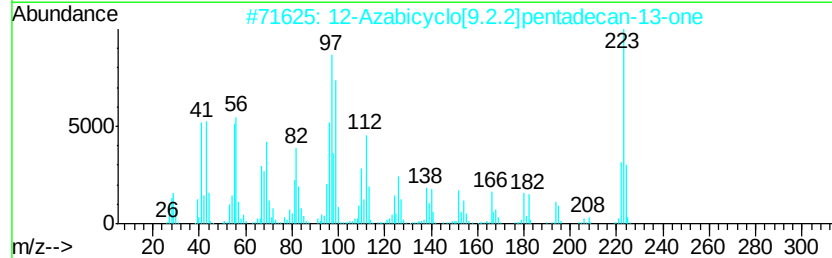
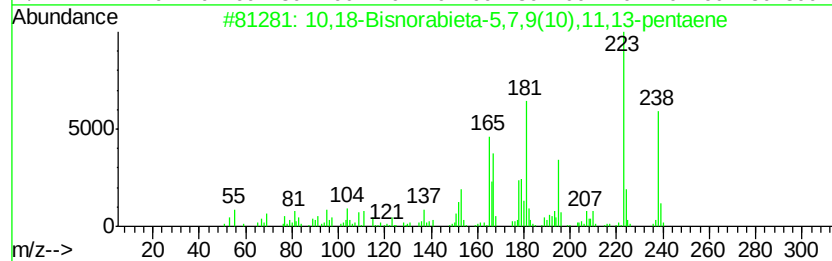
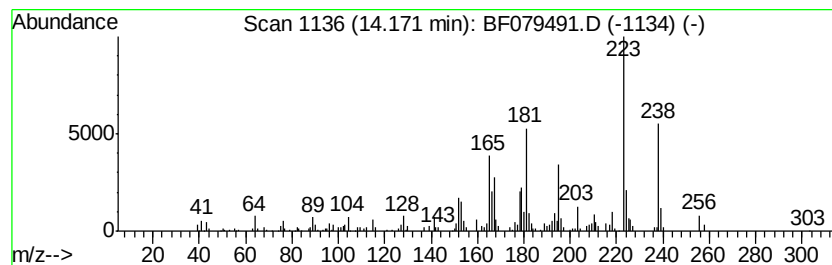
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.17	3.09 ng	191157	Phenanthrene-d10	12.59

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2			12-Azabicyclo[9.2.2]pentadecan-1...	223	C14H25NO	1000191-26-7	64
3			3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	62
4			3-(N,N-Diethylamino)carbazole	238	C16H18N2	054994-31-9	49
5			4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	45



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-30D

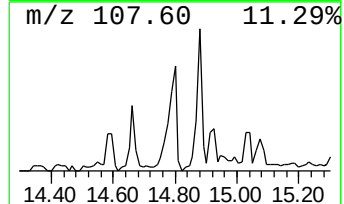
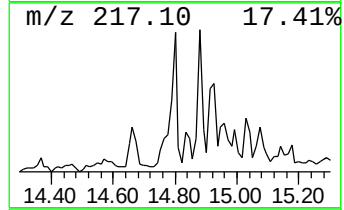
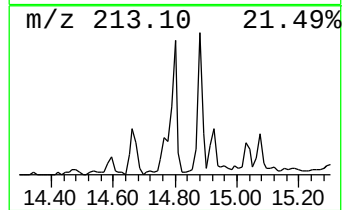
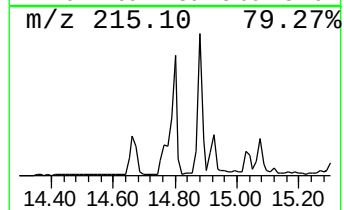
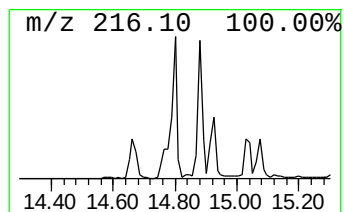
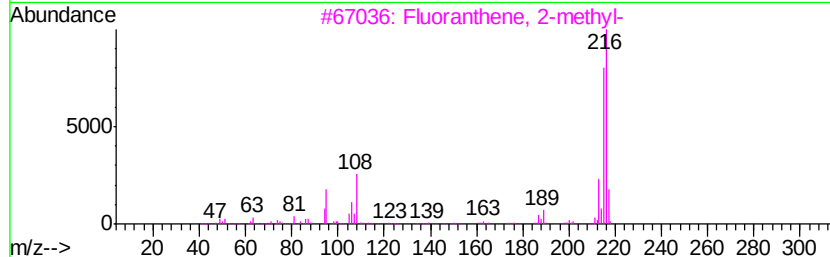
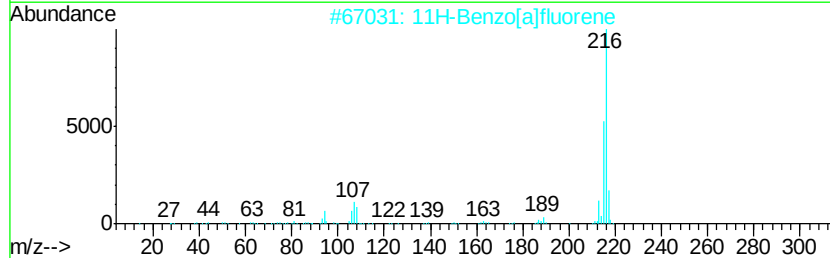
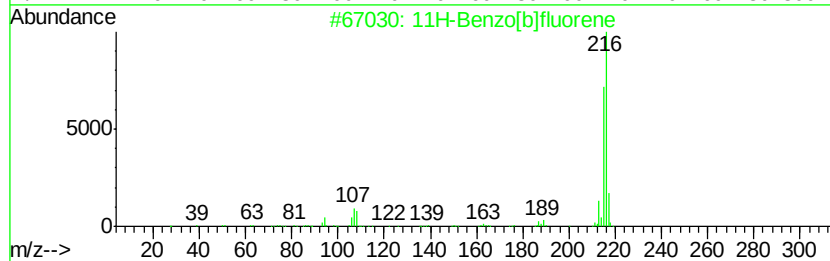
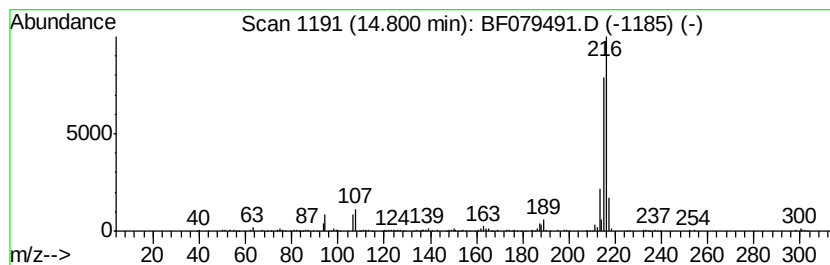
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 11H-Benzo[b]fluorene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	3.78 ng	403848	Chrysene-d12	15.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	97
2		11H-Benzo[a]fluorene	216	C17H12	000238-84-6	94
3		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	94
4		Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
5		1,3,5-Triazine-2,4,6-triamine, N...	216	C10H12N6	046731-79-7	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-30D

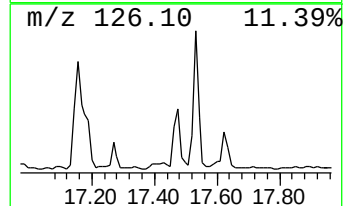
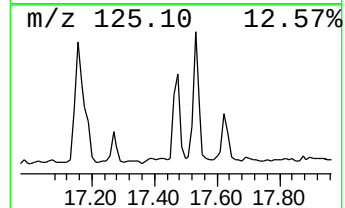
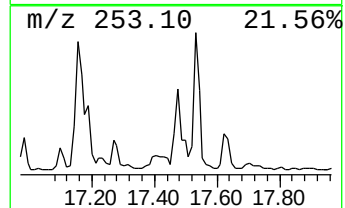
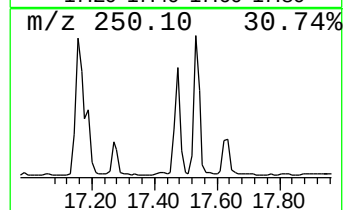
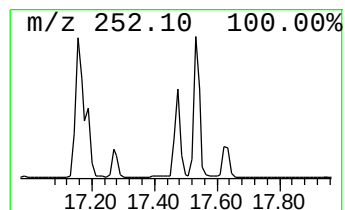
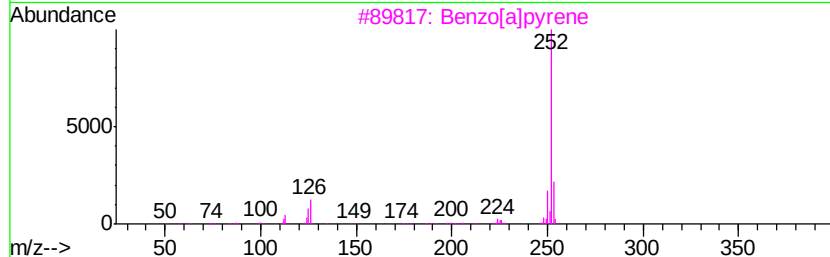
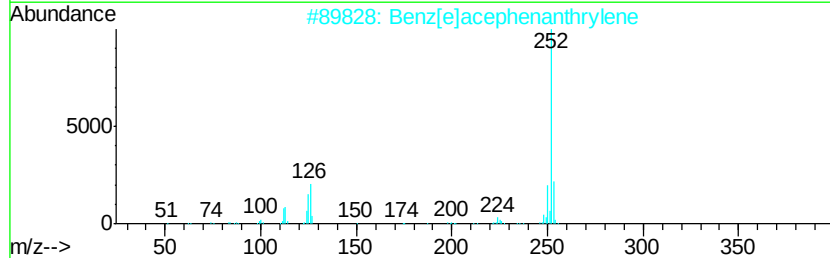
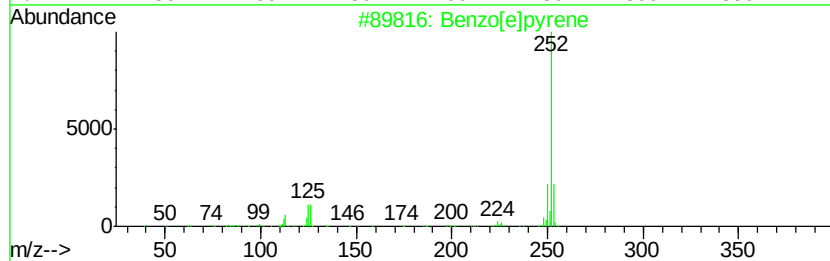
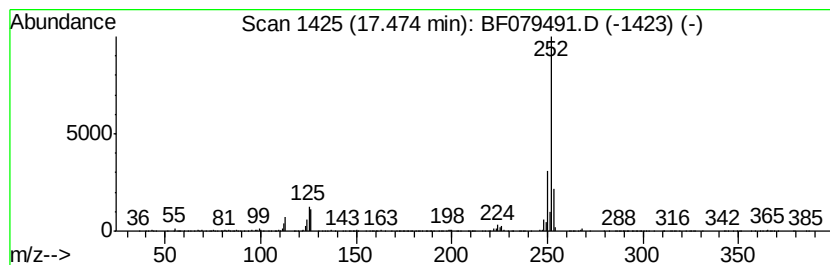
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.47	7.72 ng	438644	Perylene-d12	17.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]bpvrene	252	C20H12	000192-97-2	98
2		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	95
3		Benzo[a]bpvrene	252	C20H12	000050-32-8	94
4		Perylene	252	C20H12	000198-55-0	90
5		Benzo[j]fluoranthene	252	C20H12	000205-82-3	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052615\
Data File : BF079491.D
Acq On : 26 May 2015 22:50
Operator : TP/IZ
Sample : G2289-10 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-30D

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052615.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-methoxy...	1.43	10.4	ng	296860	1	7.02	569668	20.0
unknown6.73	6.73	14.8	ng	422637	1	7.02	569668	20.0
Phenanthrene, 2-m...	13.22	3.7	ng	226763	4	12.59	1236750	20.0
Phenanthrene, 1-m...	13.26	3.8	ng	233237	4	12.59	1236750	20.0
Anthracene, 2-met...	13.31	2.2	ng	134480	4	12.59	1236750	20.0
4H-Cyclopenta[def...	13.36	7.7	ng	475651	4	12.59	1236750	20.0
2-Phenyl naphthalene	13.60	2.8	ng	171275	4	12.59	1236750	20.0
9,10-Dimethylanth...	13.91	2.6	ng	164020	4	12.59	1236750	20.0
10,18-Bisnorabiet...	14.17	3.1	ng	191157	4	12.59	1236750	20.0
11H-Benzo[b]fluorene	14.80	3.8	ng	403848	5	15.87	2134330	20.0
Benzo[e]pyrene	17.47	7.7	ng	438644	6	17.60	1135780	20.0