

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040815\
 Data File : BF078353.D
 Acq On : 9 Apr 2015 00:22
 Operator : TP/IZ
 Sample : G1793-03 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3DA

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-BF040115.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	1.081	15	17	19	rVB	213520	239167	1.96%	0.211%
2	1.653	65	67	95	rVB	564453	1052546	8.61%	0.929%
3	4.944	352	355	358	rBV	633819	766056	6.27%	0.676%
4	5.093	365	368	373	rBV	263958	444889	3.64%	0.393%
5	5.436	395	398	401	rVB	178752	211237	1.73%	0.186%
6	6.327	473	476	477	rBV	273895	323738	2.65%	0.286%
7	6.362	477	479	482	rVB	257491	247188	2.02%	0.218%
8	6.522	491	493	497	rBV	191686	216588	1.77%	0.191%
9	6.693	507	508	514	rVB3	219786	403416	3.30%	0.356%
10	6.910	525	527	530	rVB	232346	236305	1.93%	0.209%
11	7.127	544	546	549	rVB	1479457	1783858	14.59%	1.574%
12	7.230	550	555	557	rBV2	252286	310627	2.54%	0.274%
13	7.322	561	563	566	rVB	166808	200404	1.64%	0.177%
14	7.619	587	589	591	rVB	202466	230285	1.88%	0.203%
15	8.099	629	631	633	rVB	314007	278071	2.28%	0.245%
16	8.190	633	639	641	rBV2	305894	608249	4.98%	0.537%
17	8.236	641	643	647	rVV	365507	788357	6.45%	0.696%
18	8.545	663	670	672	rVV	6771028	12222586	100.00%	10.785%
19	8.956	703	706	708	rVV	115382	220111	1.80%	0.194%
20	9.082	712	717	718	rVV2	150510	329827	2.70%	0.291%
21	9.128	718	721	724	rVV3	137261	328068	2.68%	0.289%
22	9.288	733	735	737	rVV2	124942	305803	2.50%	0.270%
23	9.379	740	743	746	rVB	2898749	4059757	33.22%	3.582%
24	9.505	751	754	756	rBV	2447849	2765056	22.62%	2.440%
25	9.962	792	794	796	rBV	1878641	1740759	14.24%	1.536%
26	10.076	801	804	808	rVB	1684883	1953797	15.99%	1.724%
27	10.156	808	811	813	rBV	409400	700236	5.73%	0.618%
28	10.248	816	819	820	rBV	901704	995952	8.15%	0.879%
29	10.282	820	822	823	rVV	359191	461791	3.78%	0.407%
30	10.385	829	831	835	rVB	426126	589365	4.82%	0.520%
31	10.488	837	840	842	rBV	915113	1288897	10.55%	1.137%
32	10.659	850	855	857	rBV	835704	1037760	8.49%	0.916%
33	10.705	857	859	861	rVV	4280715	4818519	39.42%	4.252%
34	10.796	864	867	870	rVV	252095	360078	2.95%	0.318%

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35	10.911	875	877	880	rBV2	343744	496332	4.06%	0.438%
36	11.048	886	889	891	rBV3	186348	316949	2.59%	0.280%
37	11.174	897	900	902	rBV2	133468	284283	2.33%	0.251%
38	11.219	902	904	906	rVB	439670	471444	3.86%	0.416%
39	11.299	906	911	912	rBV	368468	656880	5.37%	0.580%
40	11.334	912	914	917	rVV	2187713	2679776	21.92%	2.365%
41	11.402	917	920	922	rVV	281242	519112	4.25%	0.458%
42	11.459	922	925	927	rVV3	398252	828694	6.78%	0.731%
43	11.551	930	933	935	rVB2	311327	565137	4.62%	0.499%
44	11.608	935	938	939	rBV	702162	721596	5.90%	0.637%
45	11.631	939	940	943	rVB3	195351	235788	1.93%	0.208%
46	12.008	971	973	975	rVV	378852	686189	5.61%	0.605%
47	12.042	975	976	978	rVV	336902	382178	3.13%	0.337%
48	12.111	981	982	984	rVV	309902	333402	2.73%	0.294%
49	12.191	987	989	993	rVV4	130838	324564	2.66%	0.286%
50	12.259	993	995	998	rVB	192764	229628	1.88%	0.203%
51	12.351	1002	1003	1006	rVV	393460	503842	4.12%	0.445%
52	12.534	1013	1019	1021	rBV	5268780	8910091	72.90%	7.862%
53	12.591	1021	1024	1025	rVB	2165185	2500248	20.46%	2.206%
54	12.637	1025	1028	1030	rBV	165936	231979	1.90%	0.205%
55	12.877	1046	1049	1052	rVV2	272320	532962	4.36%	0.470%
56	13.117	1067	1070	1071	rBV	1425035	1368252	11.19%	1.207%
57	13.151	1071	1073	1075	rVB	1479469	1485455	12.15%	1.311%
58	13.208	1075	1078	1079	rBV	606868	613990	5.02%	0.542%
59	13.242	1079	1081	1086	rVV2	1750698	3119516	25.52%	2.753%
60	13.494	1100	1103	1105	rVV	823121	1077059	8.81%	0.950%
61	13.665	1116	1118	1120	rVV	168990	216425	1.77%	0.191%
62	13.711	1120	1122	1126	rVB	163418	367834	3.01%	0.325%
63	13.791	1127	1129	1131	rBV	547750	801804	6.56%	0.707%
64	13.837	1131	1133	1134	rVV	401260	510312	4.18%	0.450%
65	13.871	1134	1136	1137	rVV	305607	348859	2.85%	0.308%
66	13.905	1137	1139	1143	rVV2	214191	357401	2.92%	0.315%
67	13.997	1143	1147	1148	rVV	3347189	4568845	37.38%	4.031%
68	14.100	1154	1156	1159	rVB	1335485	1426985	11.67%	1.259%
69	14.271	1167	1171	1174	rBV	3848826	6463478	52.88%	5.703%
70	14.477	1187	1189	1193	rVB	235192	275805	2.26%	0.243%
71	14.545	1193	1195	1201	rVB2	340987	531875	4.35%	0.469%

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72	14.682	1201	1207	1209	rVB	744562	1668561	13.65%	1.472%
73	14.762	1212	1214	1216	rBV	759364	815897	6.68%	0.720%
74	14.797	1216	1217	1219	rVB	423955	531025	4.34%	0.469%
75	14.888	1223	1225	1226	rBV	669393	846506	6.93%	0.747%
76	14.945	1229	1230	1235	rVB	381376	677090	5.54%	0.597%
77	15.208	1249	1253	1256	rBV	141619	266150	2.18%	0.235%
78	15.277	1256	1259	1263	rBV2	118154	332484	2.72%	0.293%
79	15.437	1269	1273	1275	rBV2	166347	352483	2.88%	0.311%
80	15.494	1275	1278	1280	rVV2	406886	845238	6.92%	0.746%
81	15.745	1296	1300	1302	rBV2	2025069	3547556	29.02%	3.130%
82	15.791	1302	1304	1306	rVV	1313985	1600331	13.09%	1.412%
83	15.871	1309	1311	1313	rBV2	250890	367586	3.01%	0.324%
84	16.248	1341	1344	1346	rVV	430589	618753	5.06%	0.546%
85	16.363	1349	1354	1355	rVV3	268281	489457	4.00%	0.432%
86	16.397	1355	1357	1358	rVV2	175233	317399	2.60%	0.280%
87	16.420	1358	1359	1365	rVV2	379049	759565	6.21%	0.670%
88	17.026	1409	1412	1418	rBV	912685	2237421	18.31%	1.974%
89	17.140	1419	1422	1424	rVV	396121	514053	4.21%	0.454%
90	17.346	1436	1440	1442	rVB	740199	1103371	9.03%	0.974%
91	17.403	1442	1445	1448	rBV	1450898	1963291	16.06%	1.732%
92	17.460	1448	1450	1452	rBV	384605	575723	4.71%	0.508%
93	17.666	1462	1468	1473	rBV3	118823	523628	4.28%	0.462%
94	17.803	1478	1480	1483	rVV2	110202	257360	2.11%	0.227%
95	17.963	1491	1494	1499	rVV	205571	424012	3.47%	0.374%
96	18.591	1546	1549	1551	rBV2	125318	258209	2.11%	0.228%
97	18.729	1558	1561	1566	rVB2	455411	890215	7.28%	0.785%
98	19.094	1589	1593	1601	rBV	490434	1112260	9.10%	0.981%
99	19.289	1607	1610	1614	rVB	281731	530367	4.34%	0.468%
100	20.957	1750	1756	1762	rBV3	101083	442352	3.62%	0.390%

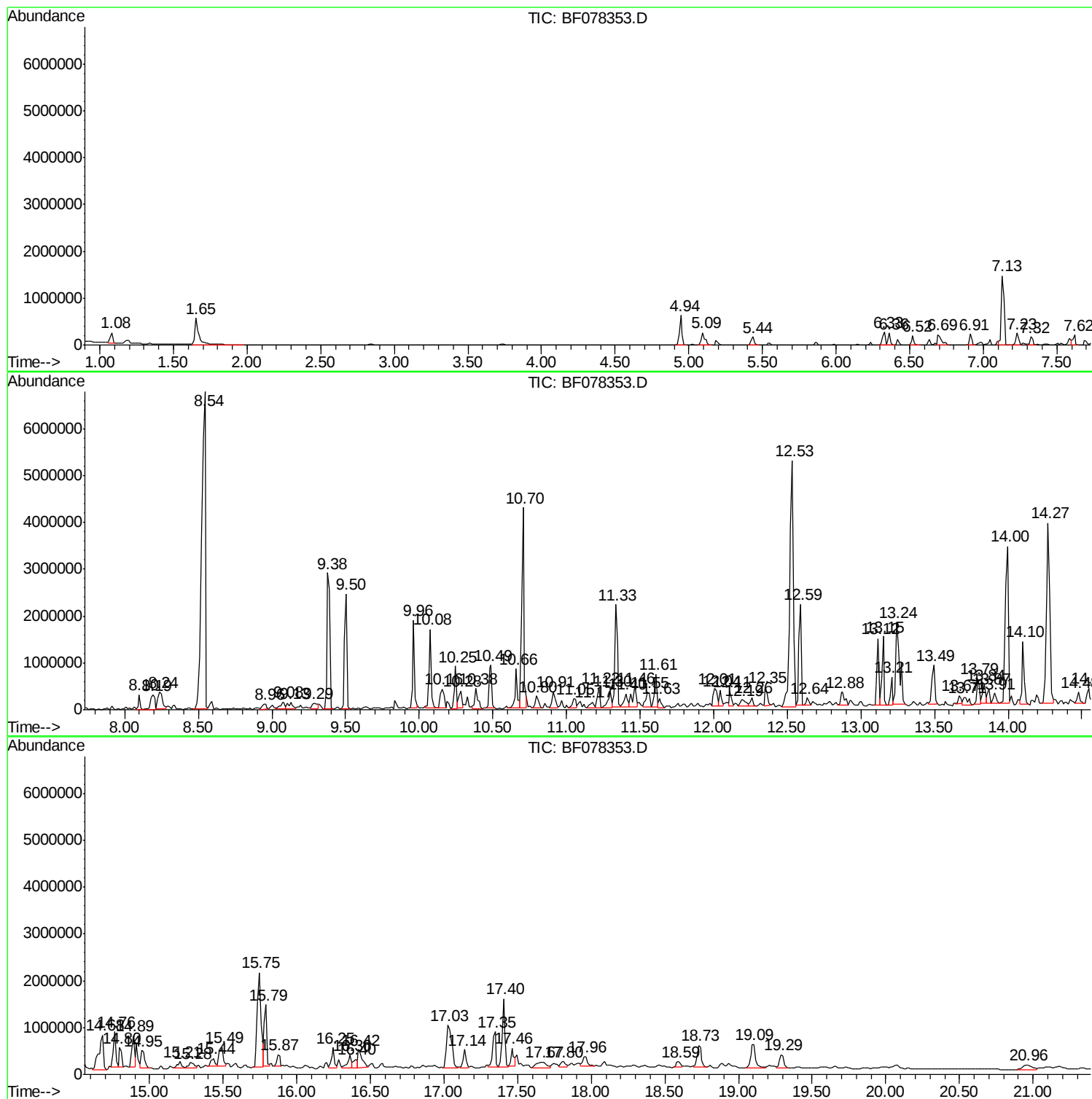
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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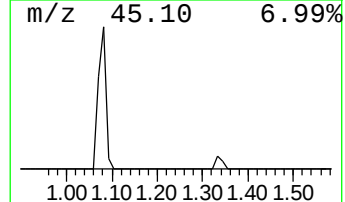
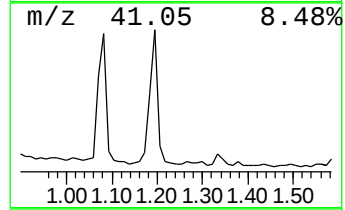
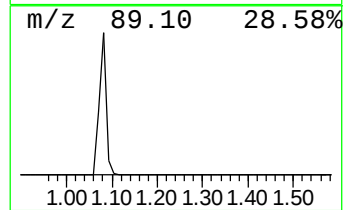
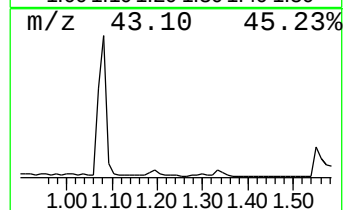
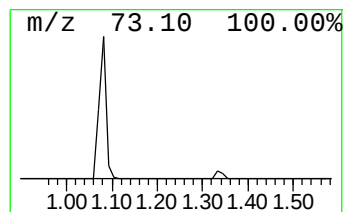
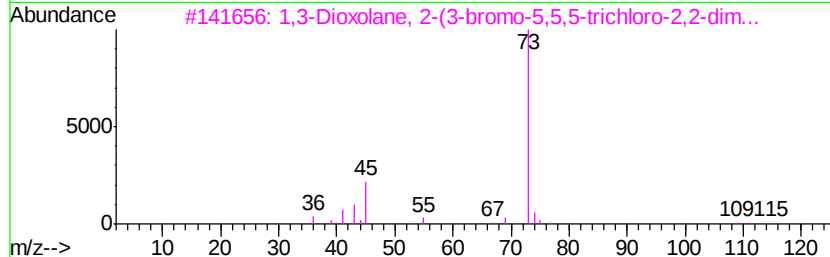
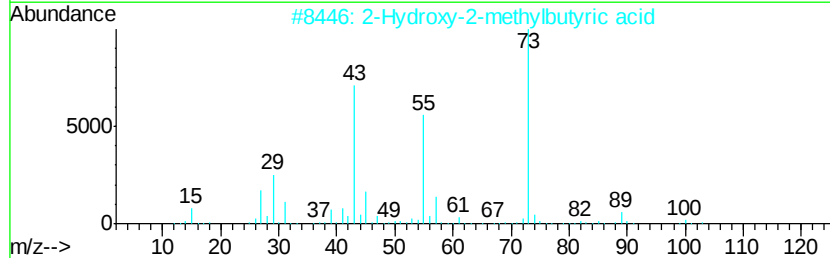
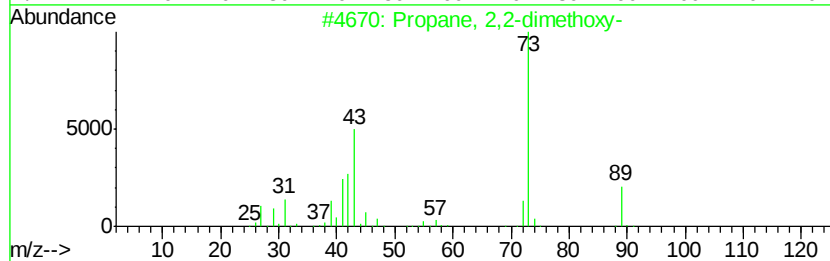
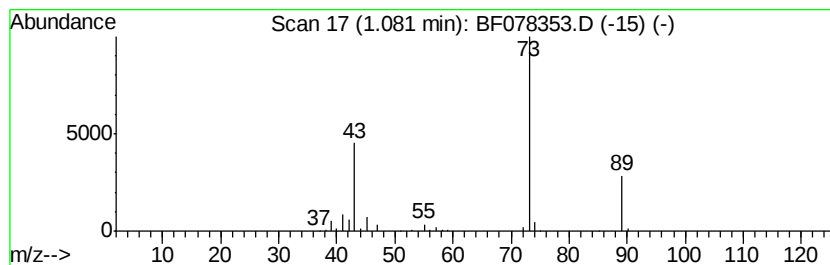
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown1.08 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.08	20.24 ng	239167	1,4-Dichlorobenzene-d4	6.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	38
2		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
3		1,3-Dioxolane, 2-(3-bromo-5,5,5-...	352	C10H16BrCl3O2	1000115-31-4	9
4		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
5		N-Ethylformamide	73	C3H7NO	000627-45-2	5



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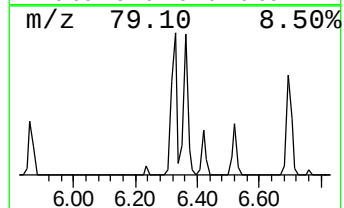
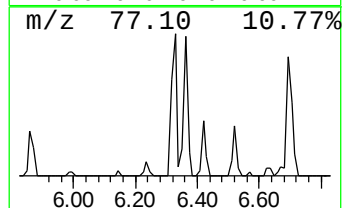
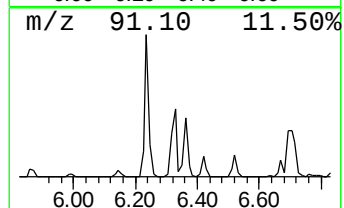
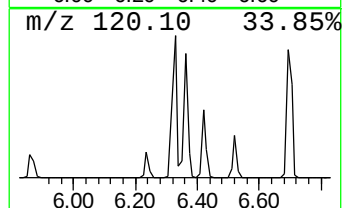
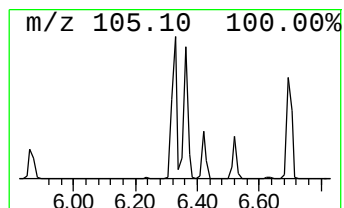
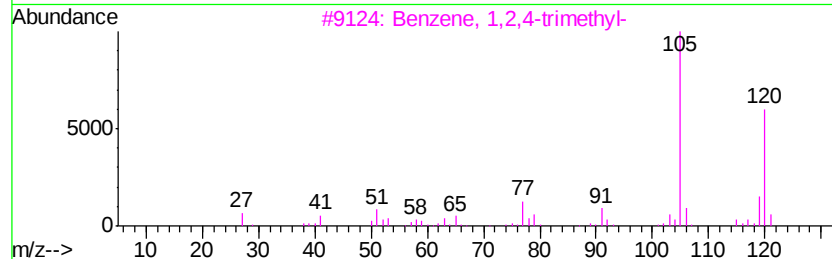
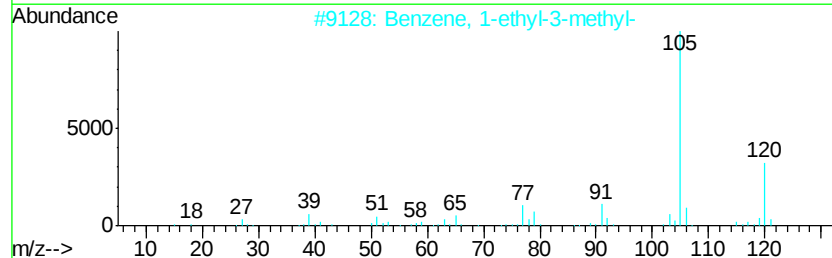
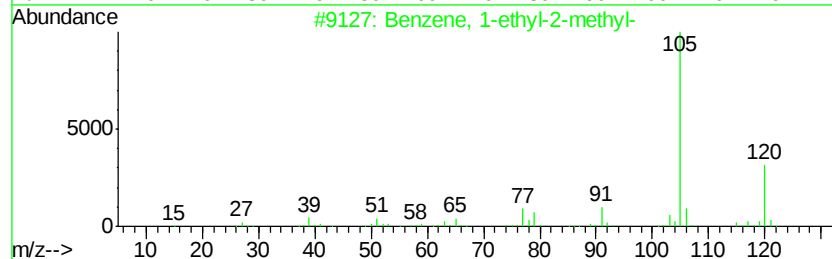
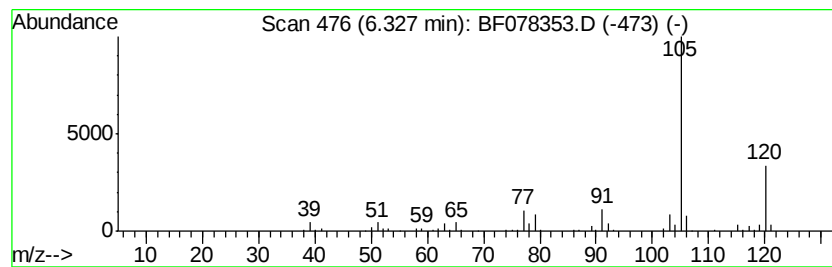
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.33	27.40 ng	323738	1,4-Dichlorobenzene-d4	6.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91



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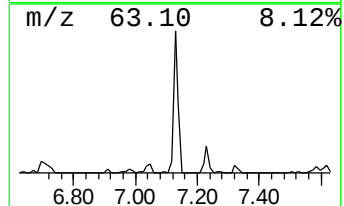
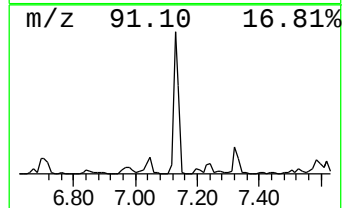
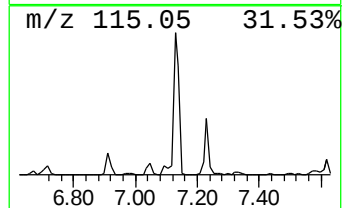
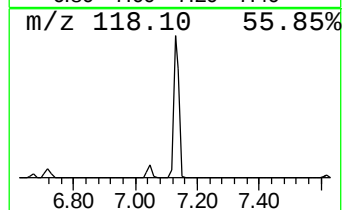
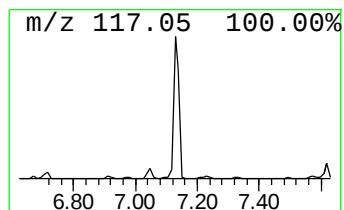
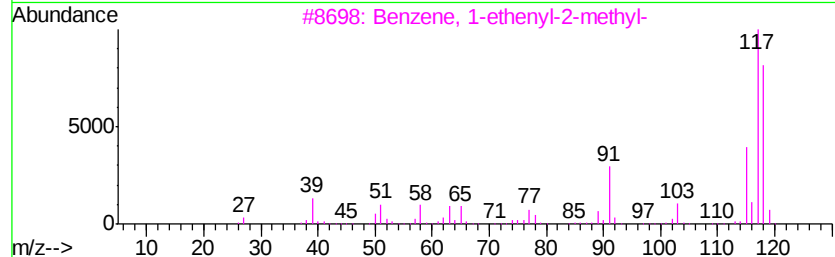
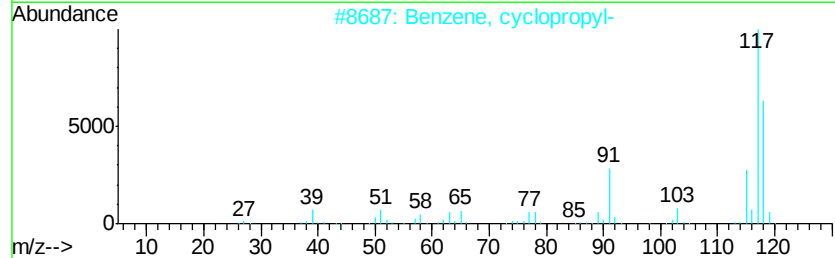
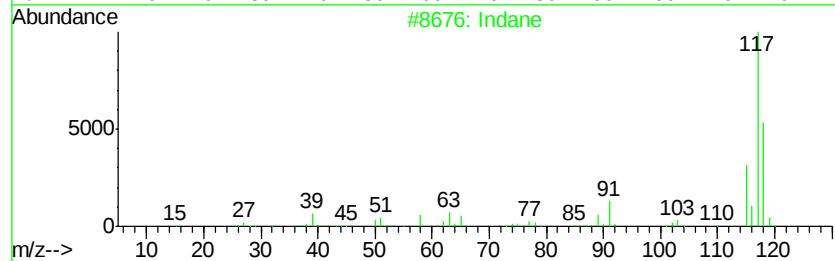
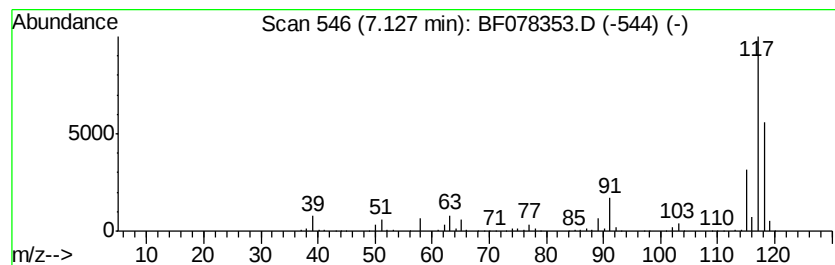
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.13	150.98 ng	1783860	1,4-Dichlorobenzene-d4	6.91

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane		118	C9H10	000496-11-7	91
2	Benzene, cyclopropyl-		118	C9H10	000873-49-4	83
3	Benzene, 1-ethenyl-2-methyl-		118	C9H10	000611-15-4	60
4	Benzene, 1-propenyl-		118	C9H10	000637-50-3	59
5	Benzene, 1,1'-(1,5-hexadiene-1,6...		234	C18H18	004439-45-6	59



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040815\
Data File : BF078353.D
Acq On : 9 Apr 2015 00:22
Operator : TP/IZ
Sample : G1793-03 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

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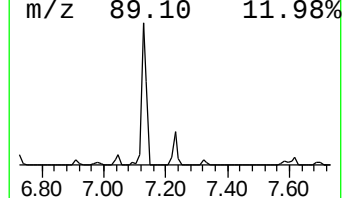
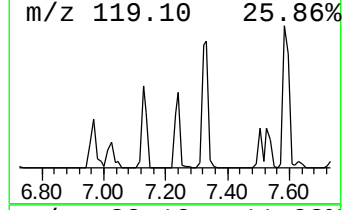
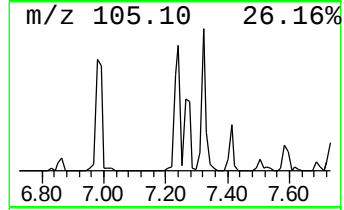
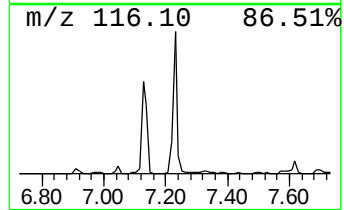
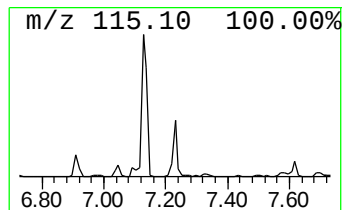
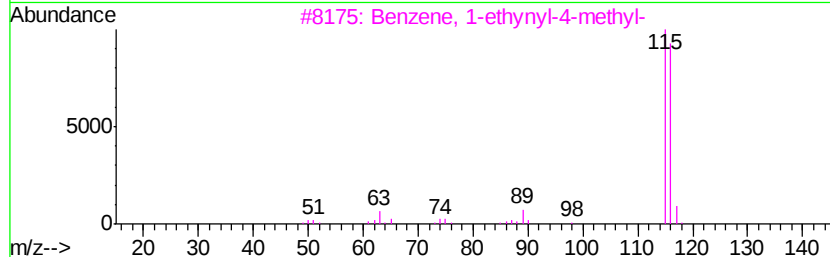
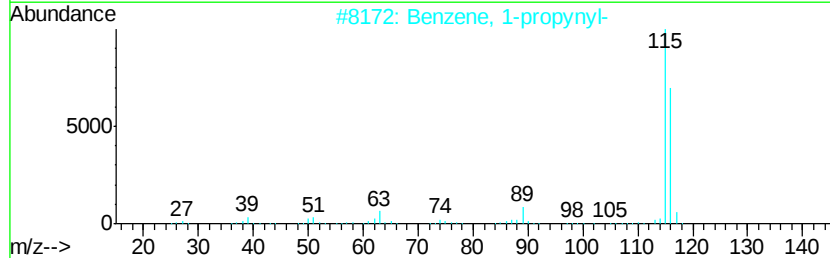
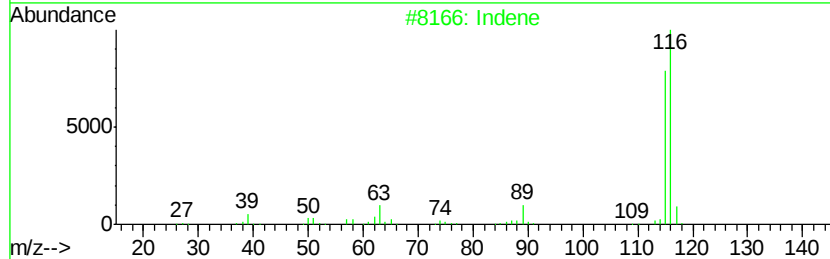
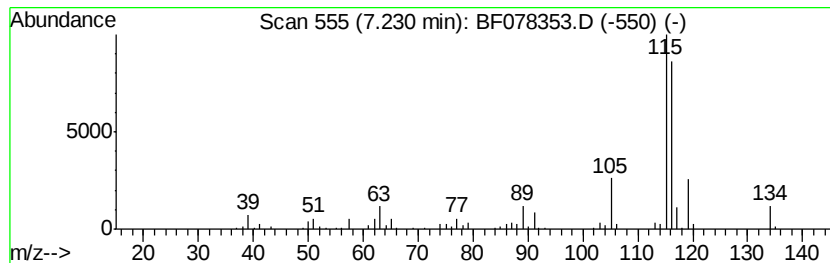
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzene, 1-propynyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.23	26.29 ng	310627	1,4-Dichlorobenzene-d4	6.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Indene	116	C9H8	000095-13-6	95
2		Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3		Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	76
4		Benzene, 1-ethynyl-2-methyl-	116	C9H8	000766-47-2	74
5		Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	58



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040815\
Data File : BF078353.D
Acq On : 9 Apr 2015 00:22
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Misc :
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ClientSampleId :
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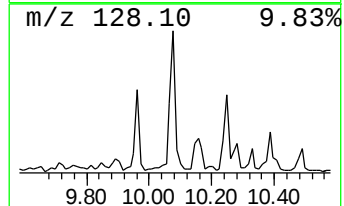
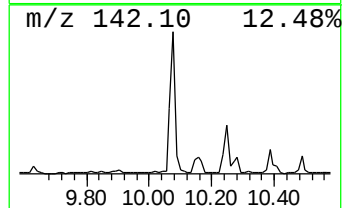
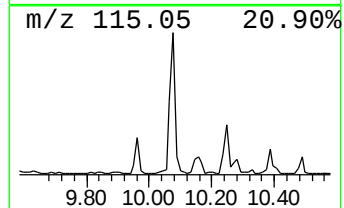
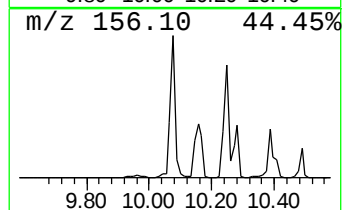
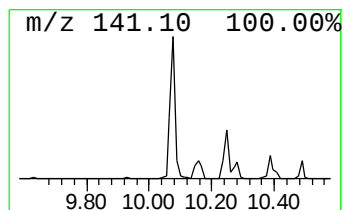
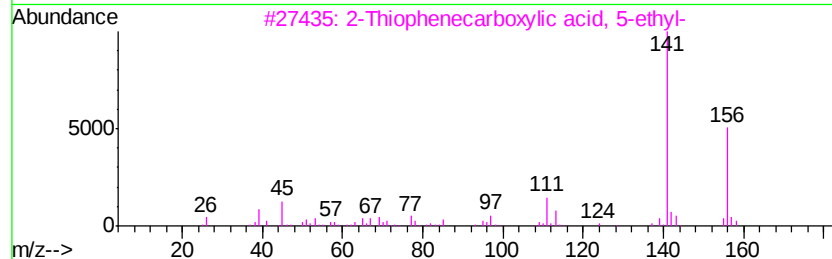
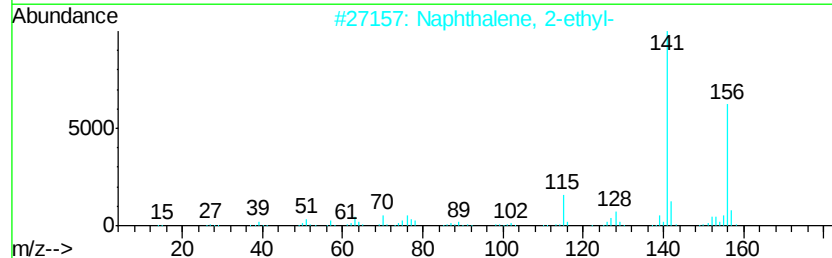
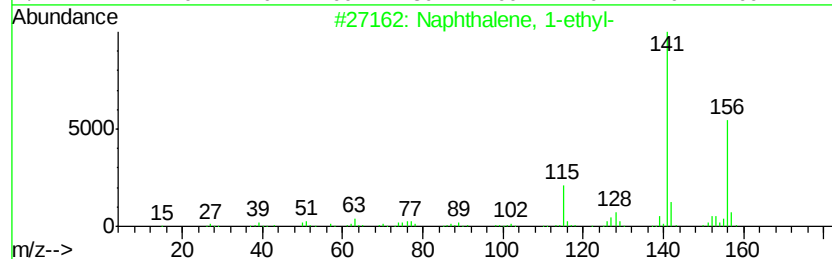
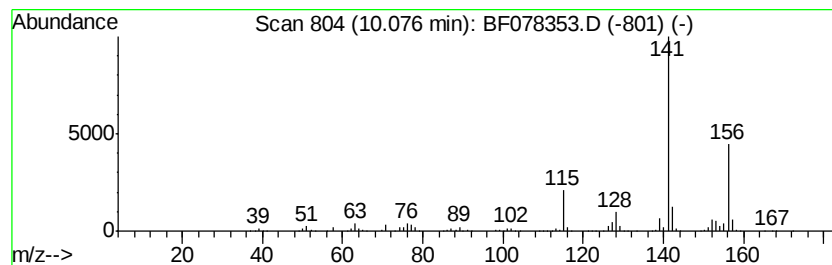
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, 1-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.08	37.65 ng	1953800	Acenaphthene-d10	10.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	96
2		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95
3		2-Thiophenecarboxylic acid, 5-et...	156	C7H8O2S	023229-72-3	64
4		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	43
5		Methyl phenylphosphinic acid	156	C7H9O2P	004271-13-0	42



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040815\
Data File : BF078353.D
Acq On : 9 Apr 2015 00:22
Operator : TP/IZ
Sample : G1793-03 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

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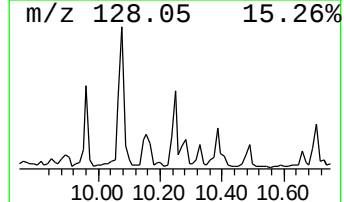
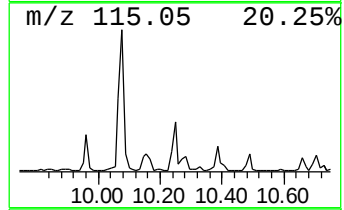
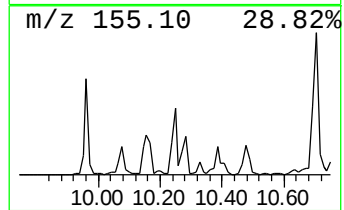
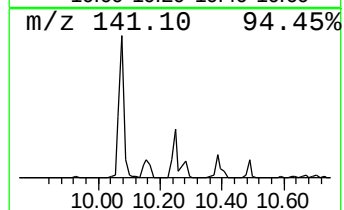
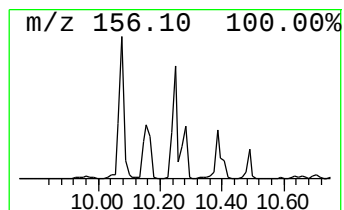
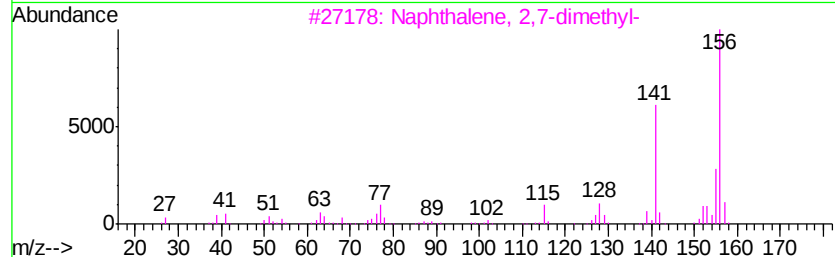
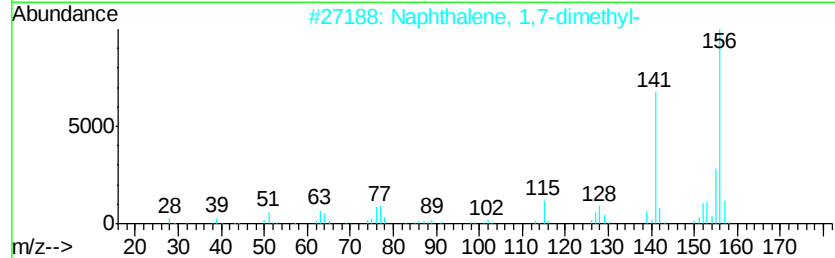
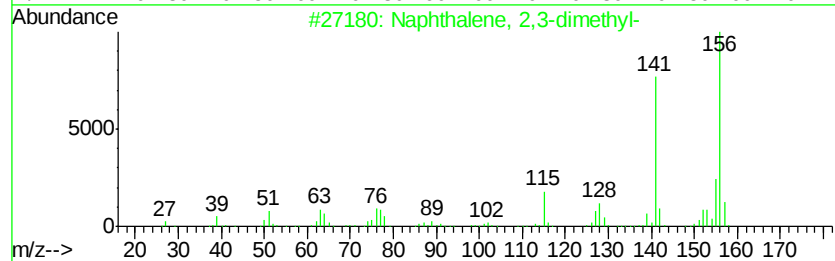
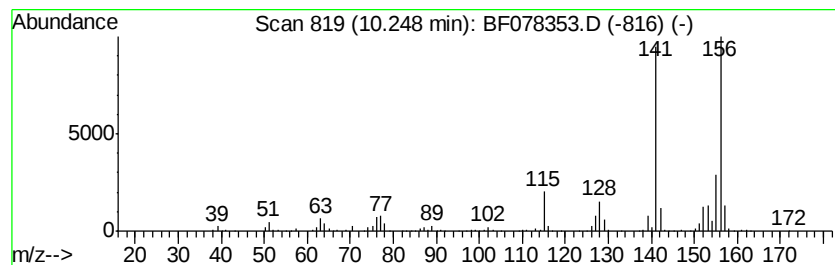
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Naphthalene, 2,3-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.25	19.19 ng	995952	Acenaphthene-d10	10.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040815\
Data File : BF078353.D
Acq On : 9 Apr 2015 00:22
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Misc :
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ClientSampleId :
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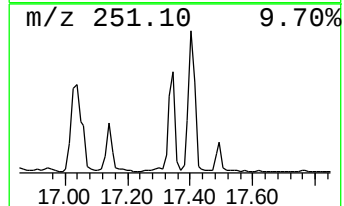
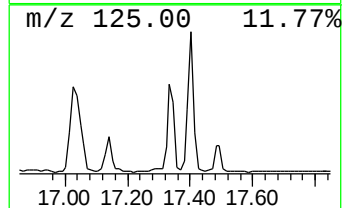
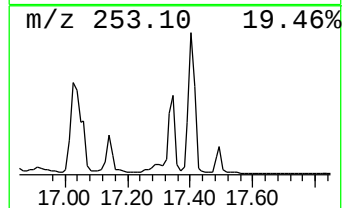
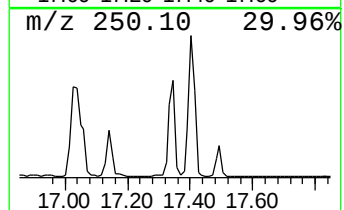
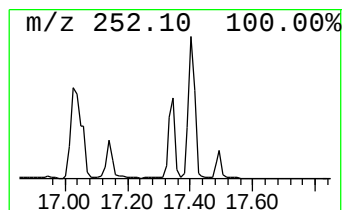
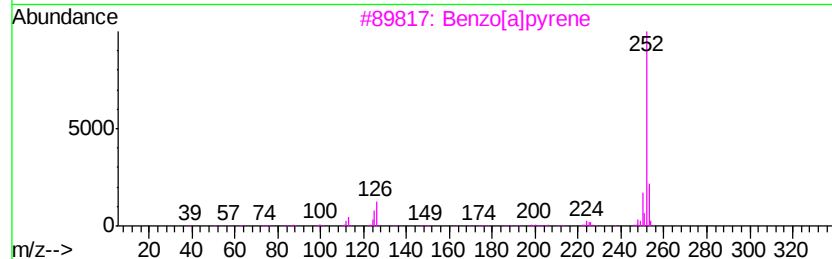
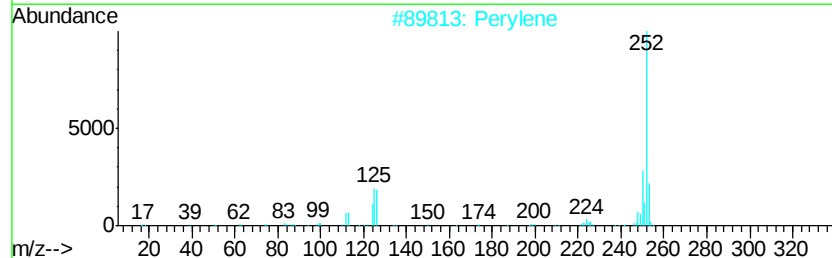
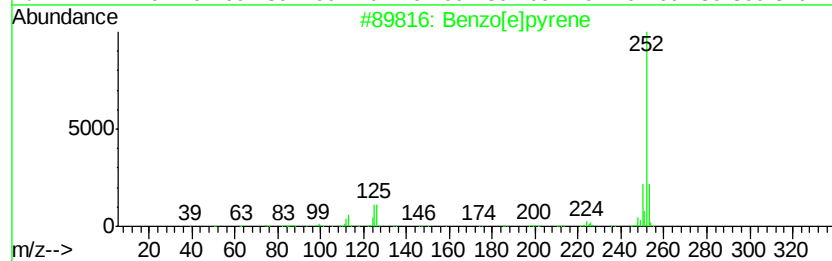
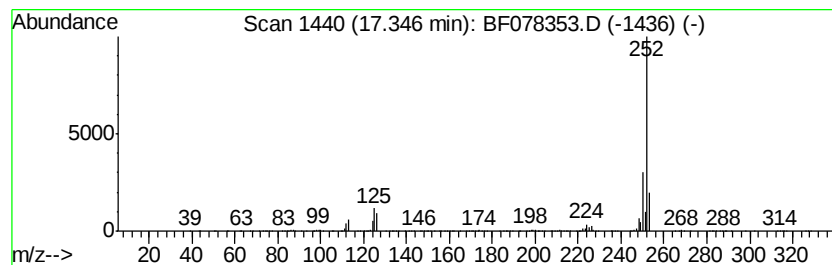
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.35	38.33 ng	1103370	Perylene-d12	17.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]bpvrene	252	C20H12	000192-97-2	96
2		Perylene	252	C20H12	000198-55-0	95
3		Benzo[a]bpvrene	252	C20H12	000050-32-8	93
4		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	90
5		Benzo[j]fluoranthene	252	C20H12	000205-82-3	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040815\
Data File : BF078353.D
Acq On : 9 Apr 2015 00:22
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ClientSampleId :
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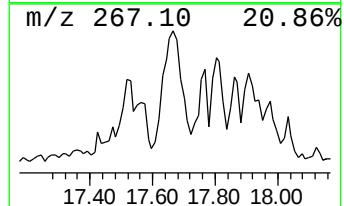
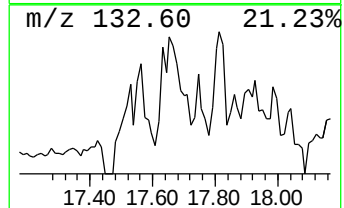
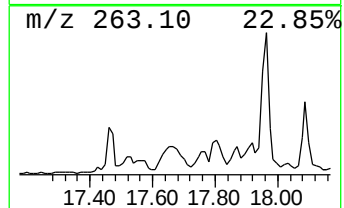
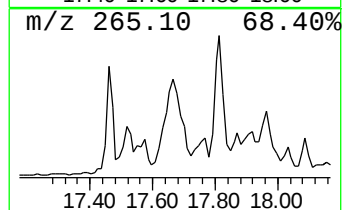
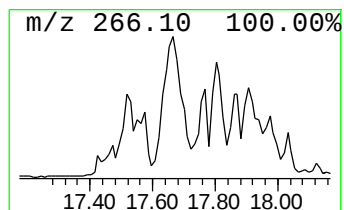
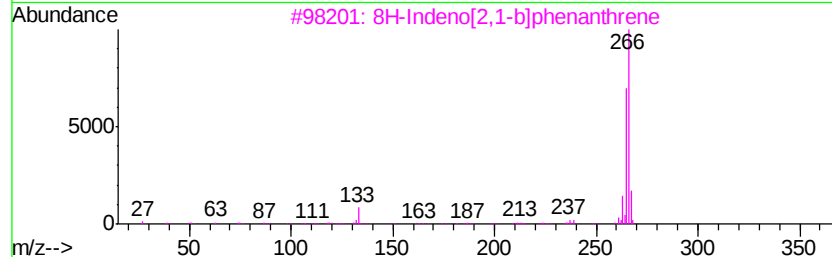
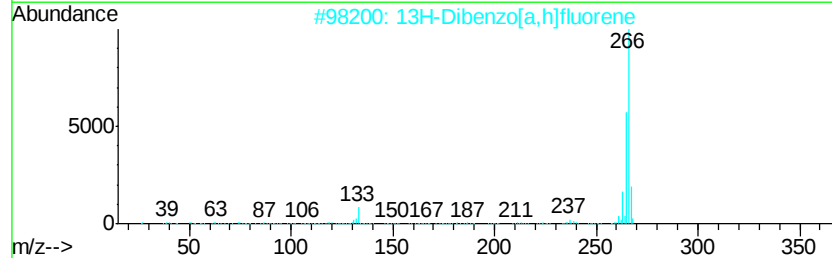
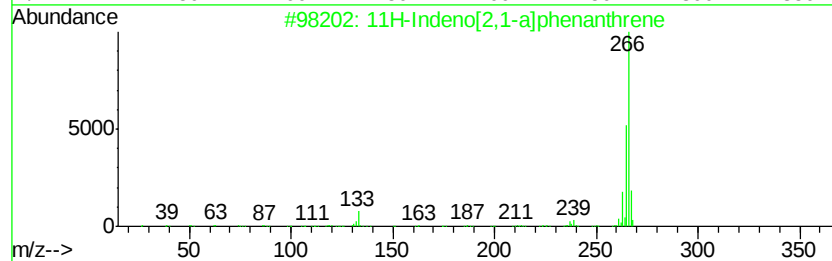
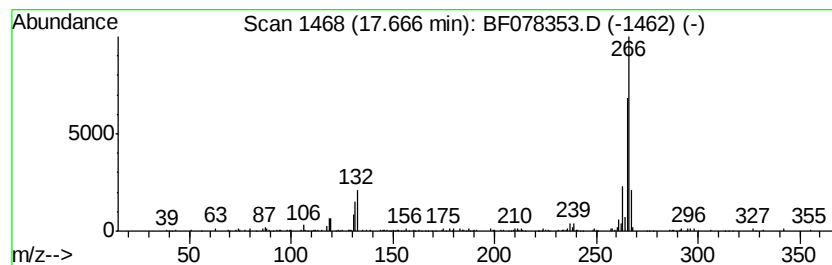
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 11H-Indeno[2,1-a]phenanthrene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.67	18.19 ng	523628	Perylene-d12	17.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Indeno[2,1-a]phenanthrene	266	C21H14	000220-97-3	89
2		13H-Dibenzof[a,h]fluorene	266	C21H14	000239-85-0	86
3		8H-Indeno[2,1-b]phenanthrene	266	C21H14	000241-28-1	76
4		4H-1-Benzopyran-4-one, 3-hydroxy...	266	C17H14O3	078396-37-9	64
5		Thiazole, 4-phenyl-2-(4-tolylami...	266	C16H14N2S	016098-04-7	52



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040815\
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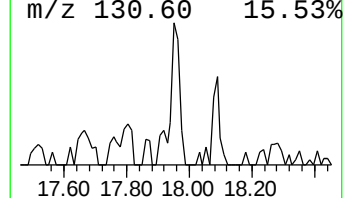
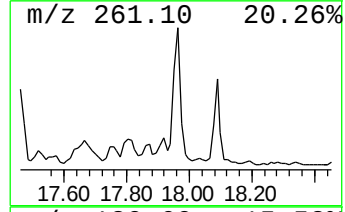
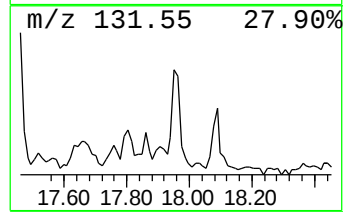
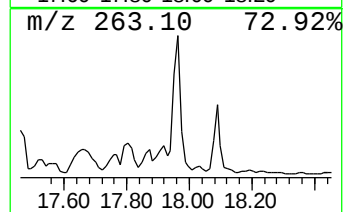
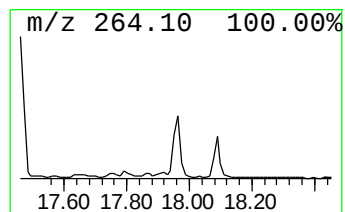
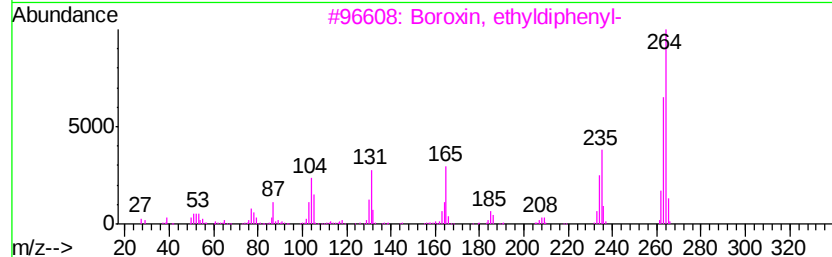
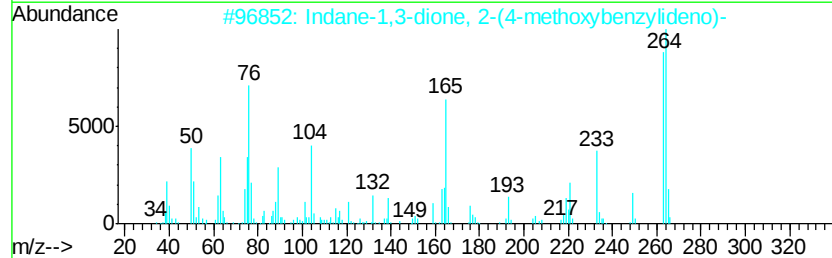
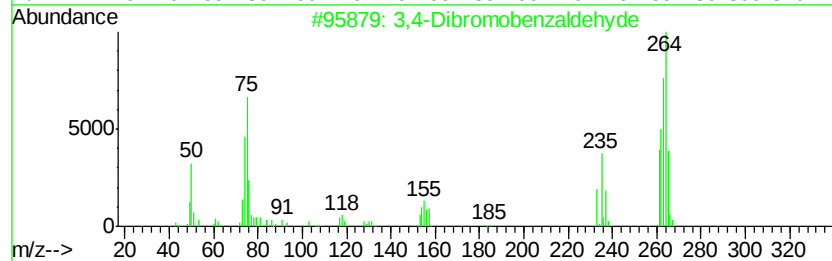
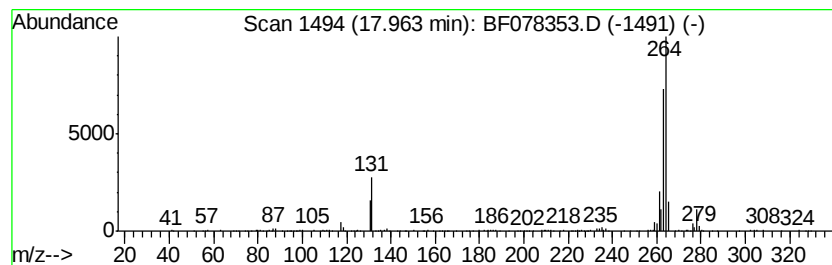
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 3,4-Dibromobenzaldehyde Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.96	14.73 ng	424012	Perylene-d12	17.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,4-Dibromobenzaldehyde	262	C7H4Br2O	074003-55-7	58
2		Indane-1,3-dione, 2-(4-methoxybe...	264	C17H12O3	007421-76-3	47
3		Boroxin, ethyldiphenyl-	264	C14H15B3O3	1000151-82-0	45
4		9,9-Dichloro-9-sila-9,10-dihydro...	264	C13H10Cl2Si	076470-26-3	38
5		7,9-Diamino-5,6,8,10-tetraaza-be...	264	C13H8N6O	052559-29-2	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040815\
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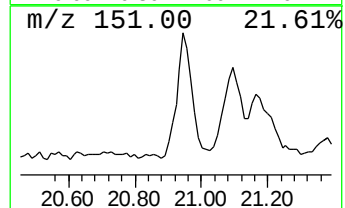
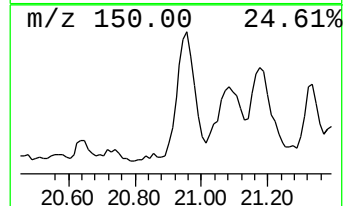
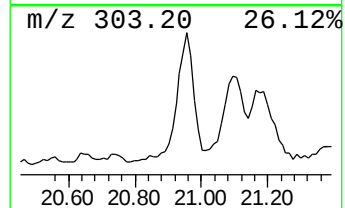
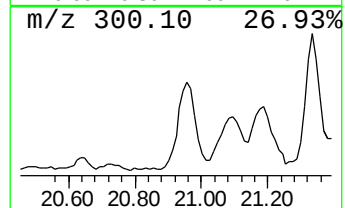
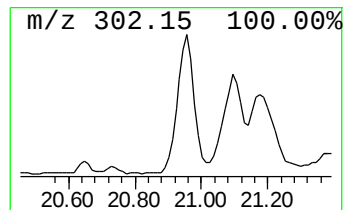
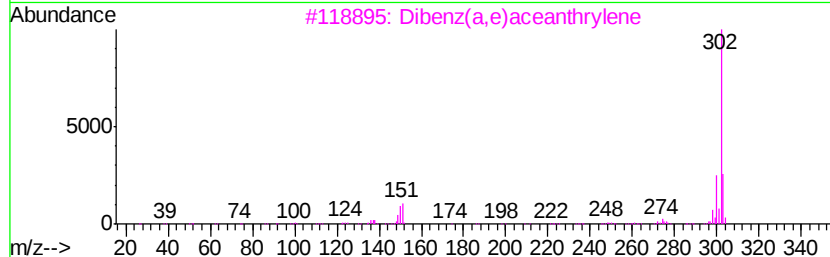
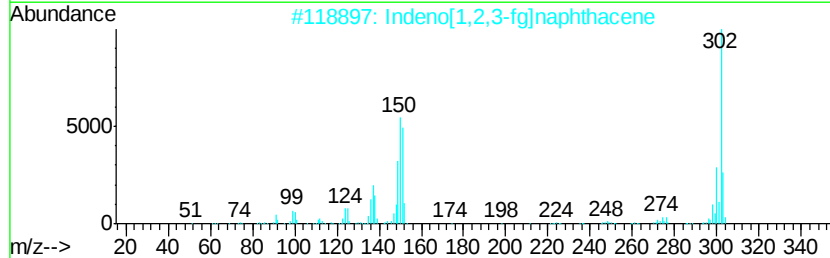
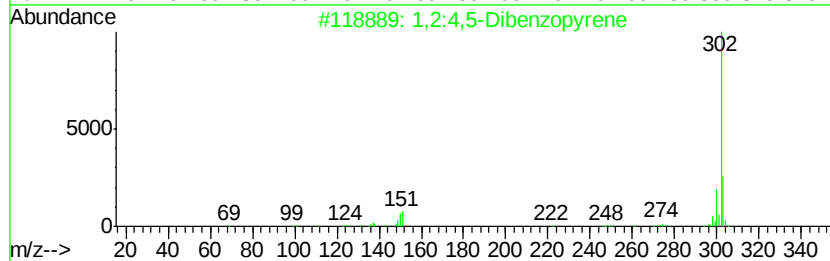
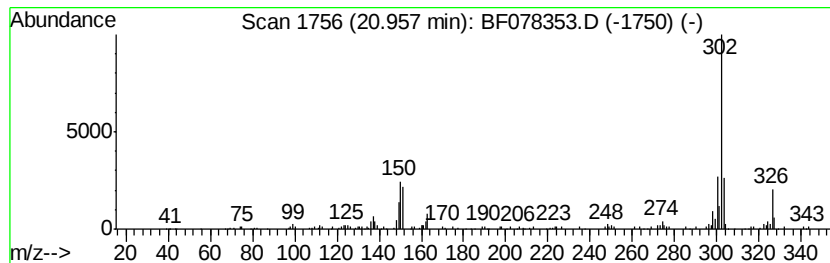
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 1,2:4,5-Dibenzopyrene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.96	15.37 ng	442352	Perylene-d12	17.46

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2:4,5-Dibenzopyrene	302	C24H14	000192-65-4	97
2		Indeno[1,2,3-fg]naphthacene	302	C24H14	000203-11-2	93
3		Dibenz(a,e)aceanthrylene	302	C24H14	005385-75-1	92
4		3,4:8,9-Dibenzopyrene	302	C24H14	000189-64-0	83
5		3,4:9,10-Dibenzopyrene	302	C24H14	000189-55-9	83



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040815\
Data File : BF078353.D
Acq On : 9 Apr 2015 00:22
Operator : TP/IZ
Sample : G1793-03 5X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-3DA

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.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
unknown1.08	1.08	20.2	ng	239167	1	6.91	236305	20.0
Benzene, 1-ethyl-...	6.33	27.4	ng	323738	1	6.91	236305	20.0
Indane	7.13	151.0	ng	1783860	1	6.91	236305	20.0
Benzene, 1-propynyl-	7.23	26.3	ng	310627	1	6.91	236305	20.0
Naphthalene, 1-et...	10.08	37.6	ng	1953800	3	10.66	1037760	20.0
Naphthalene, 2,3-...	10.25	19.2	ng	995952	3	10.66	1037760	20.0
Benzo[e]pyrene	17.35	38.3	ng	1103370	6	17.46	575723	20.0
11H-Indeno[2,1-a]...	17.67	18.2	ng	523628	6	17.46	575723	20.0
3,4-Dibromobenzal...	17.96	14.7	ng	424012	6	17.46	575723	20.0
1,2:4,5-Dibenzopy...	20.96	15.4	ng	442352	6	17.46	575723	20.0