

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040815\
 Data File : BF078342.D
 Acq On : 8 Apr 2015 19:07
 Operator : TP/IZ
 Sample : G1793-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-6D

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	20	rVB	1202845	1533717	70.74%	4.966%
2	1.195	23	27	30	rVB	298020	498217	22.98%	1.613%
3	1.344	38	40	43	rVB	121012	133633	6.16%	0.433%
4	1.550	57	58	65	rBV	122491	196805	9.08%	0.637%
5	1.653	65	67	97	rVB	843815	1599362	73.76%	5.178%
6	4.601	323	325	333	rVB	105108	143521	6.62%	0.465%
7	5.184	373	376	386	rBV	597431	872062	40.22%	2.824%
8	6.510	489	492	494	rBV	905709	998255	46.04%	3.232%
9	6.624	499	502	505	rBV	1053603	986643	45.51%	3.195%
10	6.910	525	527	530	rBV	256080	235982	10.88%	0.764%
11	7.093	541	543	545	rBV	754927	730924	33.71%	2.367%
12	7.127	545	546	548	rVB	138217	101976	4.70%	0.330%
13	7.607	586	588	591	rVB	620876	577743	26.65%	1.871%
14	8.179	636	638	641	rBV	35318	66654	3.07%	0.216%
15	8.225	641	642	645	rVV	50014	89667	4.14%	0.290%
16	8.488	662	665	666	rBV	363007	388663	17.93%	1.258%
17	8.522	666	668	670	rVV	1617960	2168197	100.00%	7.020%
18	9.208	726	728	730	rVB	58318	48571	2.24%	0.157%
19	9.368	740	742	745	rBV	516306	669772	30.89%	2.169%
20	9.493	750	753	755	rBV	511482	483659	22.31%	1.566%
21	9.836	781	783	786	rBV	1194596	1041404	48.03%	3.372%
22	9.951	791	793	796	rBV	258688	286670	13.22%	0.928%
23	10.065	801	803	807	rVB	183320	187336	8.64%	0.607%
24	10.145	807	810	813	rBV	57188	87447	4.03%	0.283%
25	10.236	816	818	820	rBV	112412	128697	5.94%	0.417%
26	10.271	820	821	822	rVV	65268	51884	2.39%	0.168%
27	10.316	824	825	827	rVV	23920	43532	2.01%	0.141%
28	10.351	827	828	829	rVV	114776	86106	3.97%	0.279%
29	10.385	829	831	836	rVB	41219	66209	3.05%	0.214%
30	10.476	836	839	842	rBV	142724	148005	6.83%	0.479%
31	10.648	852	854	856	rBV	455130	453028	20.89%	1.467%
32	10.693	856	858	859	rVV	734985	785659	36.24%	2.544%
33	10.853	869	872	875	rVB2	36090	45370	2.09%	0.147%
34	10.899	875	876	880	rBV2	44089	64752	2.99%	0.210%

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35	11.208	900	903	905	rVB	93957	101550	4.68%	0.329%
36	11.288	909	910	912	rVV	38397	47046	2.17%	0.152%
37	11.322	912	913	916	rVV	354540	412540	19.03%	1.336%
38	11.402	918	920	921	rVV	29719	42358	1.95%	0.137%
39	11.436	921	923	924	rVV	28486	48178	2.22%	0.156%
40	11.459	924	925	929	rVV2	48011	52522	2.42%	0.170%
41	11.551	931	933	935	rVV	261230	261384	12.06%	0.846%
42	11.596	935	937	938	rVV	96264	95496	4.40%	0.309%
43	11.631	938	940	942	rVB	491649	630297	29.07%	2.041%
44	12.008	969	973	974	rBV	60075	84958	3.92%	0.275%
45	12.042	974	976	977	rVB	42371	51452	2.37%	0.167%
46	12.111	980	982	984	rVB	37415	44762	2.06%	0.145%
47	12.351	1001	1003	1006	rBV	52703	65711	3.03%	0.213%
48	12.477	1012	1014	1015	rBV	385805	390410	18.01%	1.264%
49	12.511	1015	1017	1019	rVV	1643082	1699706	78.39%	5.503%
50	12.568	1019	1022	1025	rVV	538698	524475	24.19%	1.698%
51	12.865	1046	1048	1051	rVV2	66979	70037	3.23%	0.227%
52	13.105	1067	1069	1070	rBV	173060	163537	7.54%	0.530%
53	13.139	1070	1072	1074	rVB	166511	180171	8.31%	0.583%
54	13.197	1074	1077	1078	rBV	78658	75736	3.49%	0.245%
55	13.231	1078	1080	1082	rVV	319853	366119	16.89%	1.185%
56	13.265	1082	1083	1085	rVB	101370	79509	3.67%	0.257%
57	13.482	1099	1102	1104	rBV	147686	149044	6.87%	0.483%
58	13.791	1126	1129	1130	rBV	46214	67172	3.10%	0.217%
59	13.825	1130	1132	1134	rVV	51946	73349	3.38%	0.237%
60	13.894	1136	1138	1142	rVV2	24147	48685	2.25%	0.158%
61	13.974	1142	1145	1147	rVV	740042	780169	35.98%	2.526%
62	14.008	1147	1148	1151	rVV3	27900	54802	2.53%	0.177%
63	14.088	1153	1155	1158	rVB	170669	176709	8.15%	0.572%
64	14.248	1166	1169	1171	rBV	1099799	1129969	52.12%	3.659%
65	14.465	1186	1188	1191	rVV	902304	1006127	46.40%	3.258%
66	14.534	1191	1194	1196	rVV2	39769	51785	2.39%	0.168%
67	14.671	1201	1206	1208	rVB	101611	228932	10.56%	0.741%
68	14.751	1211	1213	1215	rBV	119746	127436	5.88%	0.413%
69	14.785	1215	1216	1218	rVV	74900	84433	3.89%	0.273%
70	14.877	1222	1224	1225	rBV	30834	56354	2.60%	0.182%
71	14.900	1225	1226	1228	rVV	76160	72402	3.34%	0.234%

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72	14.934	1228	1229	1234	rVB	61053	82309	3.80%	0.267%
73	15.197	1248	1252	1255	rVB2	20524	42686	1.97%	0.138%
74	15.277	1255	1259	1262	rBV4	19716	58419	2.69%	0.189%
75	15.425	1269	1272	1274	rBV2	28537	50534	2.33%	0.164%
76	15.471	1274	1276	1279	rVV2	72456	160435	7.40%	0.519%
77	15.654	1289	1292	1296	rBV3	47271	93739	4.32%	0.304%
78	15.745	1296	1300	1302	rVV2	495003	1011907	46.67%	3.276%
79	15.780	1302	1303	1305	rVV	253474	190089	8.77%	0.615%
80	15.871	1309	1311	1314	rBV	52335	57873	2.67%	0.187%
81	16.077	1324	1329	1330	rBV3	61363	166994	7.70%	0.541%
82	16.168	1334	1337	1339	rVV2	38996	97766	4.51%	0.317%
83	16.248	1341	1344	1346	rVV	82565	131787	6.08%	0.427%
84	16.591	1369	1374	1381	rBV	71867	271092	12.50%	0.878%
85	17.048	1411	1414	1419	rVB	199489	468265	21.60%	1.516%
86	17.163	1422	1424	1426	rVV	57906	68495	3.16%	0.222%
87	17.368	1440	1442	1445	rVV	175982	221818	10.23%	0.718%
88	17.426	1445	1447	1451	rVV	267826	372879	17.20%	1.207%
89	17.494	1451	1453	1460	rVV2	342705	523925	24.16%	1.696%
90	17.677	1466	1469	1477	rVB7	30492	114969	5.30%	0.372%
91	17.814	1477	1481	1482	rBV3	20314	49547	2.29%	0.160%
92	18.008	1496	1498	1501	rVB	42406	67484	3.11%	0.219%
93	18.134	1507	1509	1512	rVB	23595	44147	2.04%	0.143%
94	18.511	1539	1542	1545	rBV4	29149	83168	3.84%	0.269%
95	18.637	1550	1553	1556	rVB4	30219	68319	3.15%	0.221%
96	18.774	1562	1565	1570	rVB2	112190	223495	10.31%	0.724%
97	18.934	1576	1579	1582	rBV3	20609	56154	2.59%	0.182%
98	19.140	1594	1597	1602	rBV	98160	183489	8.46%	0.594%
99	19.334	1611	1614	1619	rVB	49458	95641	4.41%	0.310%
100	21.266	1776	1783	1790	rBV	49483	232131	10.71%	0.752%

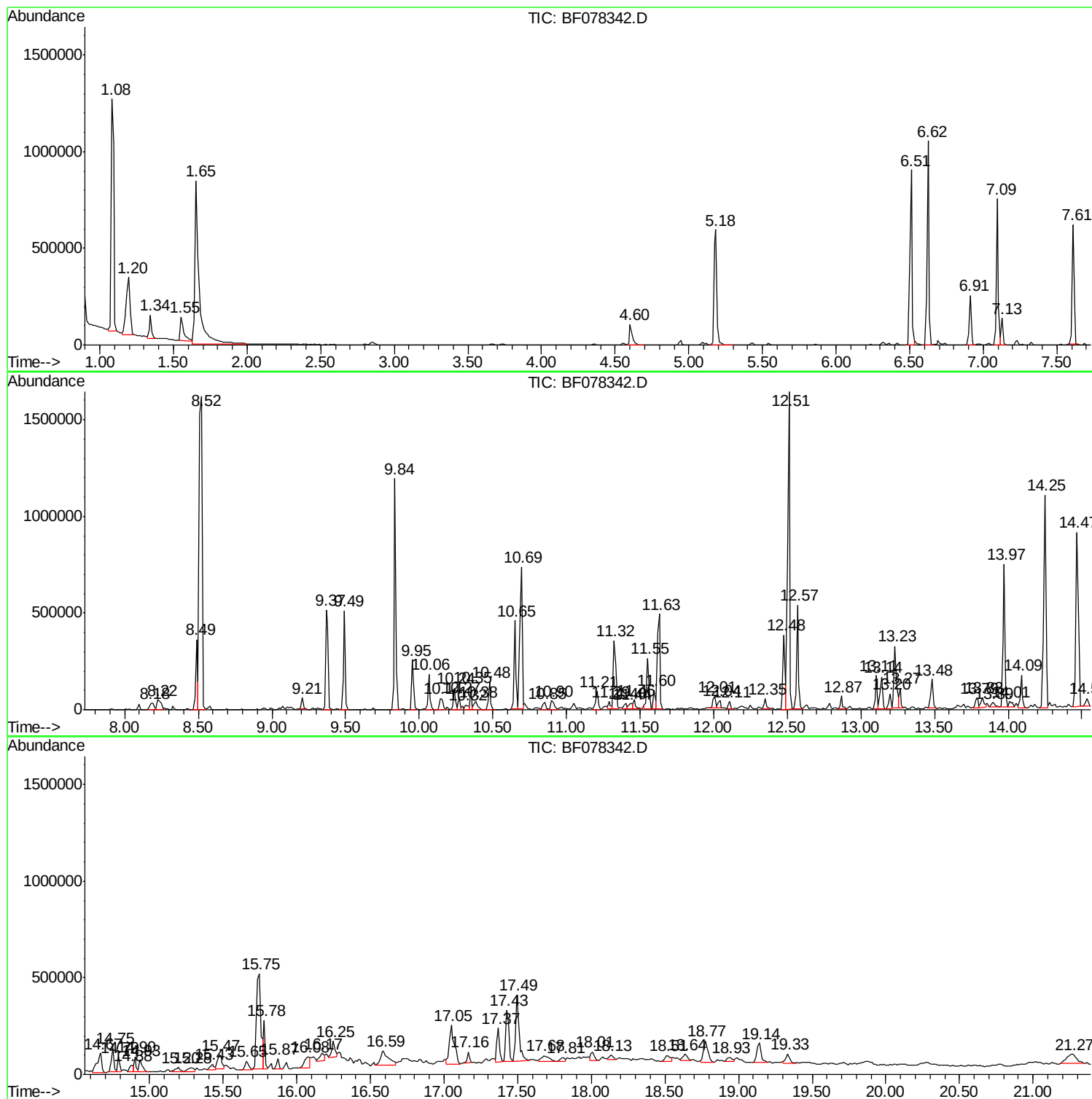
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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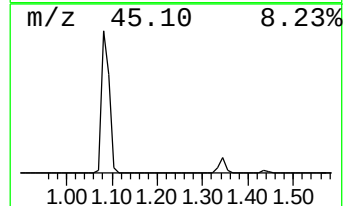
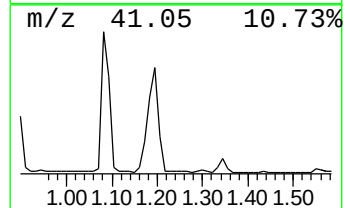
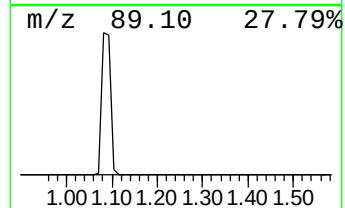
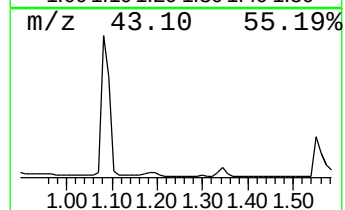
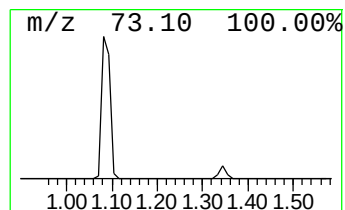
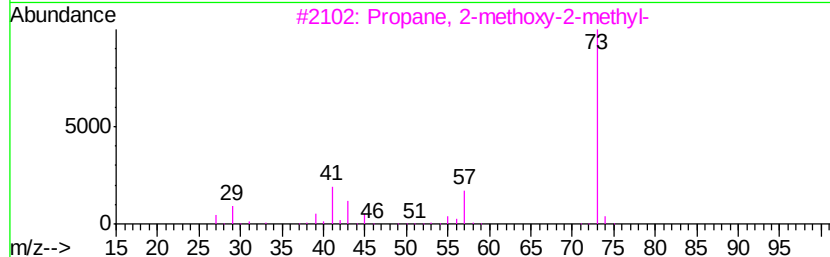
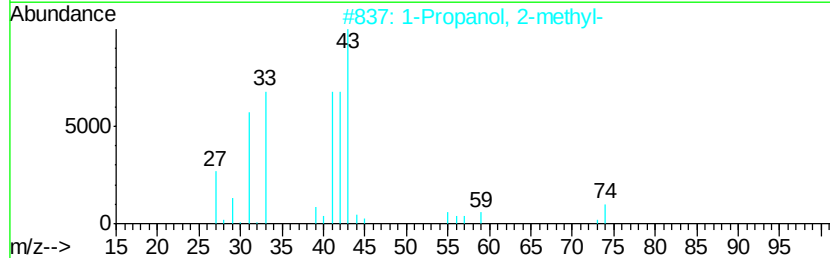
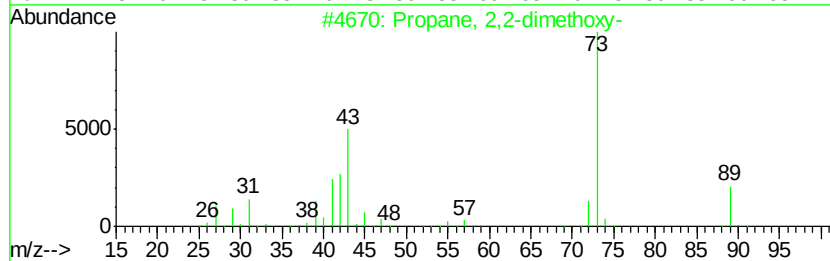
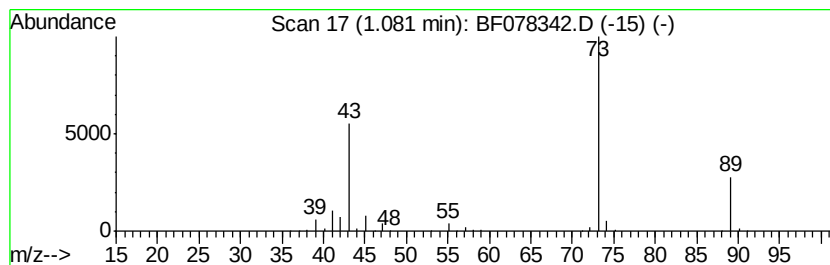
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Propane, 2,2-dimethoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.08	129.99 ng	1533720	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	72
2		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
3		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9
4		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	9
5		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9



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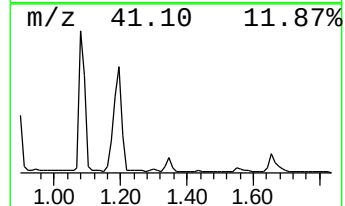
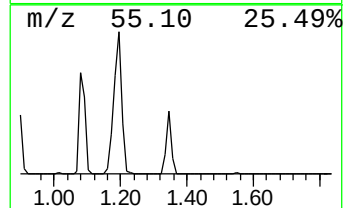
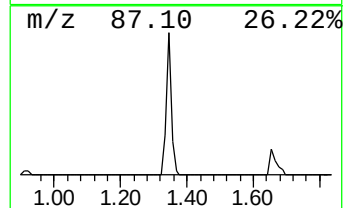
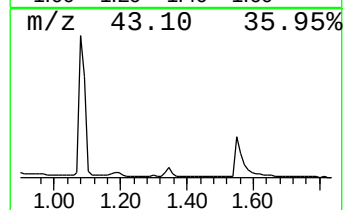
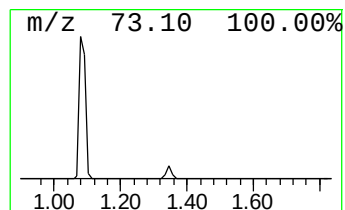
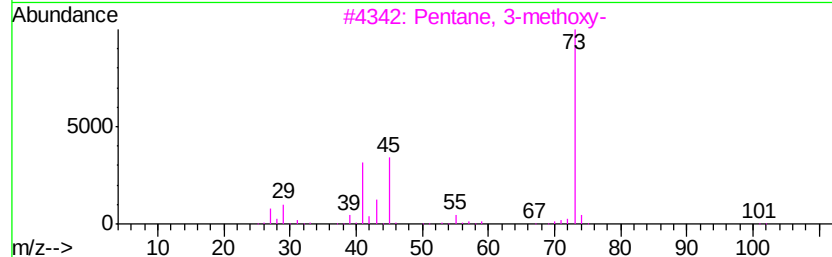
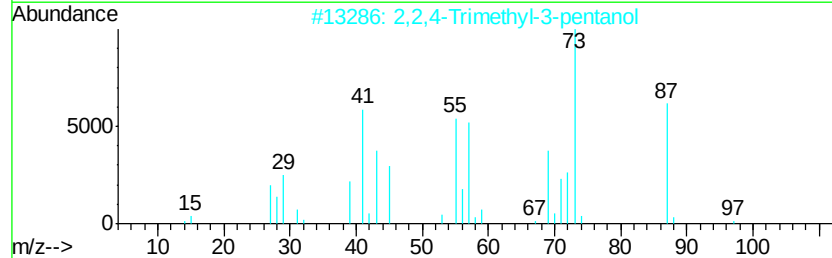
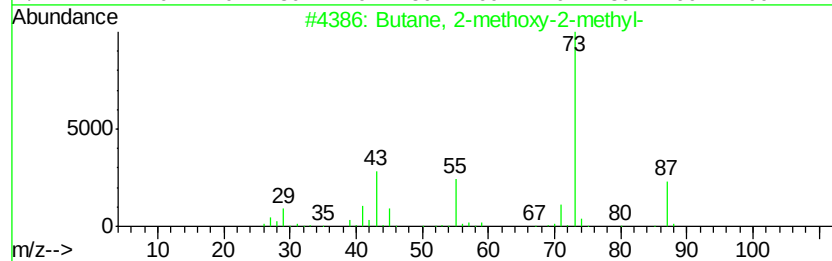
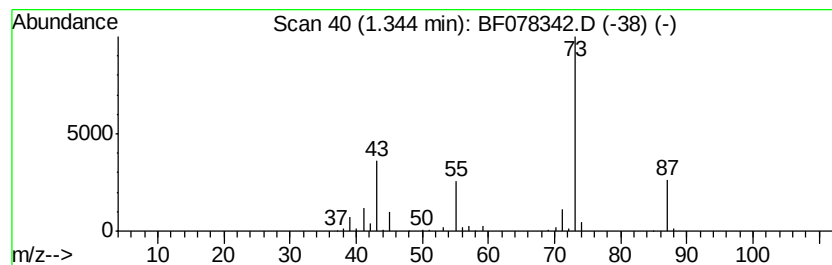
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Butane, 2-methoxy-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.34	11.33 ng	133633	1,4-Dichlorobenzene-d4	6.91

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		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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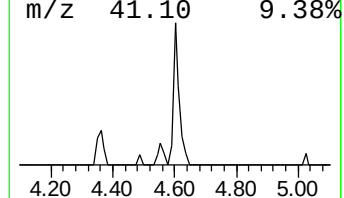
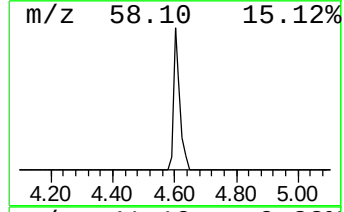
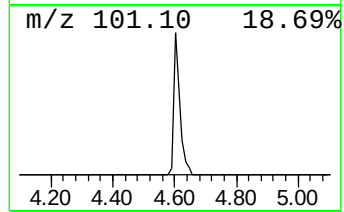
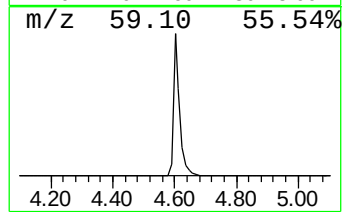
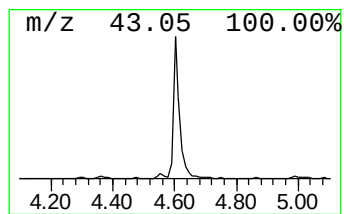
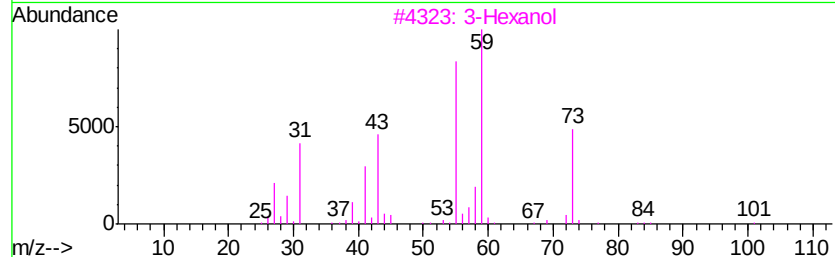
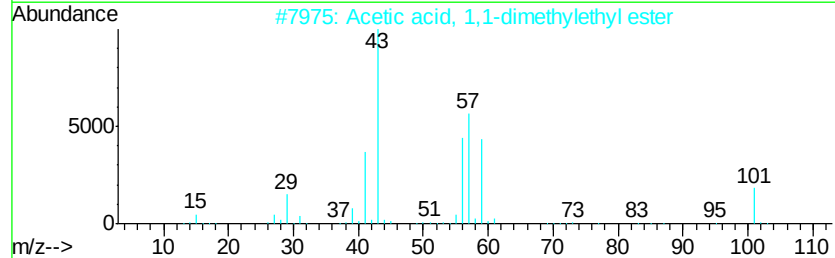
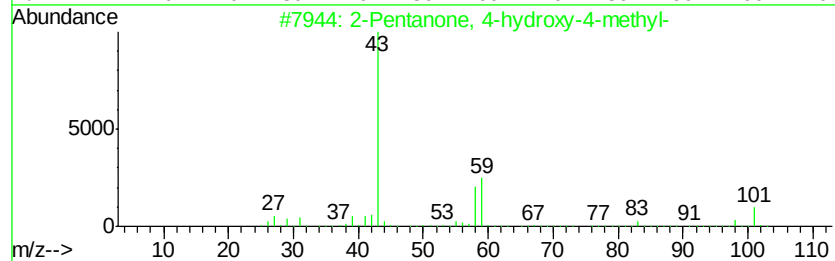
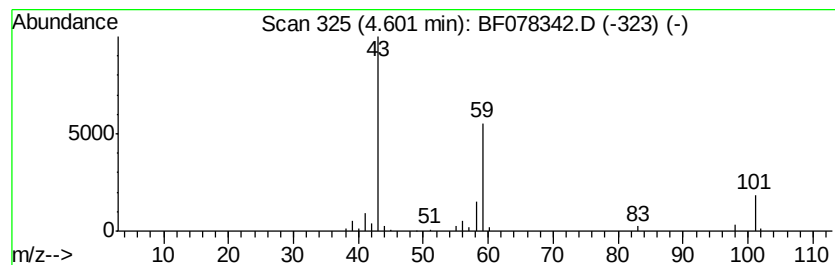
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.60	12.16 ng	143521	1,4-Dichlorobenzene-d4	6.91

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59	
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39	
3	3-Hexanol	102	C6H14O	000623-37-0	33	
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27	
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25	



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Data File : BF078342.D
Acq On : 8 Apr 2015 19:07
Operator : TP/IZ
Sample : G1793-08
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-6D

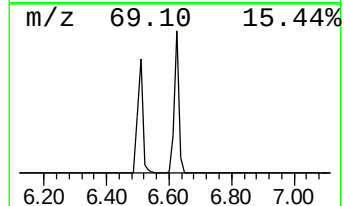
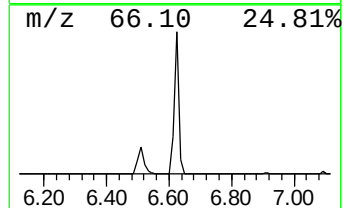
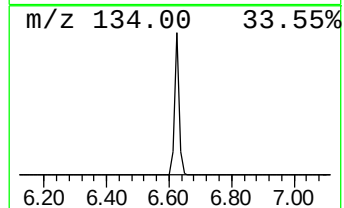
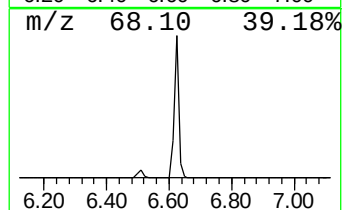
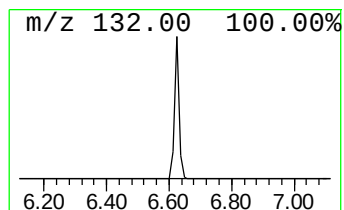
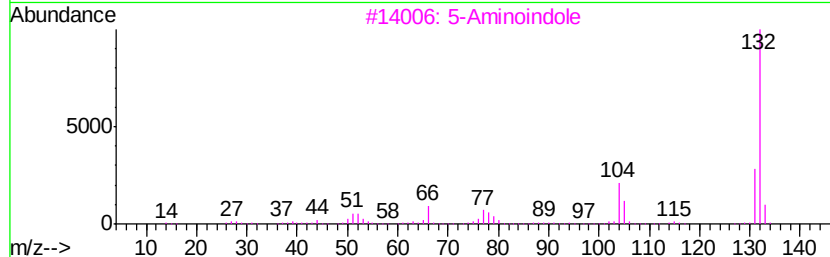
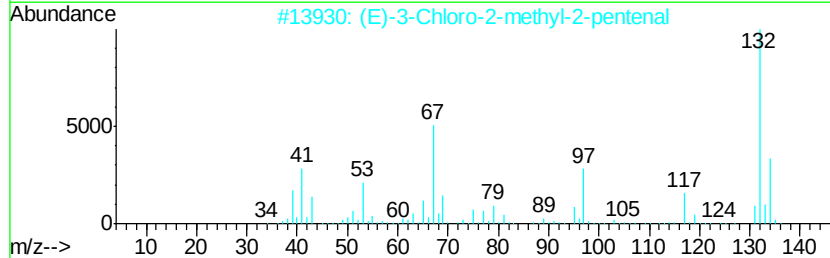
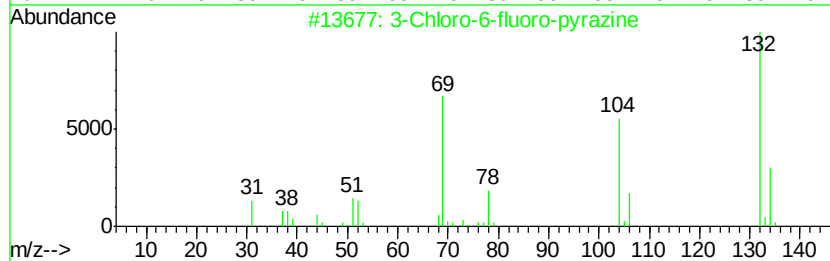
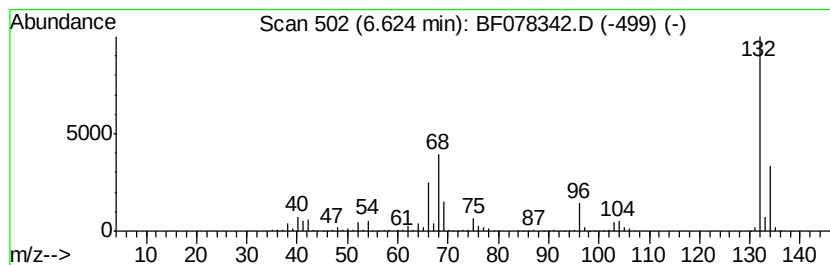
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 unknown6.62 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	83.62 ng	986643	1,4-Dichlorobenzene-d4	6.91

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040815\
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Sample : G1793-08
Misc :
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ClientSampleId :
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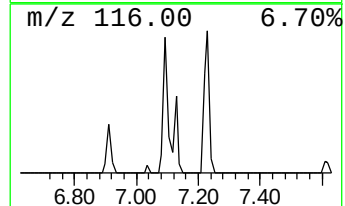
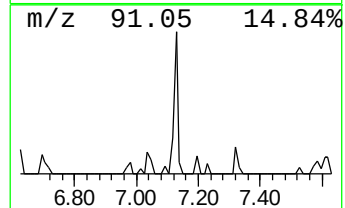
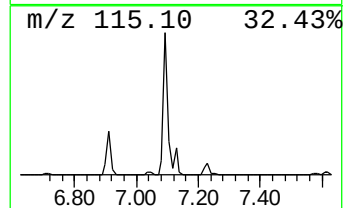
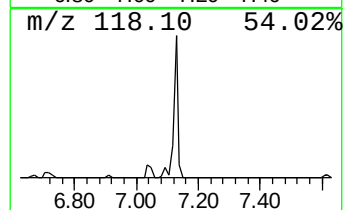
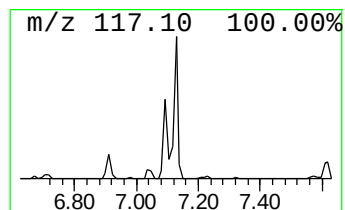
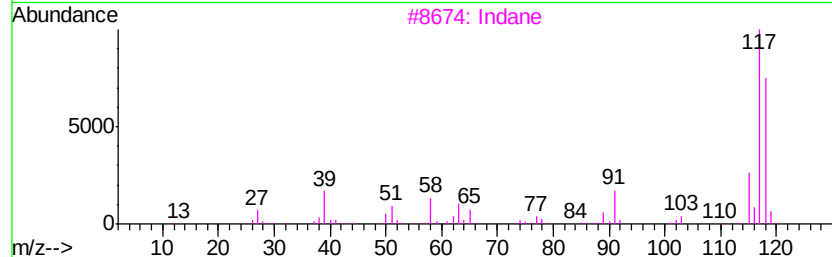
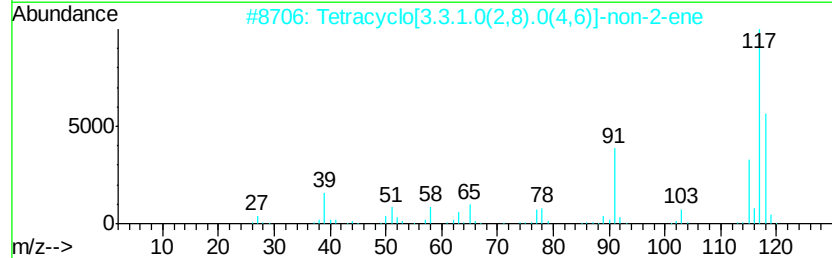
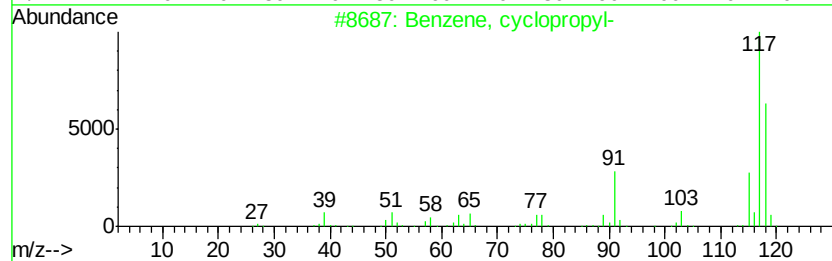
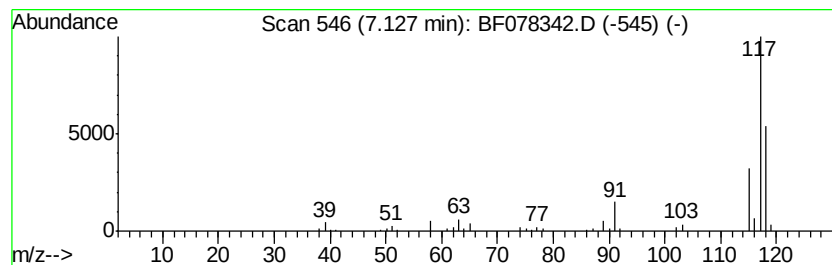
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, cyclopropyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.13	8.64 ng	101976	1,4-Dichlorobenzene-d4	6.91

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, cyclopropyl-	118	C9H10	000873-49-4	74
2		Tetracyclo[3.3.1.0(2,8).0(4,6)]-	118	C9H10	1000191-13-7	72
3		Indane	118	C9H10	000496-11-7	64
4		trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	59
5		Deltacyclene	118	C9H10	007785-10-6	59



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040815\
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Acq On : 8 Apr 2015 19:07
Operator : TP/IZ
Sample : G1793-08
Misc :
ALS Vial : 4 Sample Multiplier: 1

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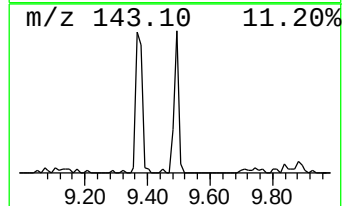
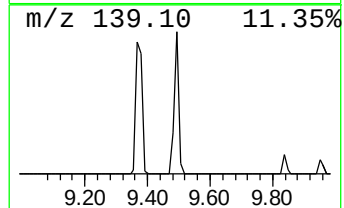
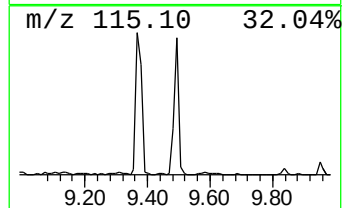
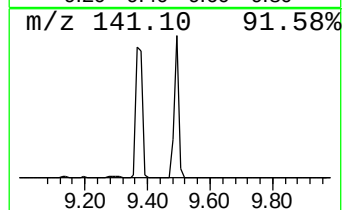
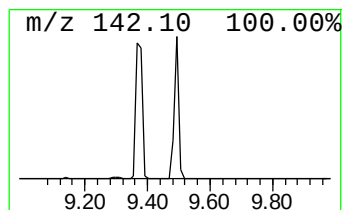
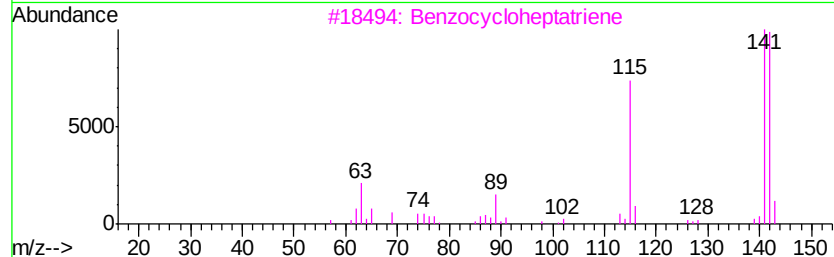
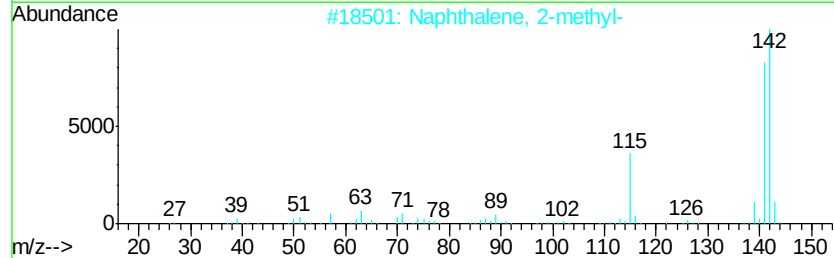
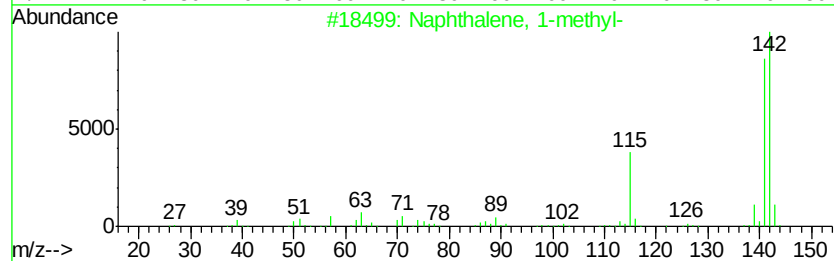
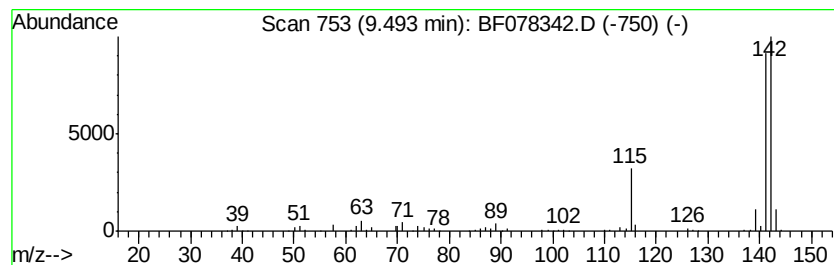
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.49	24.89 ng	483659	Naphthalene-d8	8.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Benzocycloheptatriene	142	C11H10	000264-09-5	91
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5		1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	86



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Sample : G1793-08
Misc :
ALS Vial : 4 Sample Multiplier: 1

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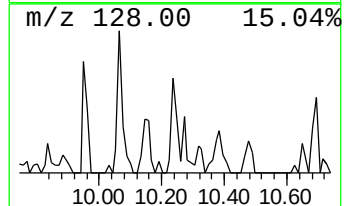
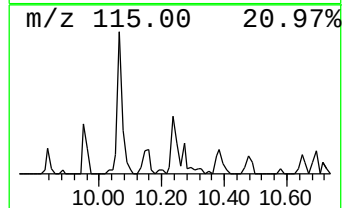
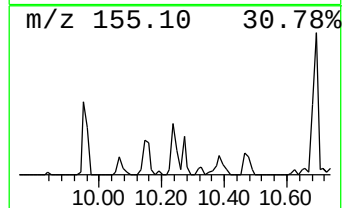
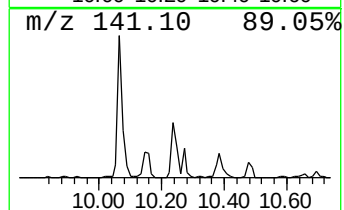
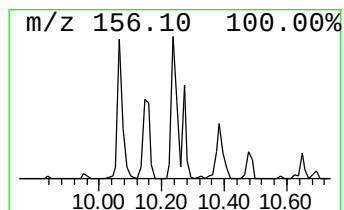
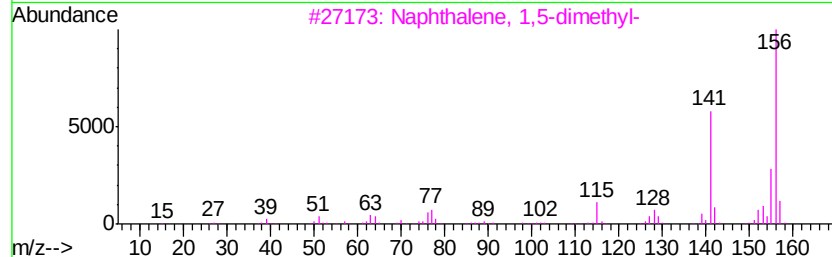
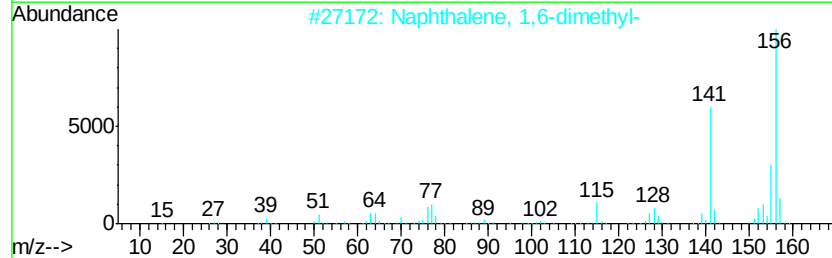
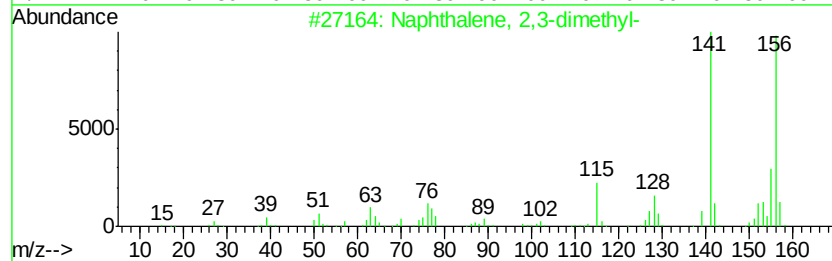
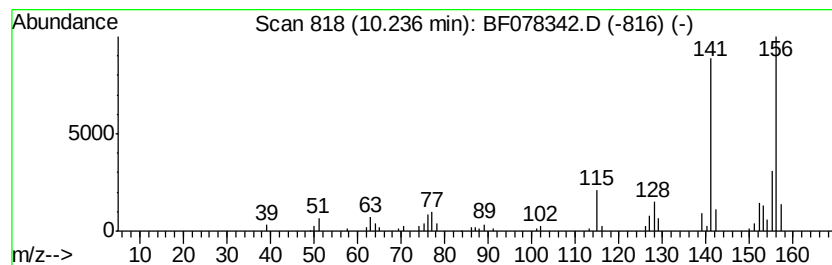
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Naphthalene, 2,3-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.24	5.68 ng	128697	Acenaphthene-d10	10.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
3			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
4			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
5			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97



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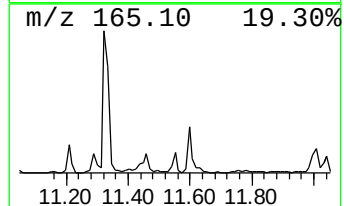
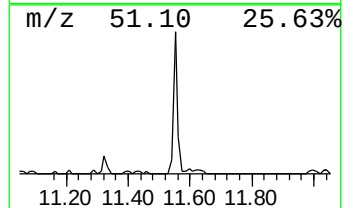
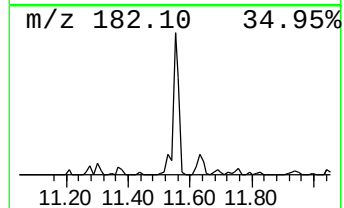
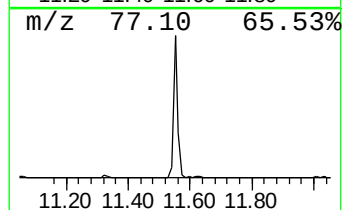
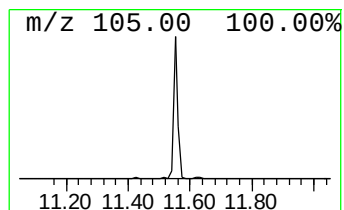
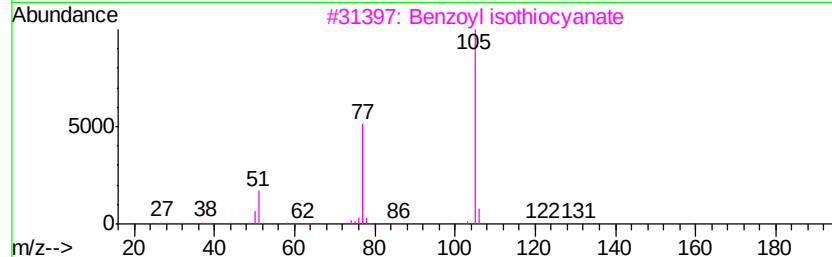
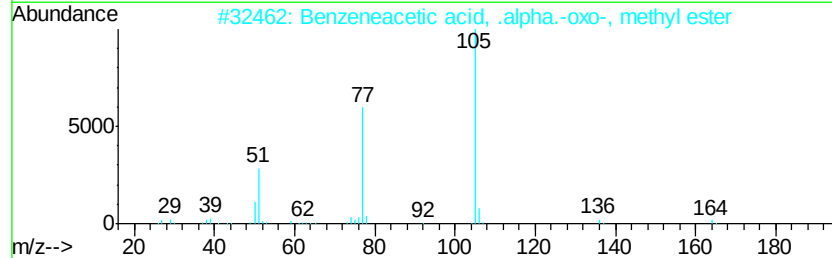
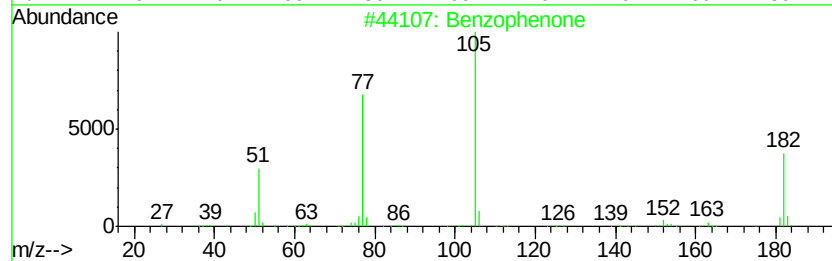
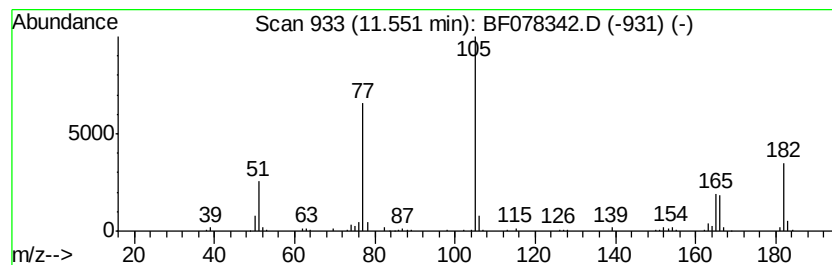
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzophenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.55	11.54 ng	261384	Acenaphthene-d10	10.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	81
2			Benzeneacetic acid, .alpha.-oxo-...	164	C9H8O3	015206-55-0	49
3			Benzoyl isothiocyanate	163	C8H5NOS	000532-55-8	47
4			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	47
5			Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	47



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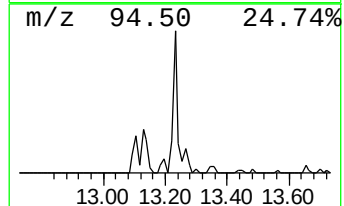
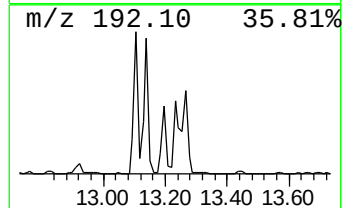
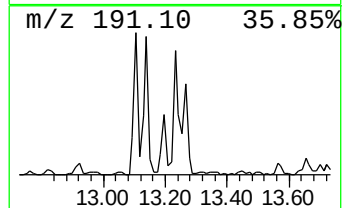
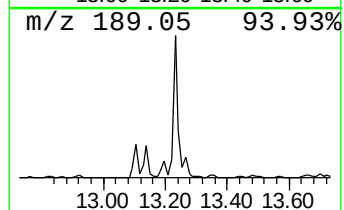
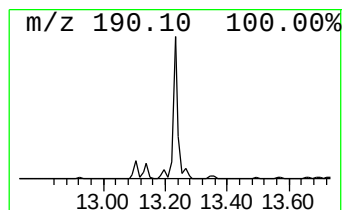
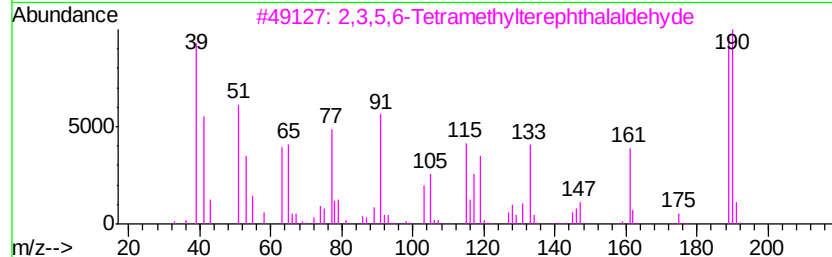
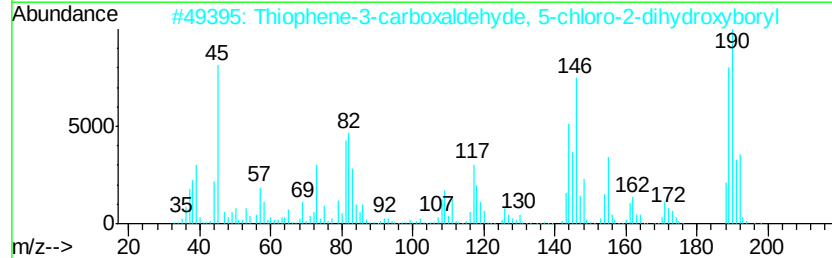
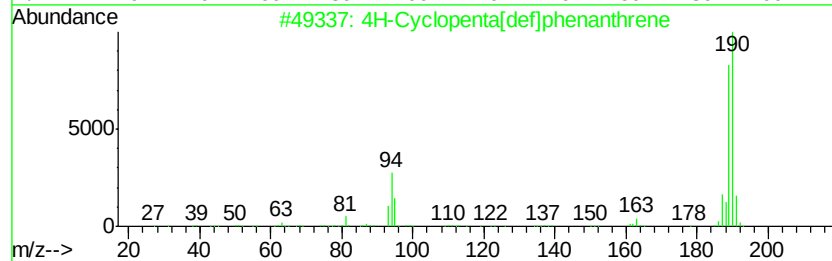
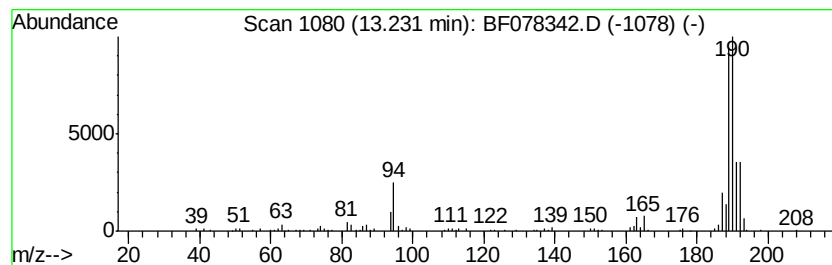
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 4H-Cyclopenta[def]phenanthrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.23	18.76 ng	366119	Phenanthrene-d10	12.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	53
3		2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	47
4		1,4-Naphthalenedione, 2,5-dihydr...	190	C10H6O4	004923-55-1	17
5		Hydrazinecarboxaldehyde, 2-[3-(m...	190	C4H6N4O5	038362-25-3	14



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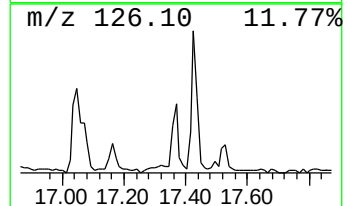
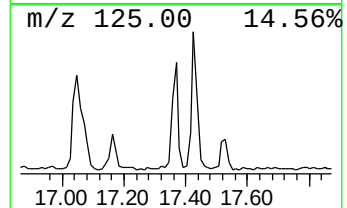
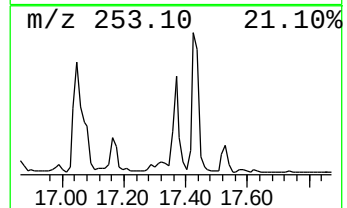
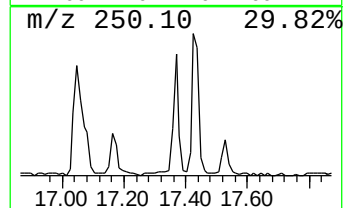
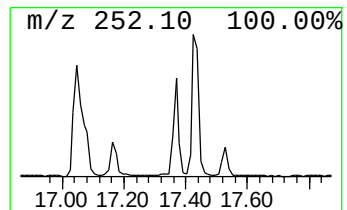
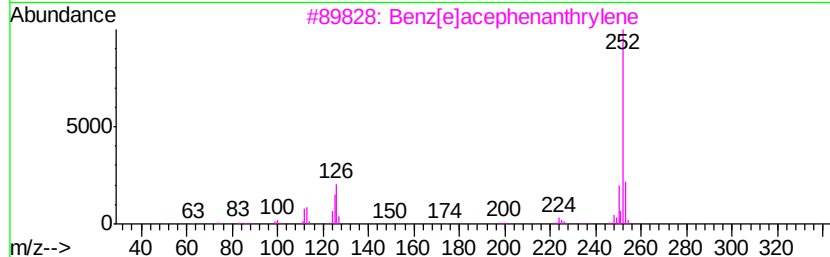
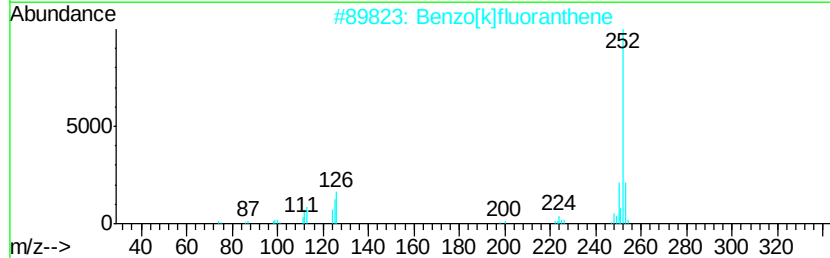
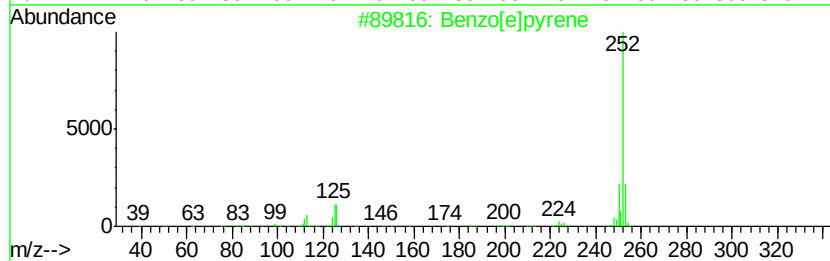
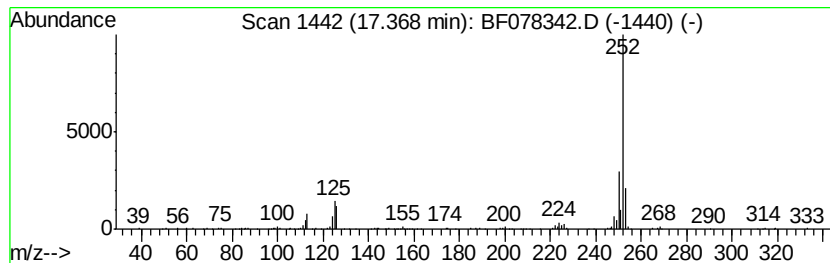
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Benzo[e]pyrene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.37	8.47 ng	221818	Perylene-d12	17.49

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3		Benzo[e]acephenanthrylene	252	C20H12	000205-99-2	94
4		Benzo[a]pyrene	252	C20H12	000050-32-8	93
5		Perylene	252	C20H12	000198-55-0	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040815\
Data File : BF078342.D
Acq On : 8 Apr 2015 19:07
Operator : TP/IZ
Sample : G1793-08
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-6D

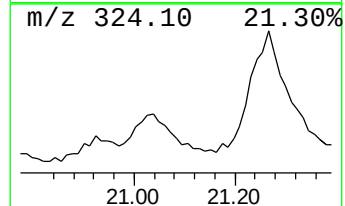
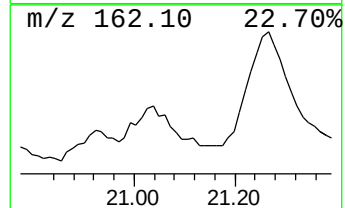
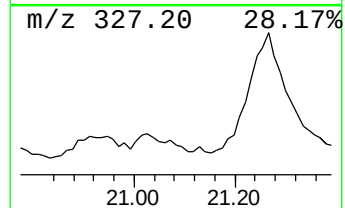
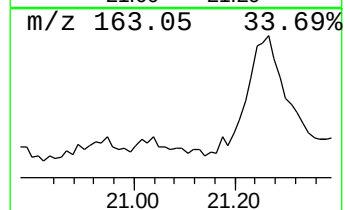
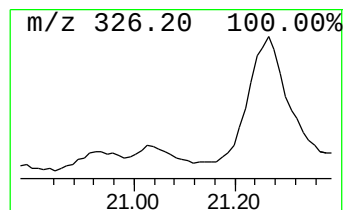
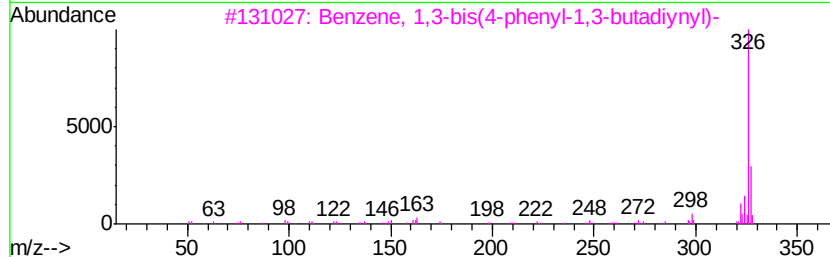
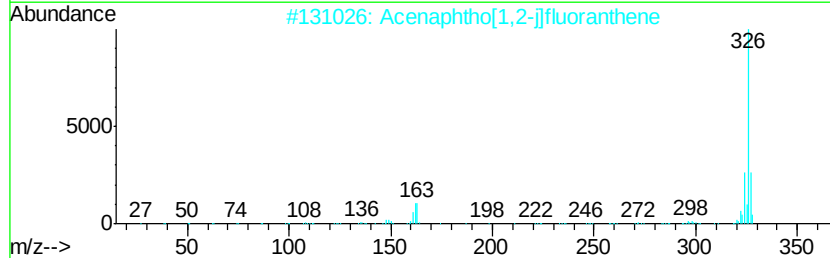
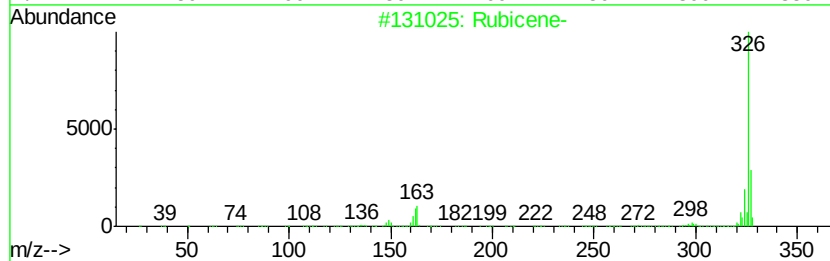
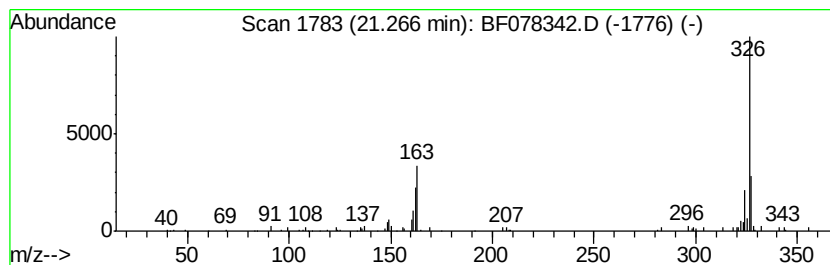
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Rubicene- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.27	8.86 ng	232131	Perylene-d12	17.49

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Rubicene-	326	C26H14	000197-61-5	83
2			Acenaphtho[1,2-i]fluoranthene	326	C26H14	000193-21-5	64
3			Benzene, 1,3-bis(4-phenyl-1,3-butadienyl)-	326	C26H14	037902-13-9	49
4			Bis(pentamethylcyclopentadienyl)iron	326	C20H30Fe	012126-50-0	47
5			9-Anthracenamine, N-[(4-nitrophenyl)imino]-	326	C21H14N2O2	127387-74-0	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040815\
Data File : BF078342.D
Acq On : 8 Apr 2015 19:07
Operator : TP/IZ
Sample : G1793-08
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-6D

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
Propane, 2,2-dime...	1.08	130.0	ng	1533720	1	6.91	235982	20.0
Butane, 2-methoxy...	1.34	11.3	ng	133633	1	6.91	235982	20.0
2-Pentanone, 4-hy...	4.60	12.2	ng	143521	1	6.91	235982	20.0
unknown6.62	6.62	83.6	ng	986643	1	6.91	235982	20.0
Benzene, cyclopro...	7.13	8.6	ng	101976	1	6.91	235982	20.0
Naphthalene, 1-me...	9.49	24.9	ng	483659	2	8.49	388663	20.0
Naphthalene, 2,3-...	10.24	5.7	ng	128697	3	10.65	453028	20.0
Benzophenone	11.55	11.5	ng	261384	3	10.65	453028	20.0
4H-Cyclopenta[def...	13.23	18.8	ng	366119	4	12.48	390410	20.0
Benzo[e]pyrene	17.37	8.5	ng	221818	6	17.49	523925	20.0
Rubicene-	21.27	8.9	ng	232131	6	17.49	523925	20.0