

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040317\
 Data File : BF094135.D
 Acq On : 3 Apr 2017 21:50
 Operator : SJ/MA
 Sample : I2493-10 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LSS-SP08-COMP

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-BF032217.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	2.269	61	65	70	rBV	86282	99315	8.62%	0.515%
2	3.081	201	203	211	rVB	16089	19810	1.72%	0.103%
3	3.522	275	278	285	rVB3	16391	20377	1.77%	0.106%
4	4.010	358	361	370	rBV	77749	82850	7.19%	0.430%
5	4.381	420	424	426	rBV	51416	56017	4.86%	0.291%
6	4.416	426	430	435	rVB	1064353	947339	82.25%	4.915%
7	4.640	464	468	474	rVB	46590	44313	3.85%	0.230%
8	5.363	587	591	594	rBV	55801	54161	4.70%	0.281%
9	5.393	594	596	602	rVB	26886	27341	2.37%	0.142%
10	5.446	602	605	608	rBV	51764	48509	4.21%	0.252%
11	5.487	608	612	618	rBV	1043420	1039763	90.27%	5.394%
12	5.540	618	621	624	rVB	35624	32303	2.80%	0.168%
13	5.628	632	636	640	rBV	1093723	1030716	89.49%	5.348%
14	5.693	644	647	650	rBV	150318	132637	11.52%	0.688%
15	5.887	675	680	684	rBV	904434	806717	70.04%	4.185%
16	5.957	689	692	695	rBV	50682	52326	4.54%	0.271%
17	6.063	706	710	714	rBV	898732	802094	69.64%	4.161%
18	6.181	726	730	732	rBV2	21555	20183	1.75%	0.105%
19	6.210	732	735	740	rVB	33069	35379	3.07%	0.184%
20	6.263	740	744	747	rBV	57164	55423	4.81%	0.288%
21	6.351	756	759	762	rBV	24710	25885	2.25%	0.134%
22	6.434	770	773	775	rBV	30435	26458	2.30%	0.137%
23	6.463	775	778	782	rVB2	22416	22971	1.99%	0.119%
24	6.528	782	789	799	rBV	568351	571432	49.61%	2.965%
25	6.792	831	834	837	rBV	42031	33528	2.91%	0.174%
26	6.822	837	839	844	rVB	47371	44411	3.86%	0.230%
27	7.004	865	870	873	rBV3	27393	28938	2.51%	0.150%
28	7.087	879	884	889	rBV4	50281	78028	6.77%	0.405%
29	7.134	889	892	895	rVB3	19908	19941	1.73%	0.103%
30	7.257	909	913	916	rVB2	15596	18753	1.63%	0.097%
31	7.292	916	919	923	rBV3	35798	38459	3.34%	0.200%
32	7.345	923	928	930	rVV4	17430	22556	1.96%	0.117%
33	7.392	931	936	939	rVV	1139077	1104754	95.91%	5.732%
34	7.416	939	940	945	rVV2	93099	77680	6.74%	0.403%

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 Integrator: RTE
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 Stop Thrs : 0

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	7.463	945	948	953	rVB5	16771	20245	1.76%	0.105%
36	7.822	1003	1009	1012	rBV2	39959	46300	4.02%	0.240%
37	8.257	1079	1083	1087	rBV	121490	112997	9.81%	0.586%
38	8.322	1090	1094	1100	rVV	94106	119890	10.41%	0.622%
39	8.375	1100	1103	1107	rVB	57939	54259	4.71%	0.282%
40	8.710	1155	1160	1164	rVB	1237875	1135910	98.62%	5.893%
41	9.028	1211	1214	1218	rVB2	26640	39366	3.42%	0.204%
42	9.122	1227	1230	1233	rBV	25169	24012	2.08%	0.125%
43	9.151	1233	1235	1238	rVB	18017	16937	1.47%	0.088%
44	9.210	1238	1245	1248	rBV	89981	92770	8.05%	0.481%
45	9.263	1250	1254	1260	rVB5	16280	28182	2.45%	0.146%
46	9.357	1263	1270	1275	rBV2	73228	116709	10.13%	0.606%
47	9.533	1292	1300	1304	rBV	1020773	1067568	92.69%	5.539%
48	9.569	1304	1306	1309	rVB2	34429	30747	2.67%	0.160%
49	9.669	1320	1323	1326	rBV4	13617	16755	1.45%	0.087%
50	9.781	1339	1342	1348	rBV2	28451	48938	4.25%	0.254%
51	9.845	1348	1353	1356	rVV4	20184	24261	2.11%	0.126%
52	9.939	1365	1369	1372	rBV6	14405	23487	2.04%	0.122%
53	10.081	1389	1393	1396	rBV2	25106	23384	2.03%	0.121%
54	10.122	1398	1400	1405	rVB	42352	38712	3.36%	0.201%
55	10.204	1411	1414	1418	rBV	57870	56093	4.87%	0.291%
56	10.404	1444	1448	1451	rBV4	27254	33177	2.88%	0.172%
57	10.504	1460	1465	1471	rBV	433080	475538	41.29%	2.467%
58	10.592	1476	1480	1486	rVB9	26936	49809	4.32%	0.258%
59	10.710	1497	1500	1502	rBV	51290	51318	4.46%	0.266%
60	10.957	1539	1542	1543	rBV3	16242	18480	1.60%	0.096%
61	11.110	1564	1568	1571	rVB2	59363	71063	6.17%	0.369%
62	11.233	1585	1589	1592	rBV2	47555	71322	6.19%	0.370%
63	11.263	1592	1594	1598	rVB	46241	40241	3.49%	0.209%
64	11.304	1598	1601	1606	rBV5	41092	61511	5.34%	0.319%
65	11.357	1606	1610	1612	rBV	833402	812524	70.54%	4.216%
66	11.386	1612	1615	1619	rVB	607365	562395	48.83%	2.918%
67	11.445	1622	1625	1628	rVB	141337	145376	12.62%	0.754%
68	11.651	1657	1660	1664	rBV2	71130	82088	7.13%	0.426%
69	11.869	1695	1697	1700	rVB2	54601	46330	4.02%	0.240%
70	11.980	1713	1716	1719	rBV	87957	95444	8.29%	0.495%
71	12.016	1719	1722	1726	rVB	131898	135248	11.74%	0.702%

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Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
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72	12.110	1735	1738	1741	rBV2	150848	168967	14.67%	0.877%
73	12.351	1776	1779	1785	rBV2	107026	173611	15.07%	0.901%
74	12.851	1860	1864	1868	rBV	964896	960435	83.38%	4.983%
75	13.127	1907	1911	1915	rVB	921375	998000	86.65%	5.178%
76	13.204	1921	1924	1929	rVB	430519	506395	43.96%	2.627%
77	13.333	1943	1946	1952	rVB	682790	685124	59.48%	3.555%
78	13.951	2048	2051	2054	rVB	192878	176360	15.31%	0.915%
79	14.616	2159	2164	2167	rBV2	754130	1151820	100.00%	5.976%
80	14.645	2167	2169	2171	rVV	260501	214497	18.62%	1.113%
81	15.845	2370	2373	2376	rBV	296649	397156	34.48%	2.061%
82	16.251	2439	2442	2446	rVB2	418622	431504	37.46%	2.239%

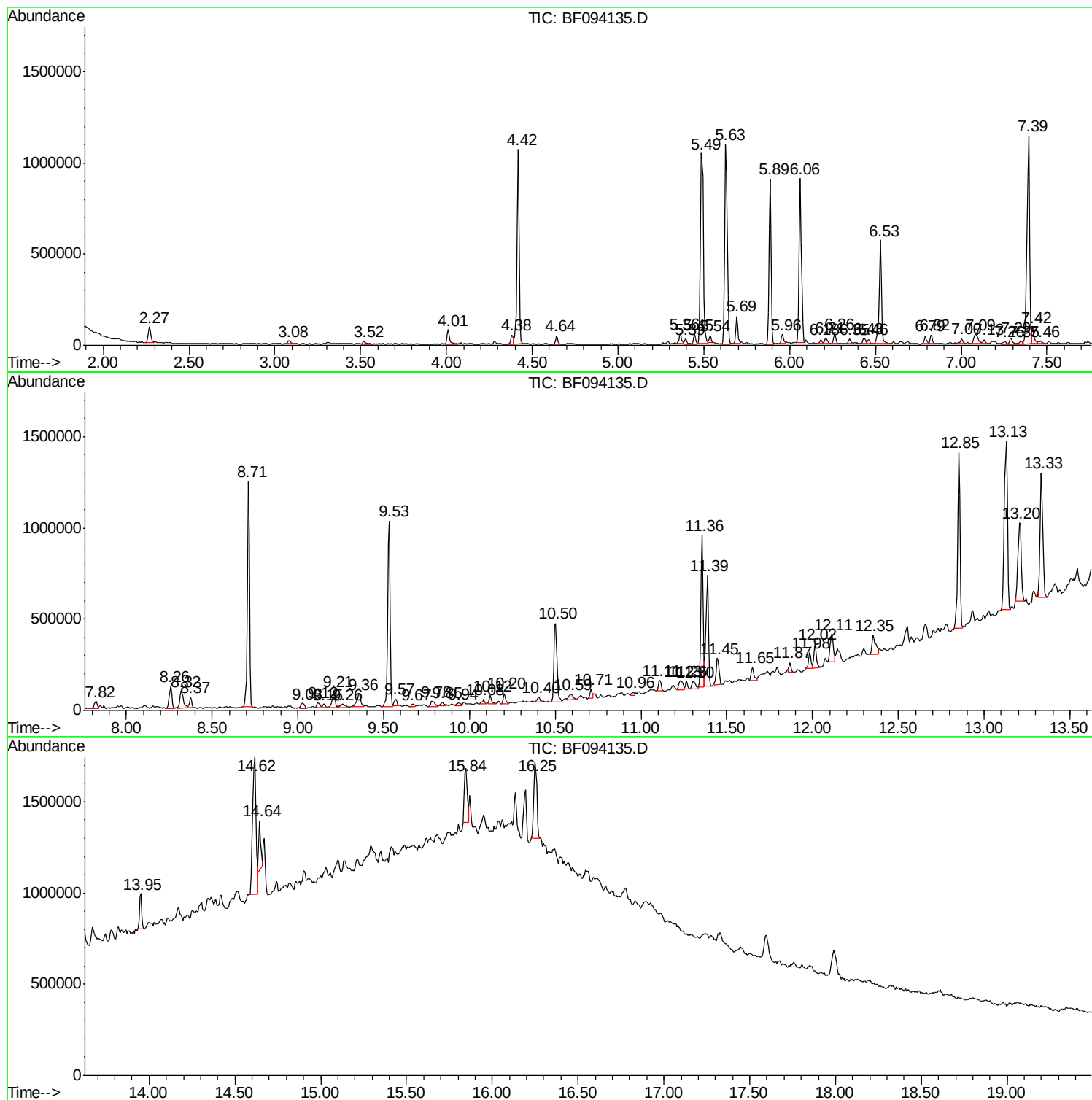
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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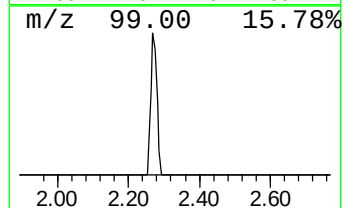
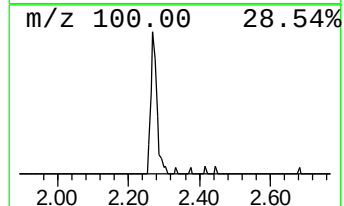
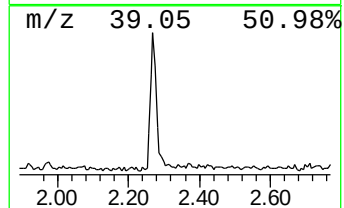
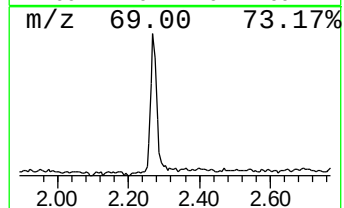
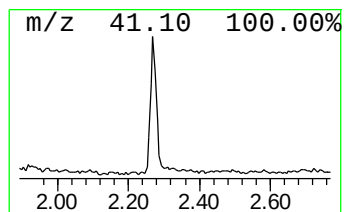
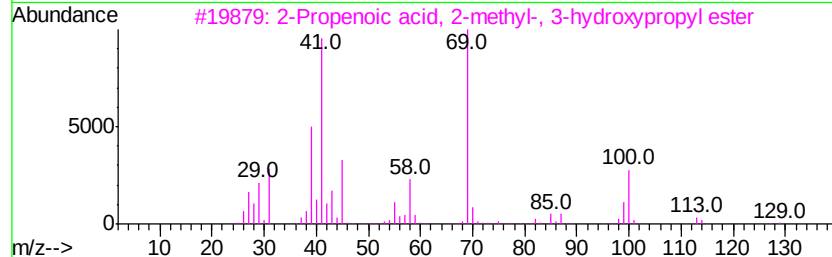
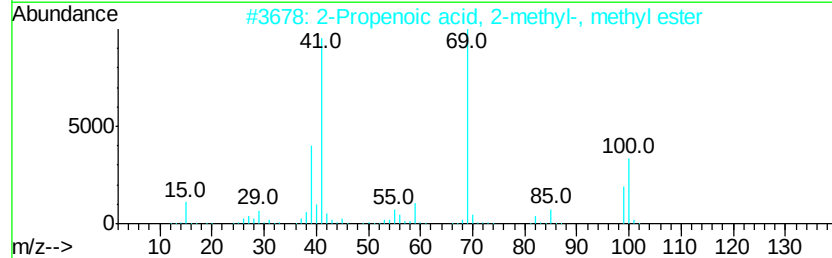
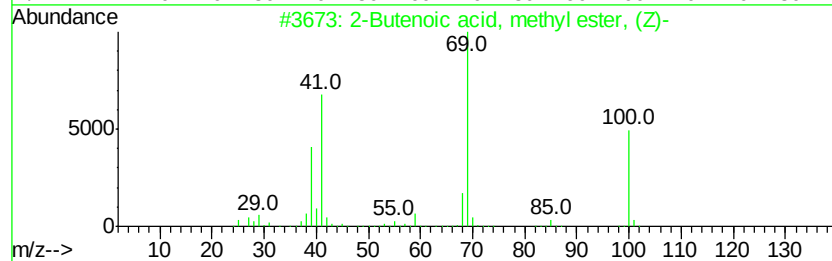
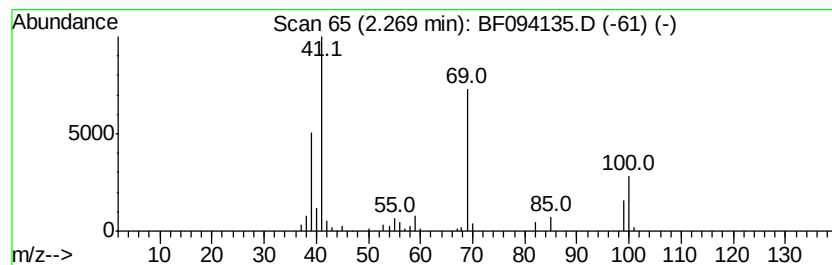
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Butenoic acid, methyl est... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.27	2.46 ng	99315	1,4-Dichlorobenzene-d4	5.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butenoic acid, methyl ester, (Z)-	100	C5H8O2	004358-59-2	87
2		2-Propenoic acid, 2-methyl-, met...	100	C5H8O2	000080-62-6	81
3		2-Propenoic acid, 2-methyl-, 3-h...	144	C7H12O3	002761-09-3	64
4		Methacrylic anhydride	154	C8H10O3	000760-93-0	53
5		Cyclopropanecarboxylic acid, met...	100	C5H8O2	002868-37-3	45



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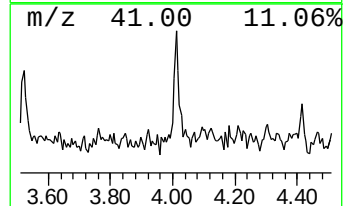
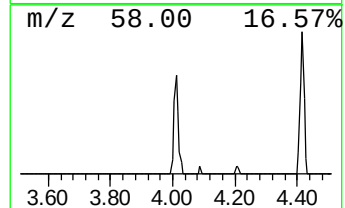
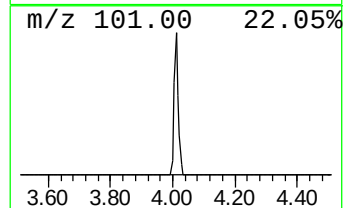
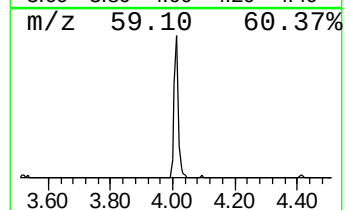
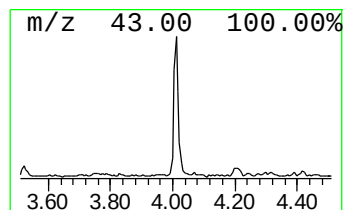
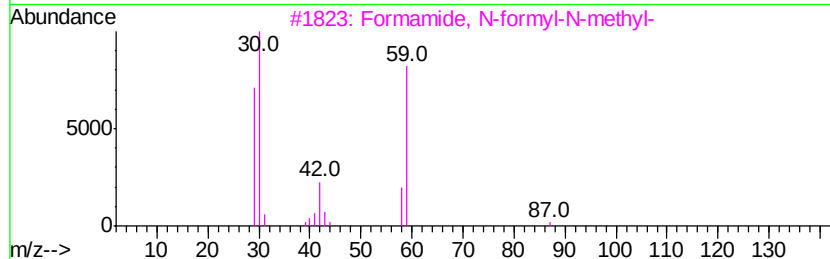
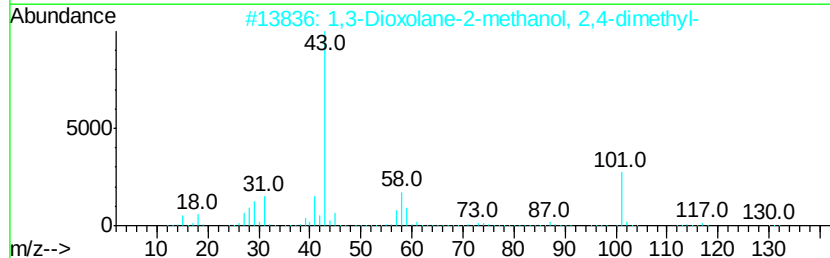
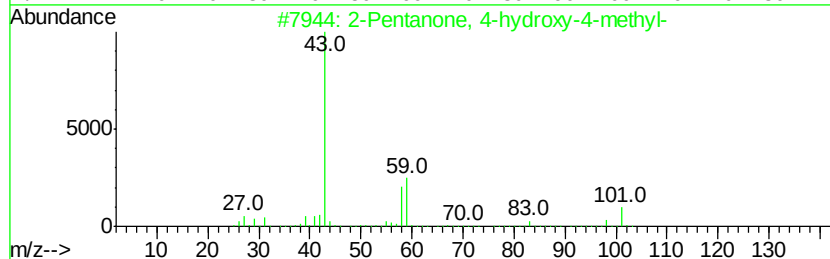
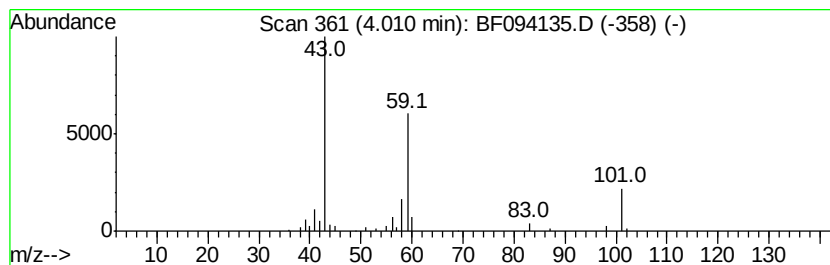
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.01	2.05 ng	82850	1,4-Dichlorobenzene-d4	5.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			1,3-Dioxolane-2-methanol, 2,4-di...	132	C6H12O3	053951-43-2	23
3			Formamide, N-formyl-N-methyl-	87	C3H5NO2	018197-25-6	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
5			Acetone	58	C3H6O	000067-64-1	9



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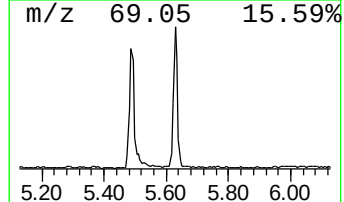
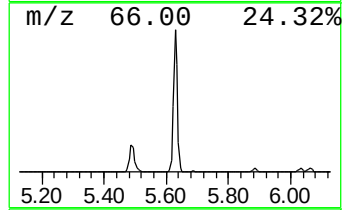
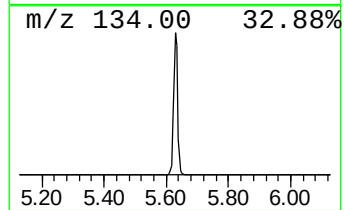
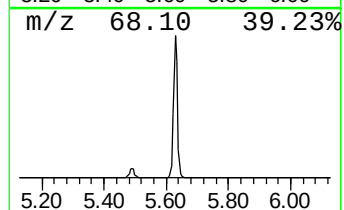
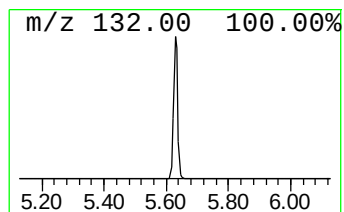
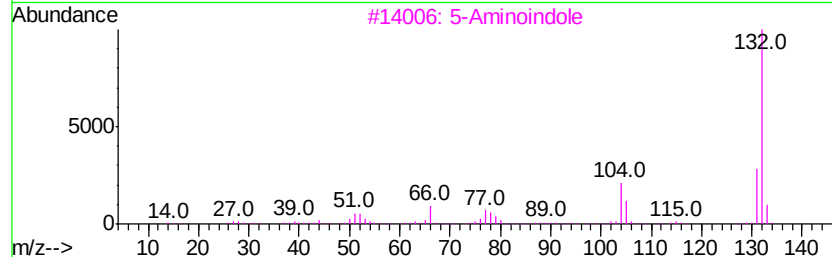
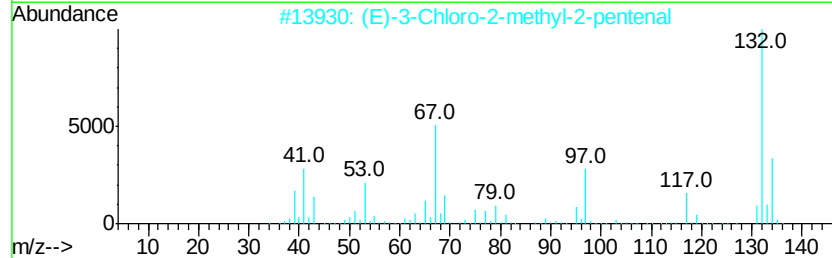
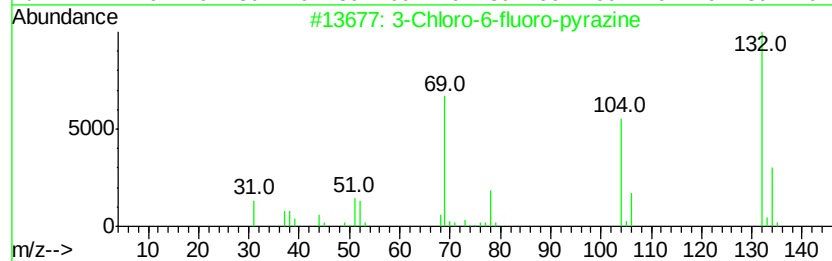
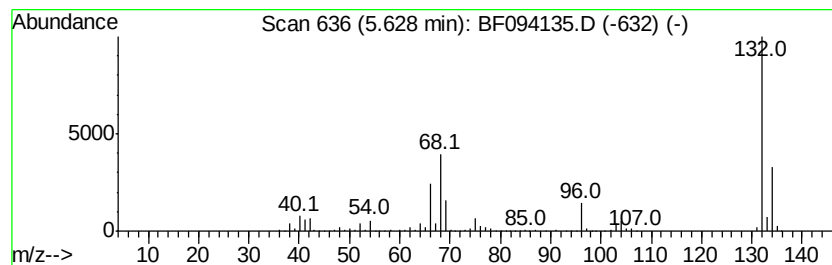
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 unknown5.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.63	25.55 ng	1030720	1,4-Dichlorobenzene-d4	5.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3		Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	(E)-3		Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
3	5		Aminoindole	132	C8H8N2	005192-03-0	12
4	Tranylcypromine-propionyl			189	C12H15NO	1000123-86-3	10
5	4		Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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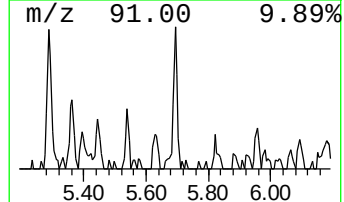
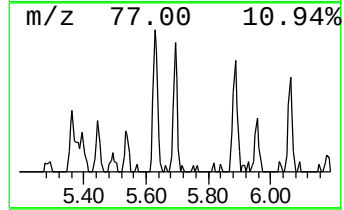
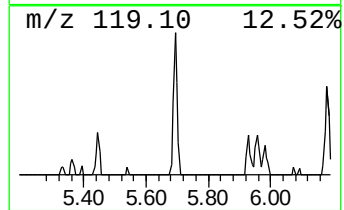
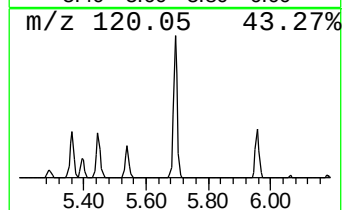
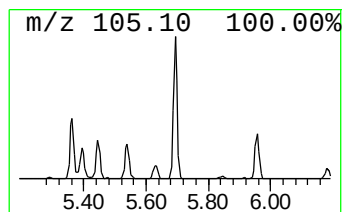
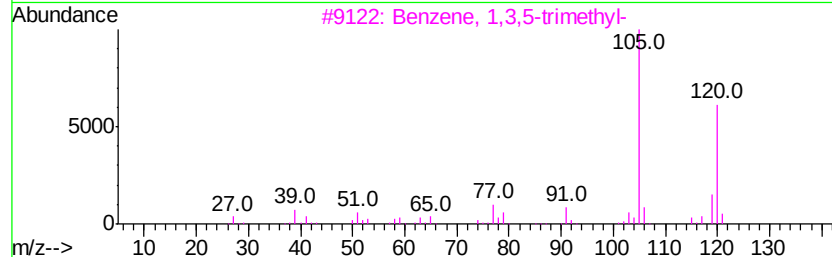
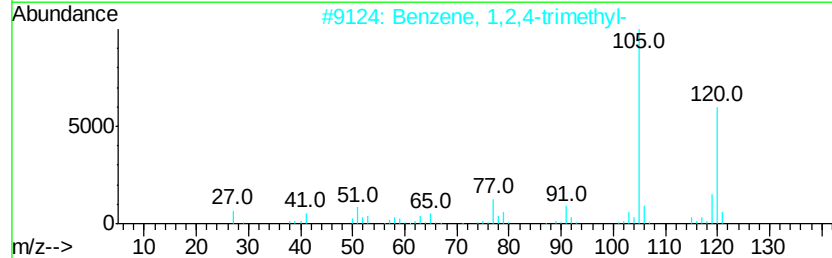
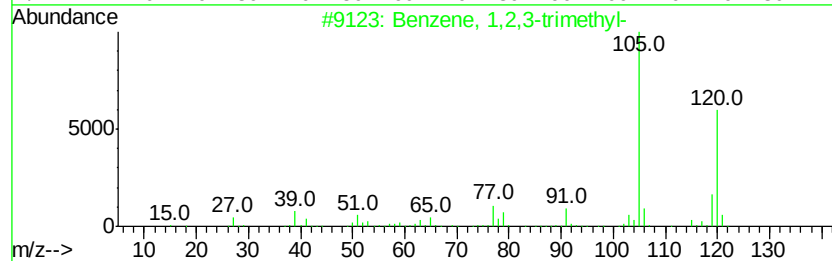
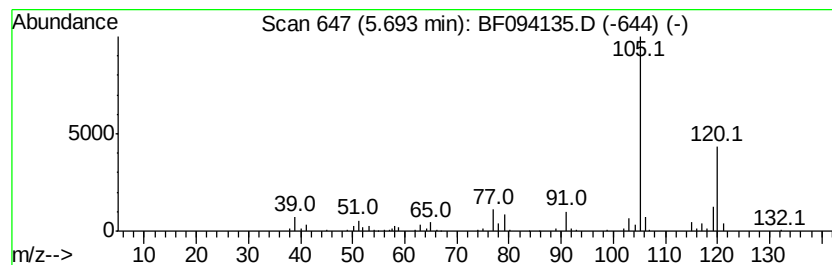
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.69	3.29 ng	132637	1,4-Dichlorobenzene-d4	5.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2		Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3		Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
4		Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5		Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040317\
Data File : BF094135.D
Acq On : 3 Apr 2017 21:50
Operator : SJ/MA
Sample : I2493-10 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

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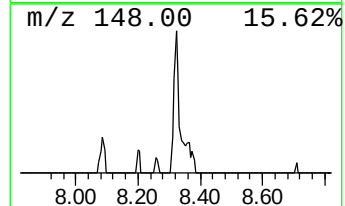
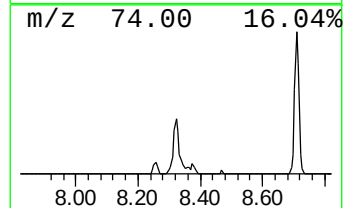
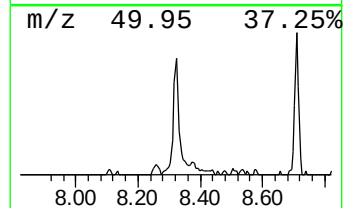
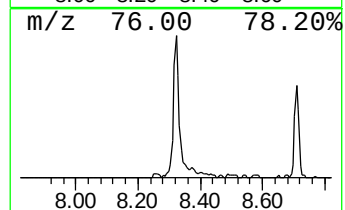
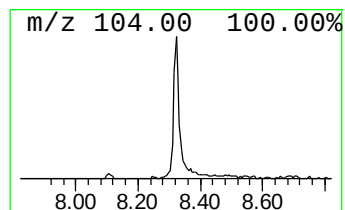
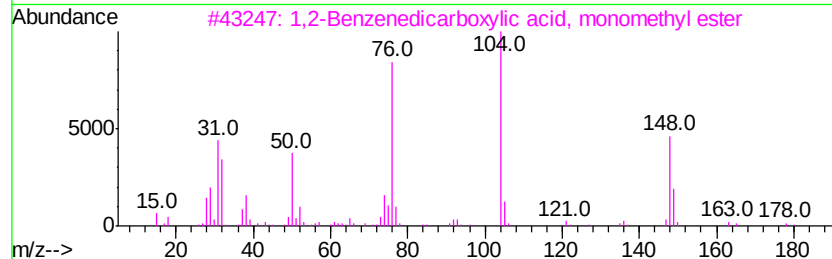
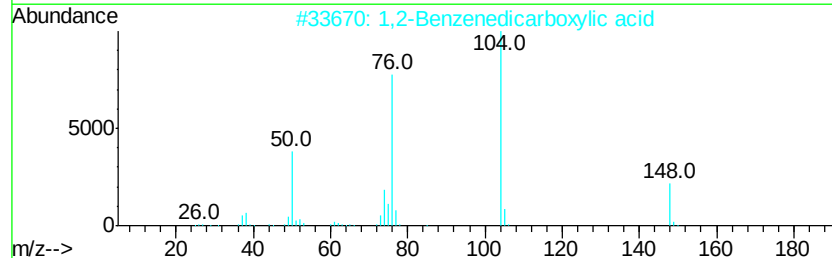
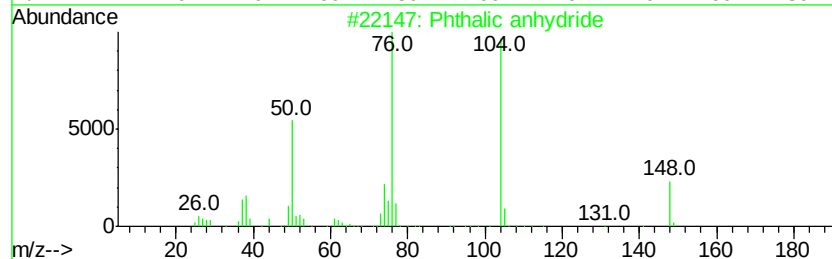
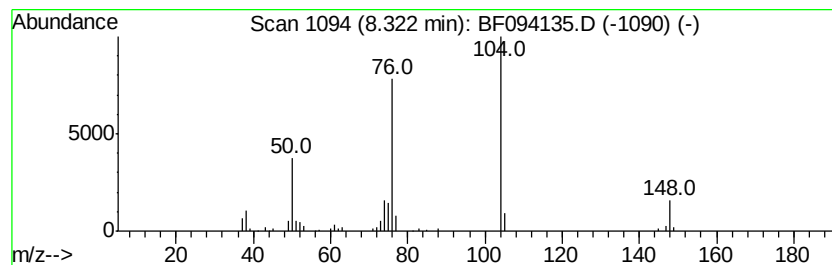
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Phthalic anhydride Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.32	2.17 ng	119890	Naphthalene-d8	7.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phthalic anhydride	148	C8H4O3	000085-44-9	95
2		1,2-Benzenedicarboxylic acid	166	C8H6O4	000088-99-3	90
3		1,2-Benzenedicarboxylic acid, mo...	180	C9H8O4	004376-18-5	83
4		Phthamic acid	165	C8H7NO3	000088-97-1	83
5		Phthalic acid, monoethyl ester	194	C10H10O4	002306-33-4	83



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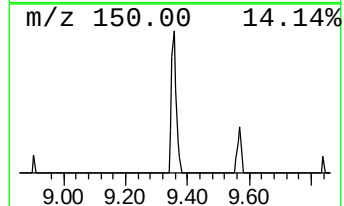
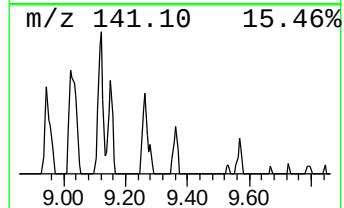
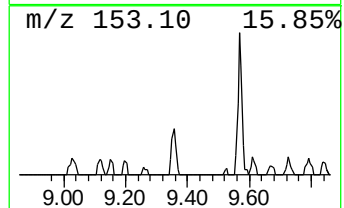
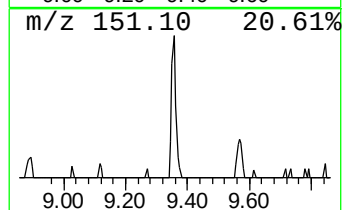
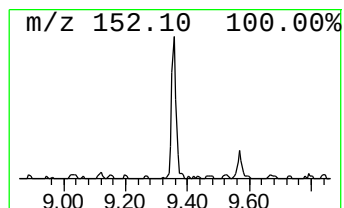
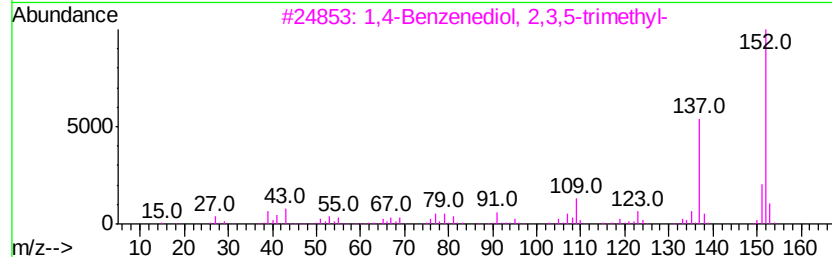
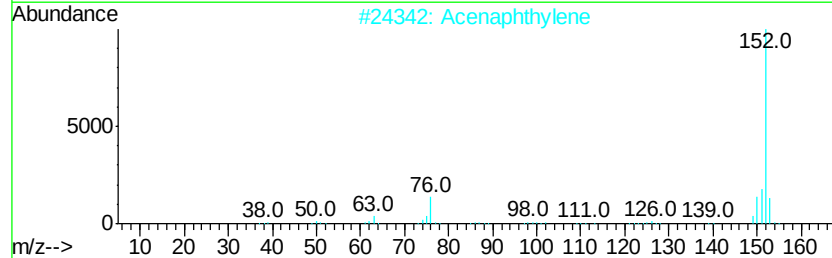
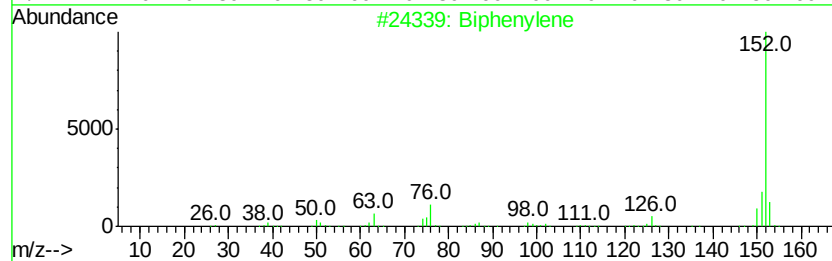
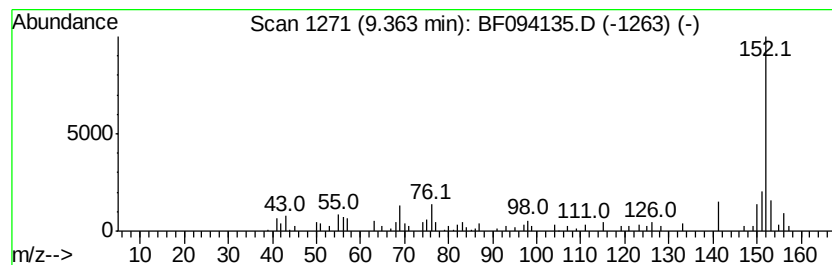
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Biphenylene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.36	2.19 ng	116709	Acenaphthene-d10	9.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Biphenylene	152	C12H8	000259-79-0	87
2		Acenaphthylene	152	C12H8	000208-96-8	72
3		1,4-Benzenediol, 2,3,5-trimethyl-	152	C9H12O2	000700-13-0	53
4		6,7,8,9-Tetrahydro-5H-[1,2,4]tri...	152	C7H12N4	052333-20-7	38
5		Tisopurine	152	C5H4N4S	005334-23-6	33



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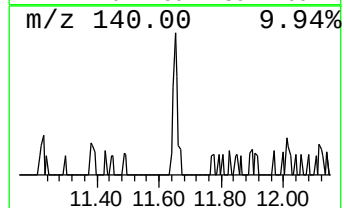
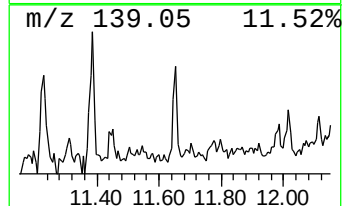
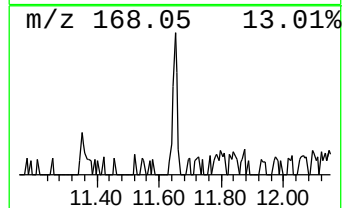
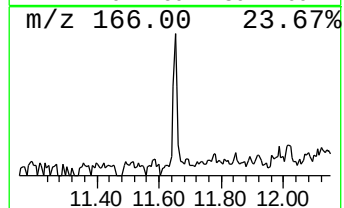
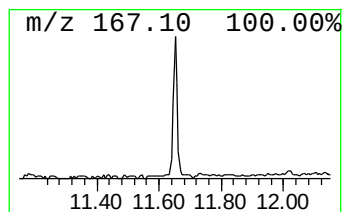
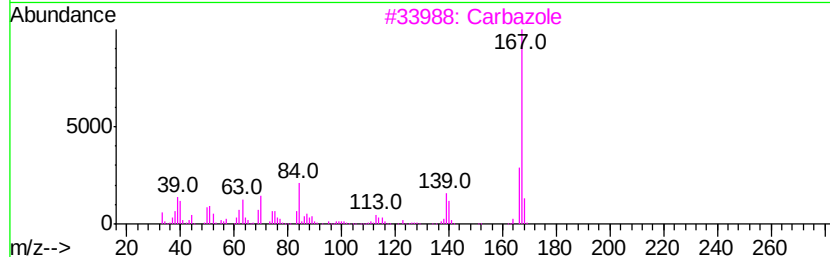
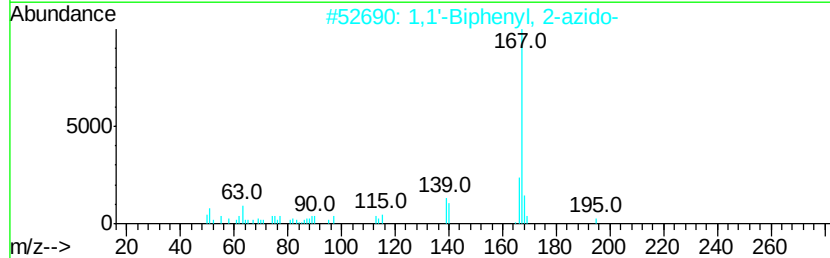
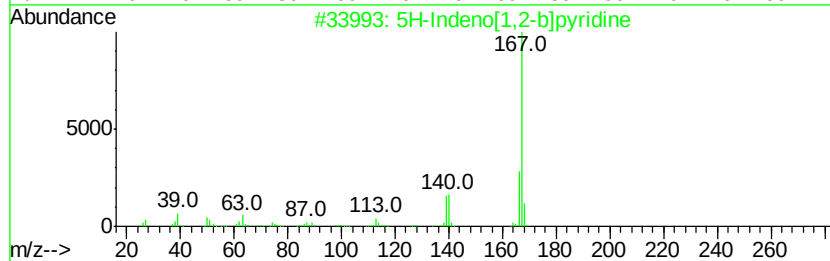
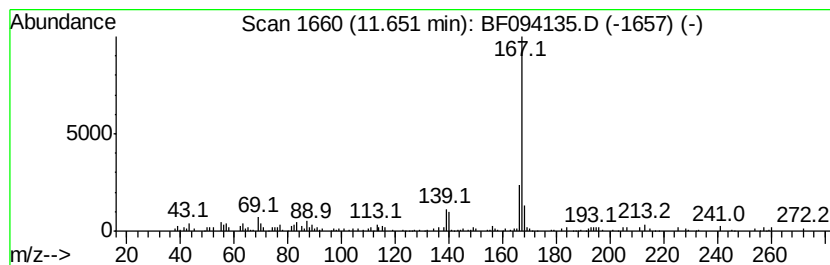
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 5H-Indeno[1,2-b]pyridine Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.65	2.02 ng	82088	Phenanthrene-d10	11.36	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	90
2	1,1'-Biphenyl, 2-azido-	195	C12H9N3	007599-23-7	87
3	Carbazole	167	C12H9N	000086-74-8	86
4	9H-Carbazole, 9-nitroso-	196	C12H8N2O	002788-23-0	86
5	D-Galactitol, 2-(acetylmethylami...	363	C16H29N08	074410-42-7	64



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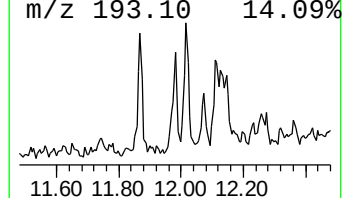
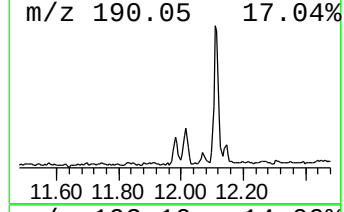
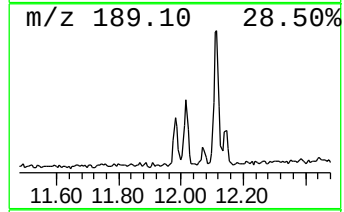
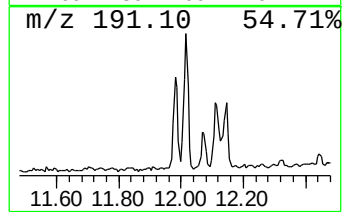
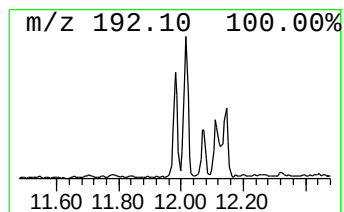
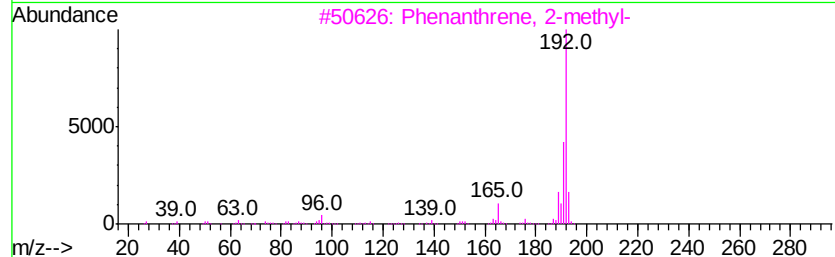
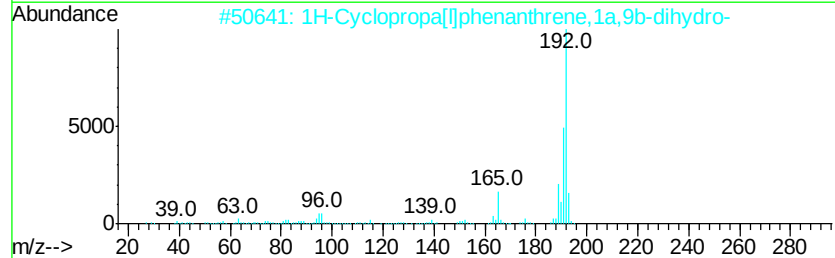
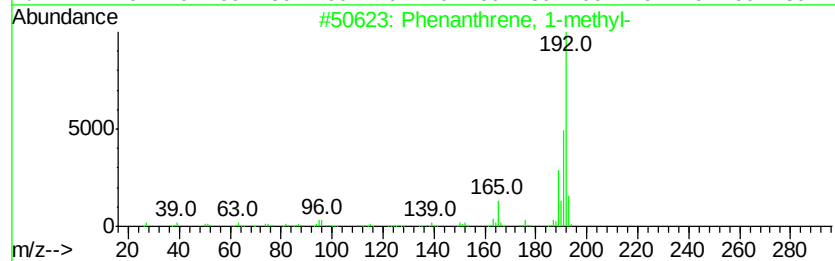
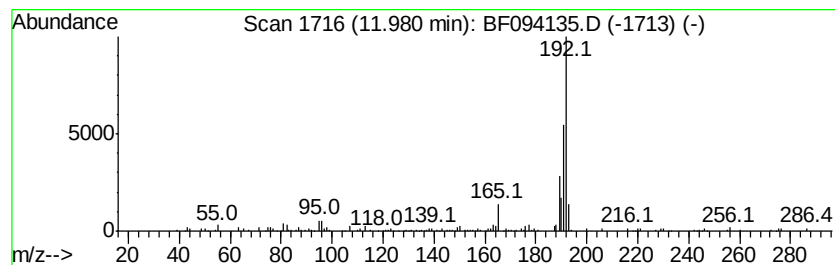
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.98	2.35 ng	95444	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
2		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	81
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	81
4		1H-Indene, 3-phenyl-	192	C15H12	001961-97-3	81
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	81



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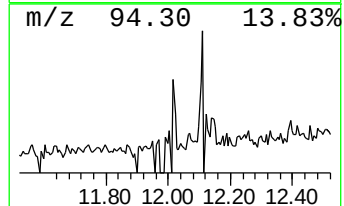
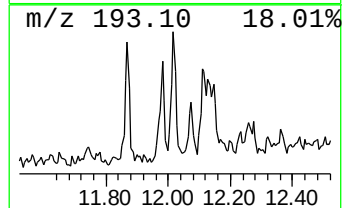
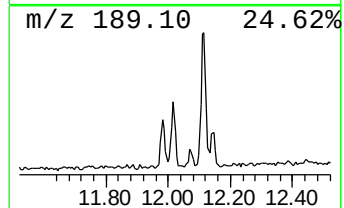
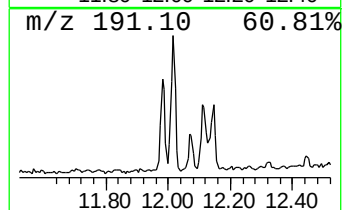
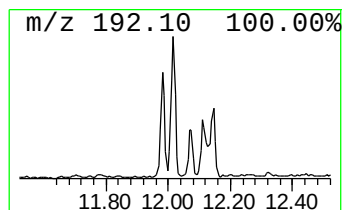
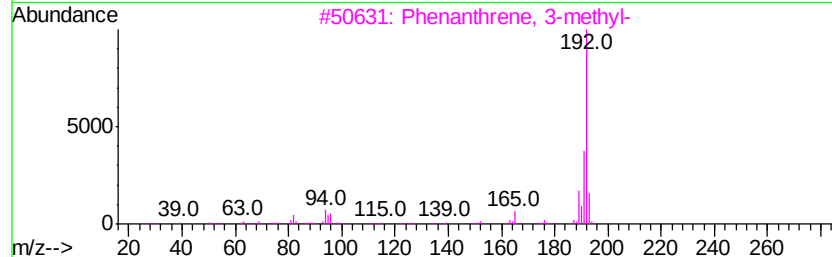
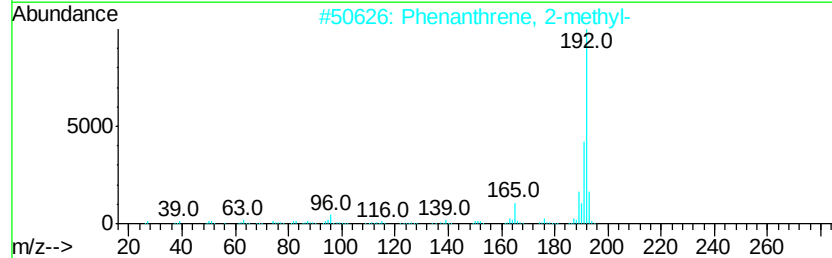
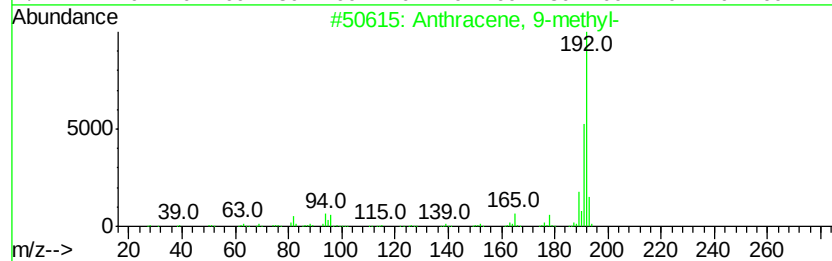
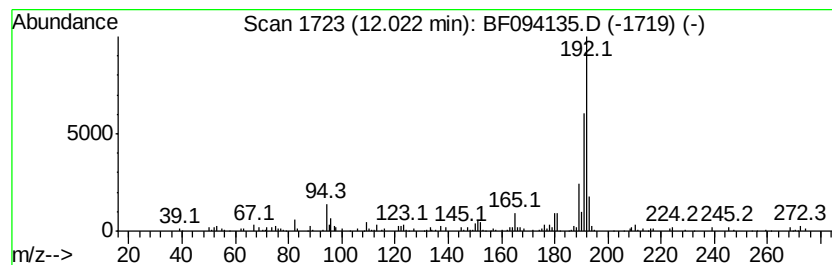
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Anthracene, 9-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.02	3.33 ng	135248	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3		Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	95
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
5		1H-Cyclopropa[l]phenanthrene, 1a, ...	192	C15H12	000949-41-7	93



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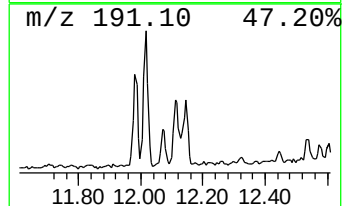
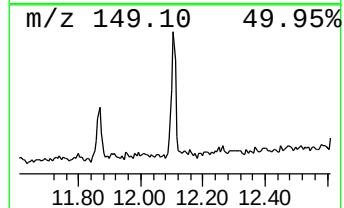
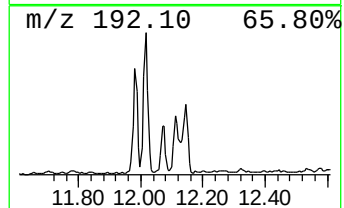
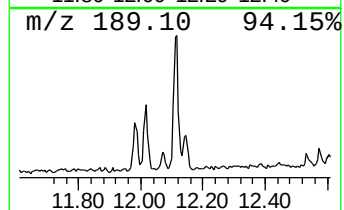
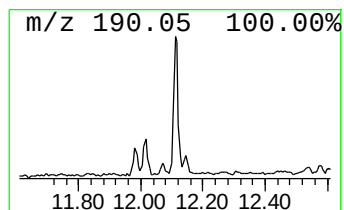
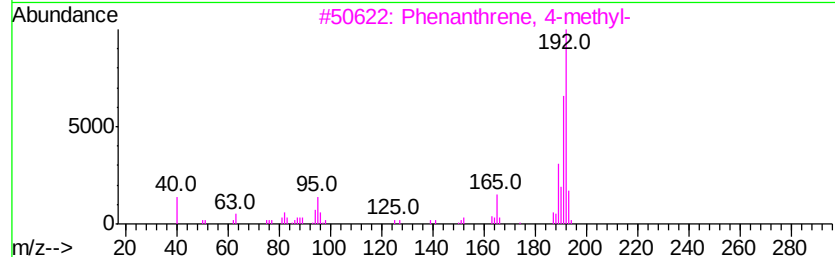
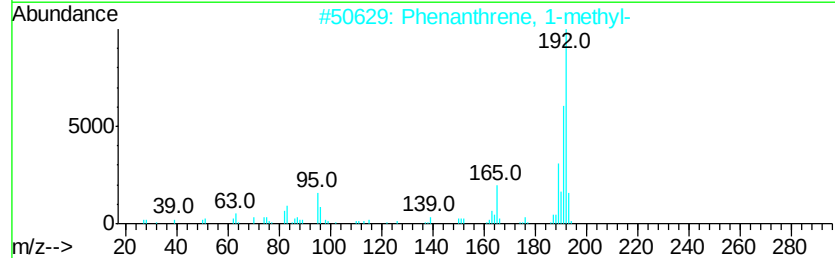
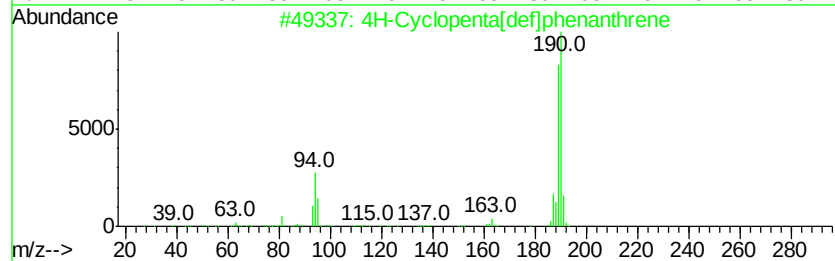
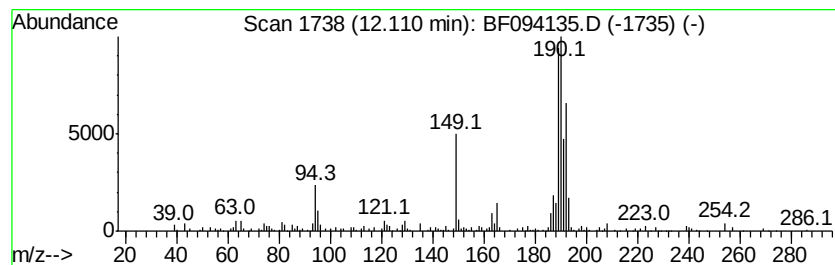
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Peak Number 12 unknown12.11 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.11	4.16 ng	168967	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	49
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	25
3		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	25
4		8H-Imidazo[4,5-E]-2,1,3-benzothi...	190	C8H6N4S	129485-68-3	22
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	15



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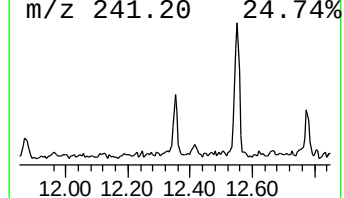
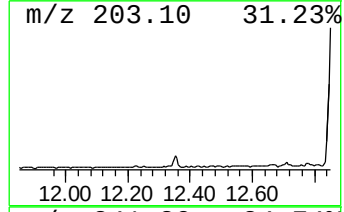
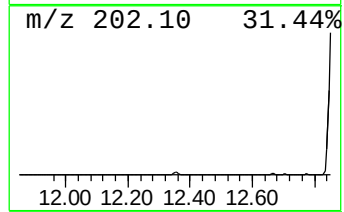
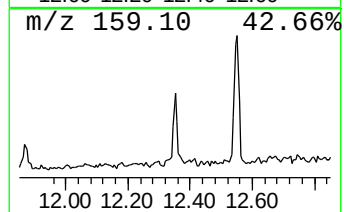
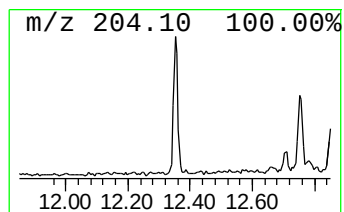
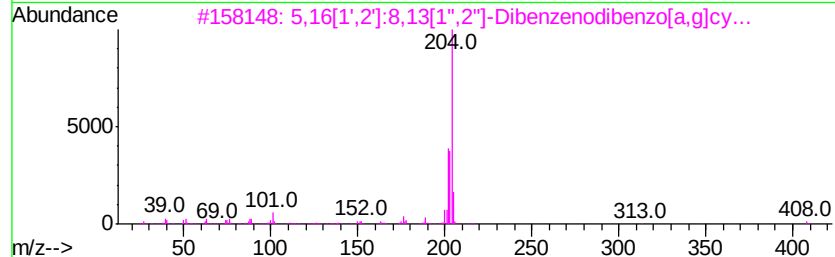
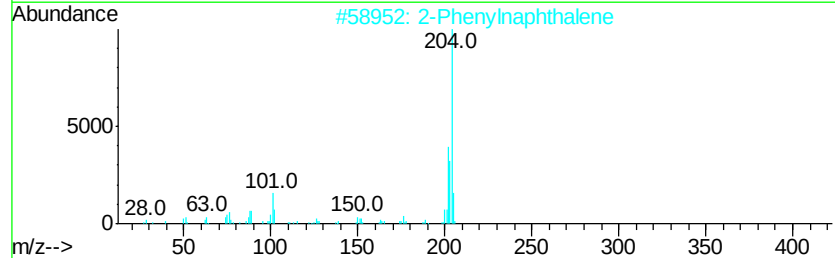
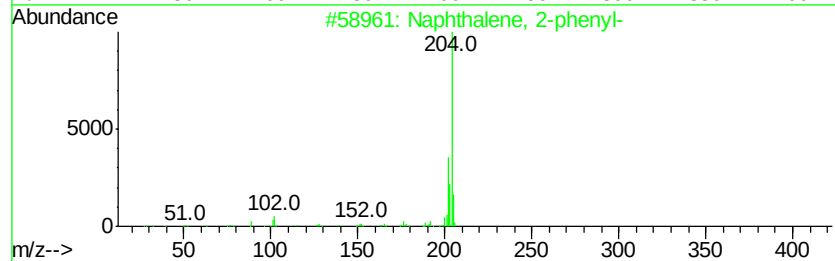
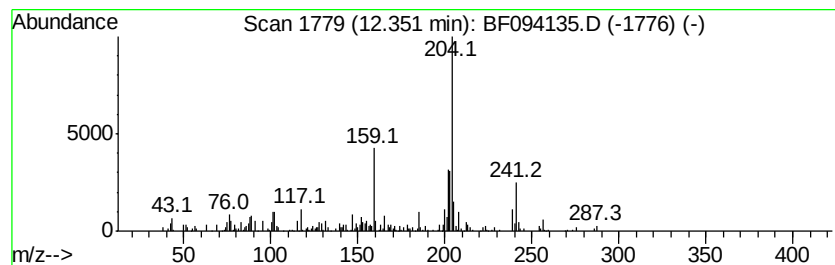
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Naphthalene, 2-phenyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.35	4.27 ng	173611	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	64
2		2-Phenylnaphthalene	204	C16H12	035465-71-5	50
3		5,16[1',2']8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	49
4		Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	42
5		9-Acridinecarbonitrile	204	C14H8N2	005326-19-2	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF040317\
Data File : BF094135.D
Acq On : 3 Apr 2017 21:50
Operator : SJ/MA
Sample : I2493-10 5X
Misc :
ALS Vial : 19 Sample Multiplier: 1

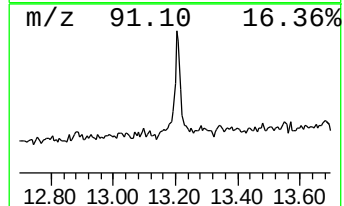
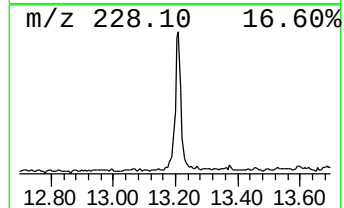
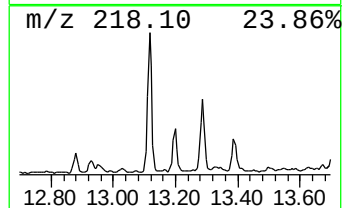
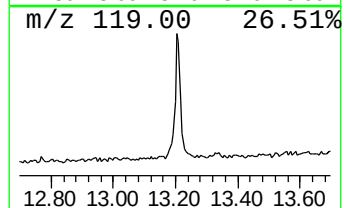
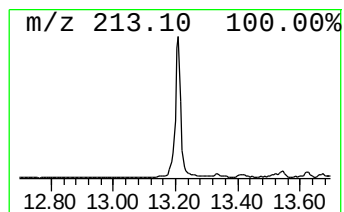
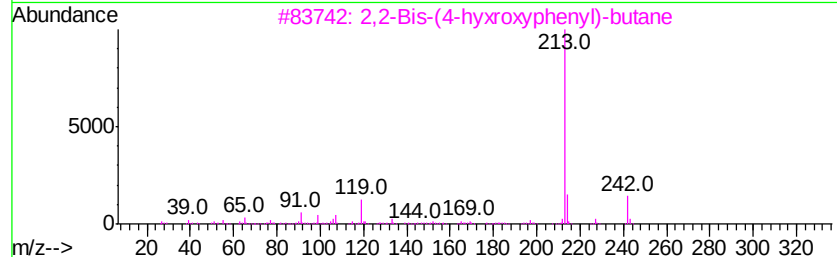
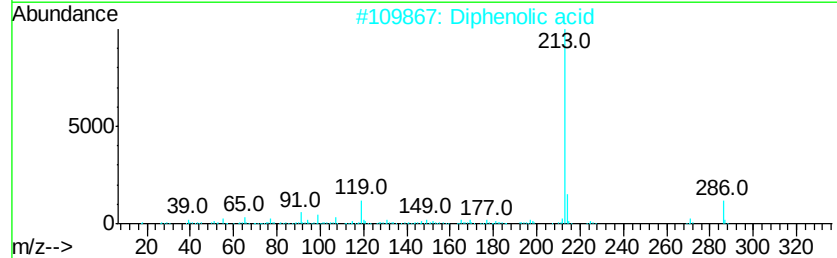
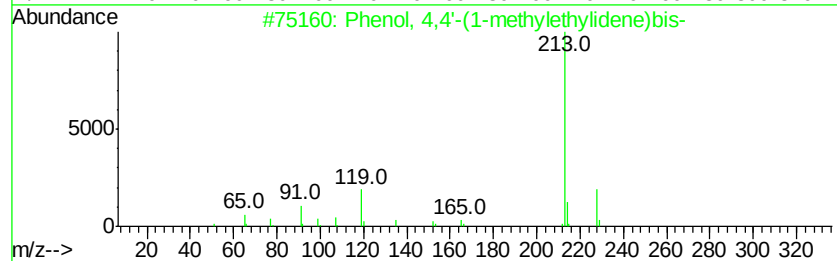
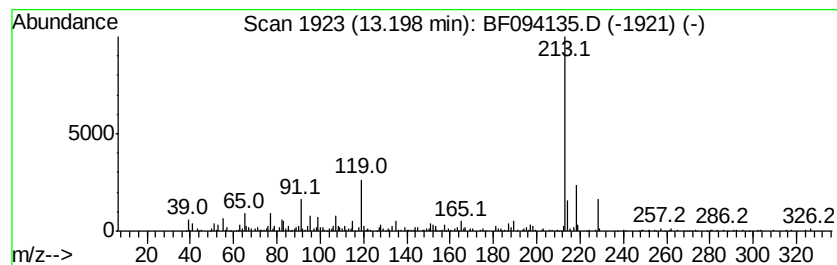
Instrument :
BNA_F
ClientSampleId :
LSS-SP08-COMP

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Phenol, 4,4'-(1-methylethyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.20	8.79 ng	506395	Chrysene-d12	14.62	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	96
2	Diphenolic acid	286	C17H18O4	000126-00-1	59
3	2,2-Bis-(4-hydroxyphenyl)-butane	242	C16H18O2	000077-40-7	59
4	2H,8H-Benzo[1,2-b:5,4-b']dipyran...	228	C14H12O3	000553-19-5	58
5	2-Bromo-7-(methylamino)tropone	213	C8H8BrNO	103539-25-9	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040317\
 Data File : BF094135.D
 Acq On : 3 Apr 2017 21:50
 Operator : SJ/MA
 Sample : I2493-10 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LSS-SP08-COMP

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Butenoic acid, ...	2.27	2.5	ng	99315	1	5.89	806717	20.0
2-Pentanone, 4-hy...	4.01	2.0	ng	82850	1	5.89	806717	20.0
unknown5.63	5.63	25.6	ng	1030720	1	5.89	806717	20.0
Benzene, 1,2,3-tr...	5.69	3.3	ng	132637	1	5.89	806717	20.0
Phthalic anhydride	8.32	2.2	ng	119890	2	7.39	1104750	20.0
Biphenylene	9.36	2.2	ng	116709	3	9.53	1067570	20.0
5H-Indeno[1,2-b]p...	11.65	2.0	ng	82088	4	11.36	812524	20.0
Phenanthrene, 1-m...	11.98	2.4	ng	95444	4	11.36	812524	20.0
Anthracene, 9-met...	12.02	3.3	ng	135248	4	11.36	812524	20.0
unknown12.11	12.11	4.2	ng	168967	4	11.36	812524	20.0
Naphthalene, 2-ph...	12.35	4.3	ng	173611	4	11.36	812524	20.0
Phenol, 4,4'-(1-m...	13.20	8.8	ng	506395	5	14.62	1151820	20.0