

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078511.D
 Acq On : 14 Apr 2015 20:57
 Operator : TP/IZ
 Sample : G1828-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB02B

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	1.495	26	27	34	rVB	28985	54927	1.76%	0.391%
2	1.598	34	36	72	rVB	1199738	3116205	100.00%	22.173%
3	4.444	283	285	291	rBV3	45179	72843	2.34%	0.518%
4	5.061	336	339	343	rBV	55901	83580	2.68%	0.595%
5	6.410	456	457	462	rBV2	97851	148047	4.75%	1.053%
6	6.524	465	467	470	rVV	127119	129689	4.16%	0.923%
7	6.799	489	491	494	rBV	94611	107731	3.46%	0.767%
8	6.993	505	508	510	rBV	72112	95925	3.08%	0.683%
9	7.507	551	553	555	rBV	110544	103206	3.31%	0.734%
10	8.387	627	630	631	rBV	115519	160782	5.16%	1.144%
11	9.599	730	736	738	rBV3	18043	48406	1.55%	0.344%
12	9.725	746	747	750	rVB	167430	232604	7.46%	1.655%
13	9.805	750	754	756	rBV	70156	90534	2.91%	0.644%
14	10.182	784	787	788	rBV	57072	58351	1.87%	0.415%
15	10.250	791	793	797	rBV4	18778	64748	2.08%	0.461%
16	10.330	798	800	802	rVB	46436	40840	1.31%	0.291%
17	10.365	802	803	806	rBV3	43672	91059	2.92%	0.648%
18	10.410	806	807	811	rVB2	78298	137623	4.42%	0.979%
19	10.479	811	813	815	rBV3	25619	48668	1.56%	0.346%
20	10.536	815	818	821	rVB2	178745	286769	9.20%	2.041%
21	10.593	821	823	826	rBV	60412	74119	2.38%	0.527%
22	10.673	828	830	832	rBV2	23784	42281	1.36%	0.301%
23	10.856	844	846	847	rBV	36174	42802	1.37%	0.305%
24	10.959	852	855	857	rVV3	94377	130918	4.20%	0.932%
25	11.005	857	859	861	rBV2	31460	45696	1.47%	0.325%
26	11.039	861	862	864	rBV	47650	58780	1.89%	0.418%
27	11.108	865	868	869	rBV	54425	75584	2.43%	0.538%
28	11.519	901	904	906	rBV	164425	230975	7.41%	1.644%
29	12.365	976	978	979	rBV	210930	183427	5.89%	1.305%
30	13.862	1107	1109	1111	rBV	216327	218948	7.03%	1.558%
31	14.137	1131	1133	1135	rVB	238747	290911	9.34%	2.070%
32	14.239	1140	1142	1144	rBV	382442	404732	12.99%	2.880%
33	14.354	1150	1152	1155	rVB	740715	888492	28.51%	6.322%
34	15.371	1237	1241	1242	rVV2	39057	74731	2.40%	0.532%

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
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35	15.405	1242	1244	1246	rVB2	33857	51086	1.64%	0.364%
36	15.542	1254	1256	1259	rBV4	28991	52661	1.69%	0.375%
37	15.622	1260	1263	1265	rBV2	352170	630362	20.23%	4.485%
38	15.748	1272	1274	1276	rBV	88148	119233	3.83%	0.848%
39	15.885	1284	1286	1288	rVB	47414	54114	1.74%	0.385%
40	15.931	1288	1290	1295	rBV2	33727	67948	2.18%	0.483%
41	16.125	1305	1307	1310	rVB	51956	72491	2.33%	0.516%
42	16.925	1374	1377	1382	rVB	377593	817411	26.23%	5.816%
43	17.040	1385	1387	1389	rVB	120829	131375	4.22%	0.935%
44	17.085	1389	1391	1394	rBV4	18116	34183	1.10%	0.243%
45	17.165	1395	1398	1399	rBV	56428	94716	3.04%	0.674%
46	17.245	1402	1405	1408	rVB	276345	346247	11.11%	2.464%
47	17.303	1408	1410	1413	rVB	408800	549315	17.63%	3.909%
48	17.371	1413	1416	1417	rBV	247663	307543	9.87%	2.188%
49	17.394	1417	1418	1420	rVB	100829	134730	4.32%	0.959%
50	17.554	1429	1432	1439	rBV5	36971	147818	4.74%	1.052%
51	17.783	1450	1452	1454	rBV	26017	48984	1.57%	0.349%
52	17.874	1458	1460	1463	rVB2	52258	84921	2.73%	0.604%
53	18.480	1510	1513	1517	rBV2	71934	159163	5.11%	1.133%
54	18.617	1522	1525	1529	rVV2	313539	574478	18.44%	4.088%
55	18.686	1529	1531	1535	rVB2	80859	153589	4.93%	1.093%
56	18.766	1536	1538	1540	rBV	62513	118963	3.82%	0.846%
57	18.811	1540	1542	1551	rVB4	73824	234295	7.52%	1.667%
58	18.971	1552	1556	1560	rBV	295563	526843	16.91%	3.749%
59	19.154	1569	1572	1574	rVB	84640	165955	5.33%	1.181%
60	20.731	1701	1710	1715	rBV2	82823	342648	11.00%	2.438%
61	20.880	1719	1723	1726	rBV	39474	97742	3.14%	0.695%

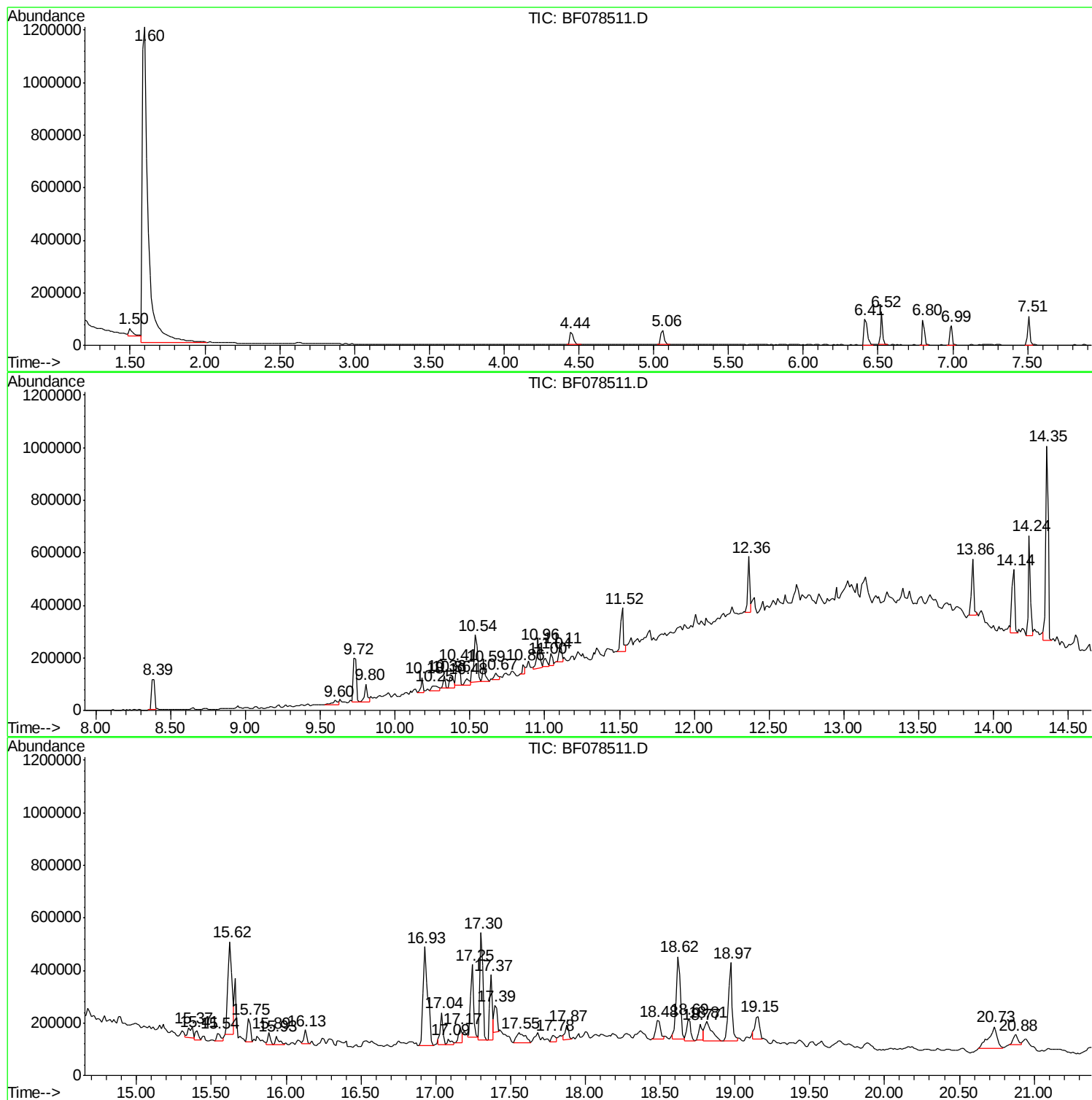
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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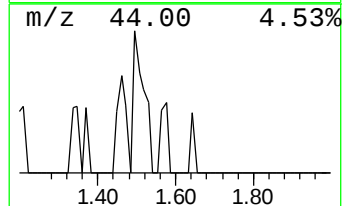
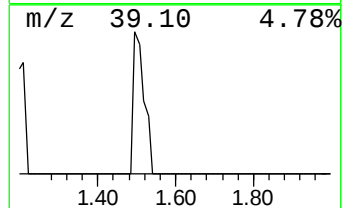
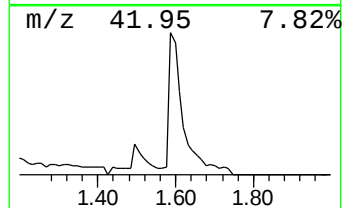
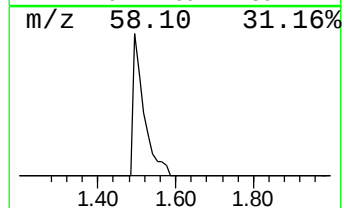
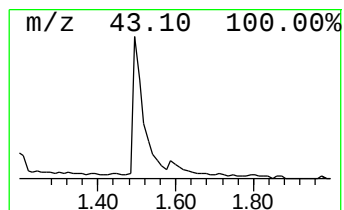
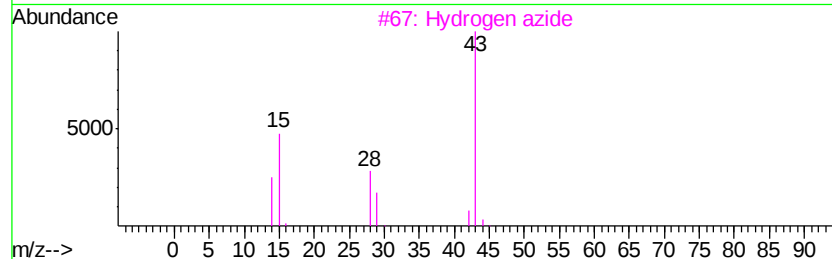
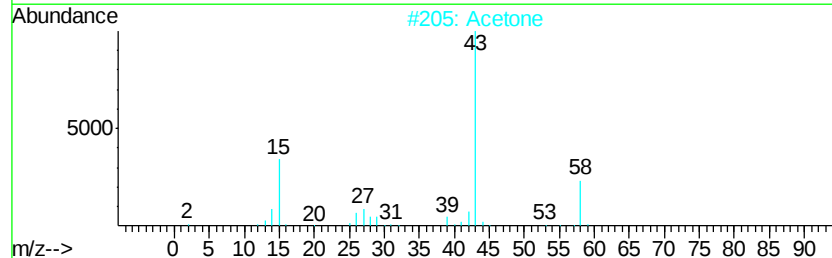
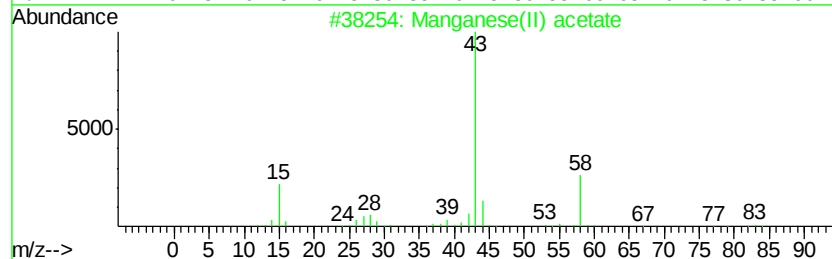
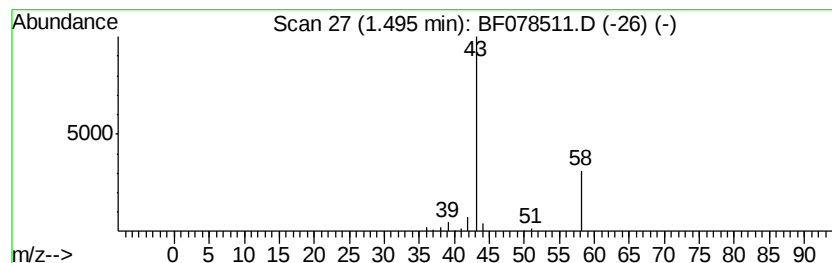
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown1.50 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.50	10.20 ng	54927	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Manganese(II) acetate	173	C4H6MnO4	006156-78-1	40
2		Acetone	58	C3H6O	000067-64-1	9
3		Hydrogen azide	43	HN3	007782-79-8	4
4		Ethene, methoxy-	58	C3H6O	000107-25-5	3
5		Diazene, dimethyl-	58	C2H6N2	000503-28-6	3



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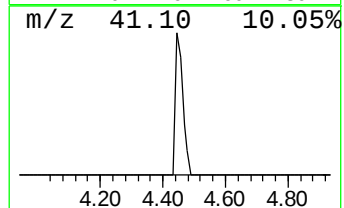
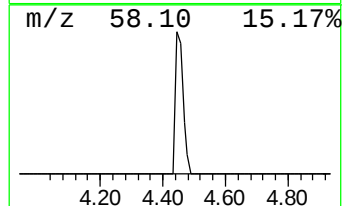
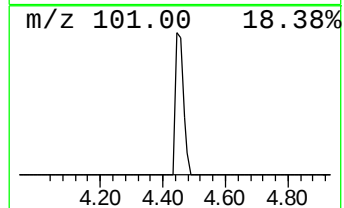
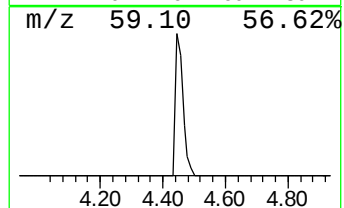
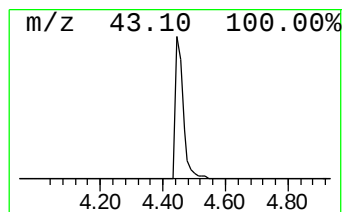
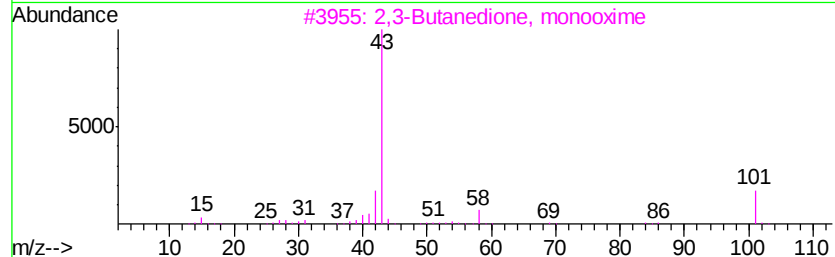
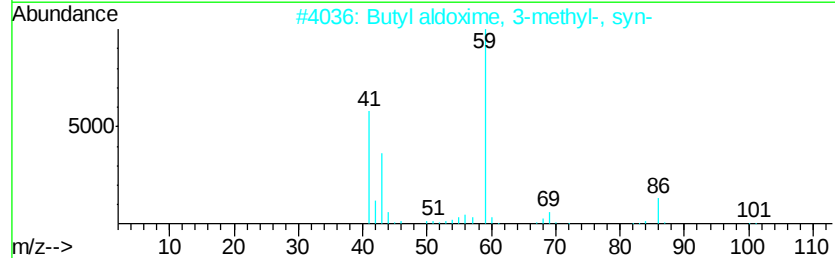
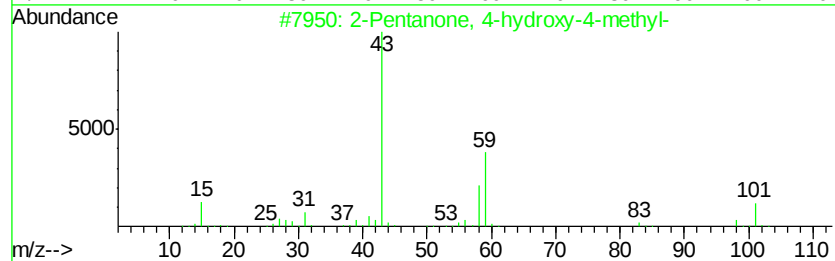
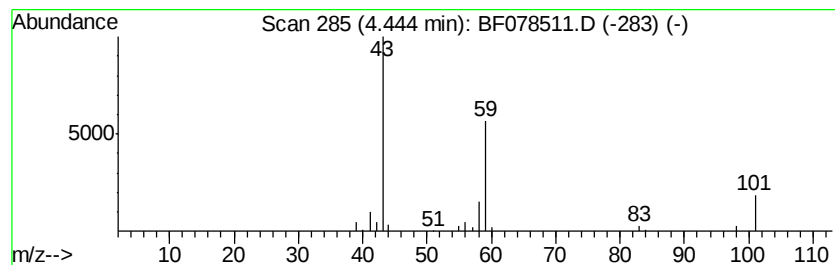
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.44	13.52 ng	72843	1,4-Dichlorobenzene-d4	6.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	37
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			3-Heptanol, 4-methyl-	130	C8H18O	014979-39-6	9



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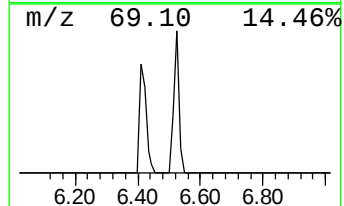
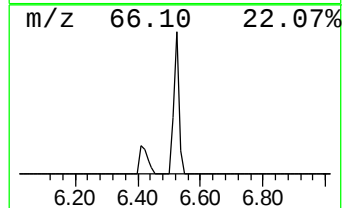
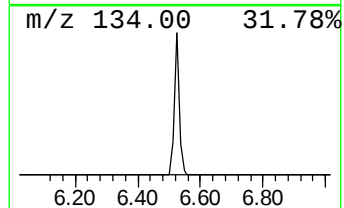
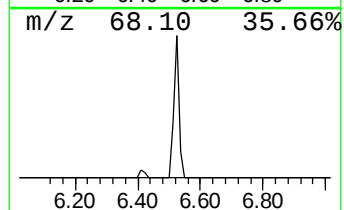
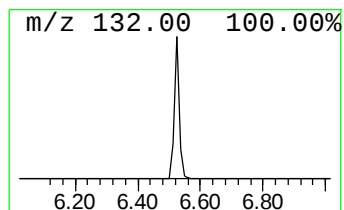
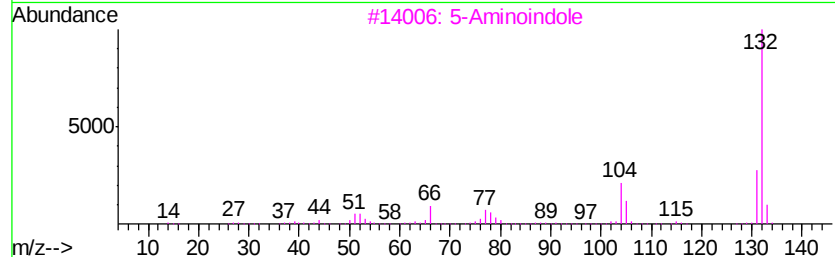
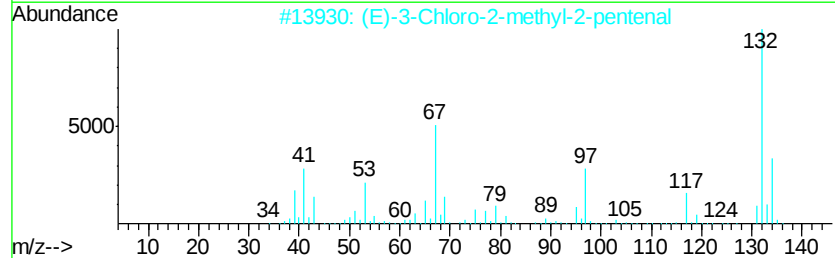
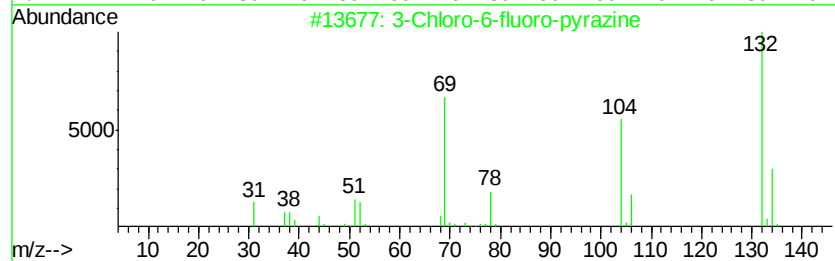
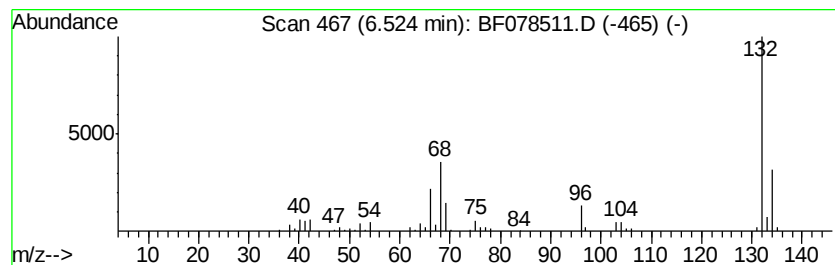
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown6.52 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.52	24.08 ng	129689	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Aminoindole	132	C8H8N2	005192-03-0	10
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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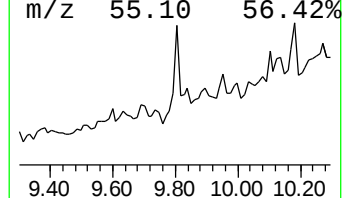
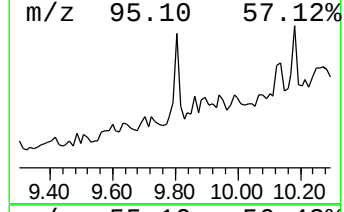
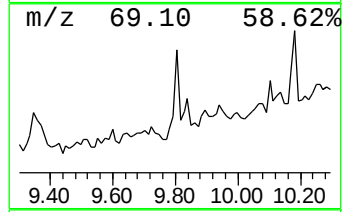
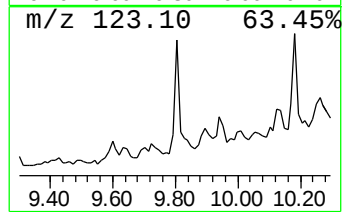
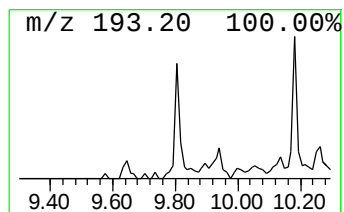
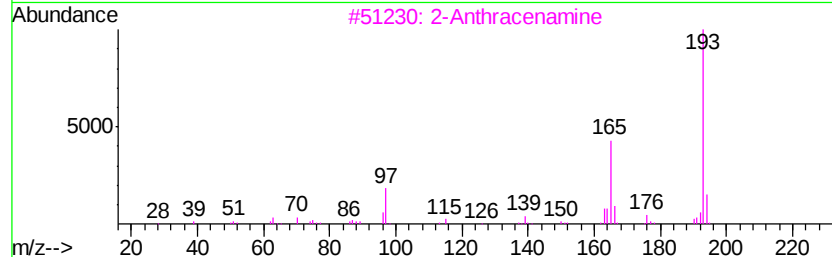
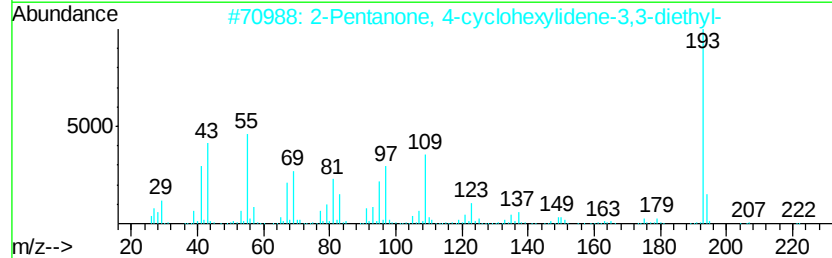
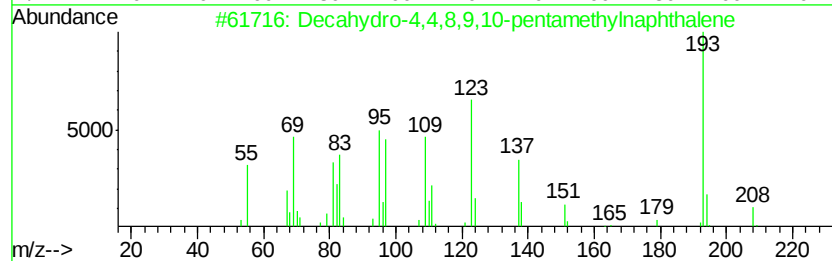
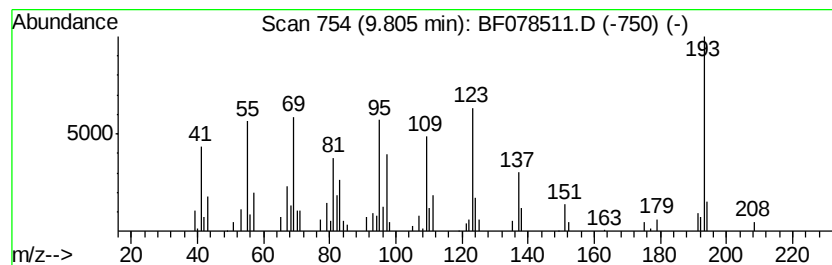
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 Decahydro-4,4,8,9,10-pentam... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.80	6.31 ng	90534	Acenaphthene-d10	10.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Decahydro-4,4,8,9,10-pentamethyl...	208	C15H28	080655-44-3	83
2		2-Pentanone, 4-cyclohexylidene-3...	222	C15H26O	313253-65-5	58
3		2-Anthracenamine	193	C14H11N	000613-13-8	35
4		Silane, tetra-2-propenyl-	192	C12H20Si	001112-66-9	27
5		2-Diethylamino-4-phenylthiooct-2...	302	C18H26N2S	1000090-57-0	27



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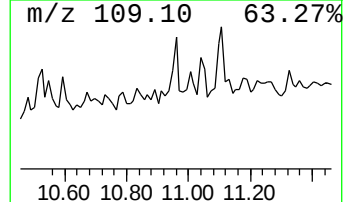
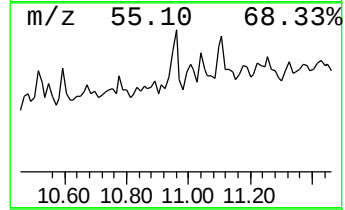
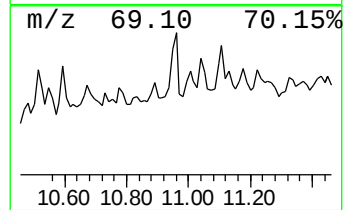
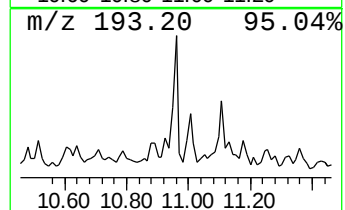
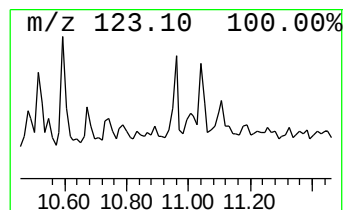
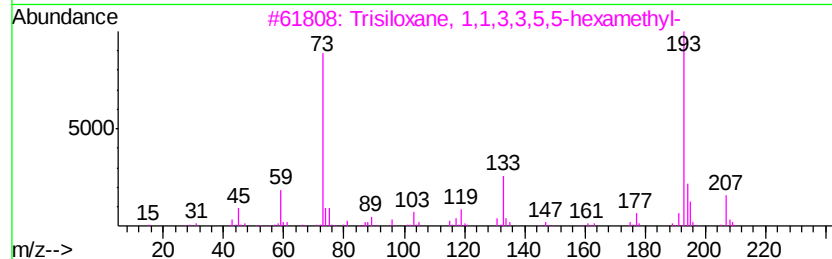
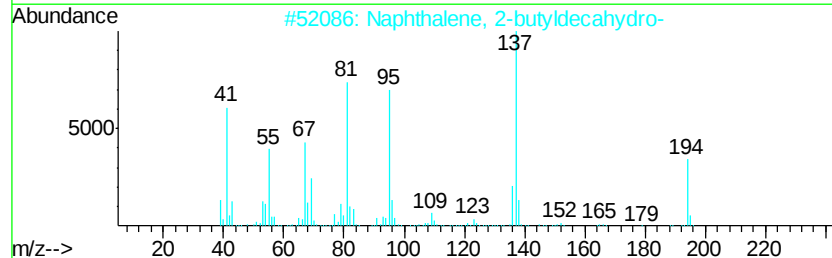
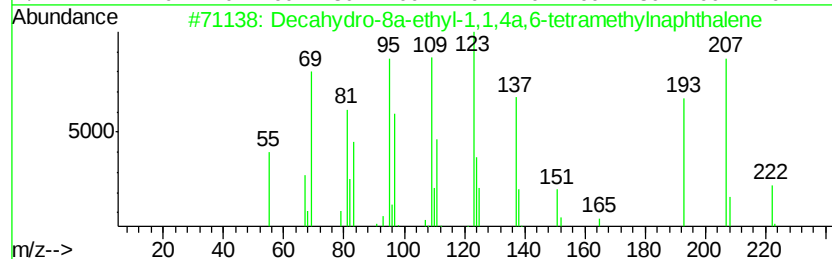
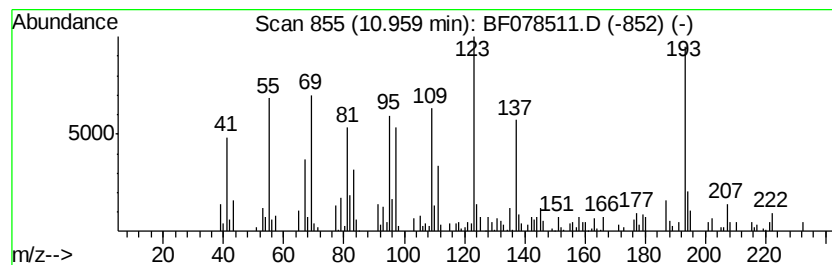
Instrument :
BNA_F
ClientSampleId :
HS-SB02B

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 7 unknown10.96 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
10.96	9.13 ng	130918	Acenaphthene-d10		10.54
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decahydro-8a-ethyl-1,1,4a,6-tetr...	222	C16H30	1000100-23-6	49
2	Naphthalene, 2-butyldecahydro-	194	C14H26	006305-52-8	35
3	Trisiloxane, 1,1,3,3,5,5-hexamet...	208	C6H20O2Si3	001189-93-1	25
4	2-(4-Fluorobenzoylamido)thiazole	222	C10H7FN2OS	313376-92-0	22
5	1-Trimethylsilylpent-1-en-4-yne	138	C8H14Si	079516-25-9	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078511.D
Acq On : 14 Apr 2015 20:57
Operator : TP/IZ
Sample : G1828-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

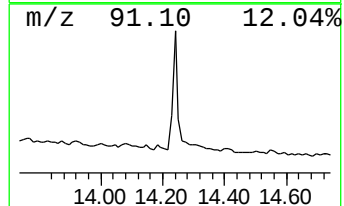
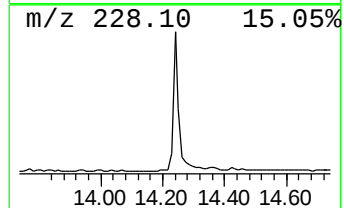
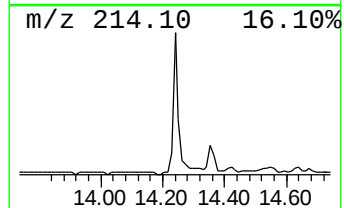
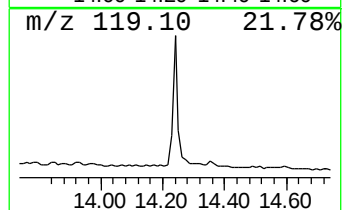
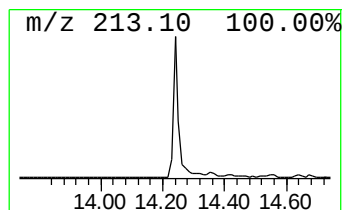
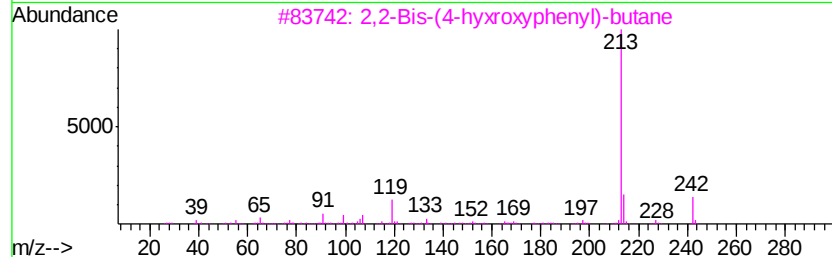
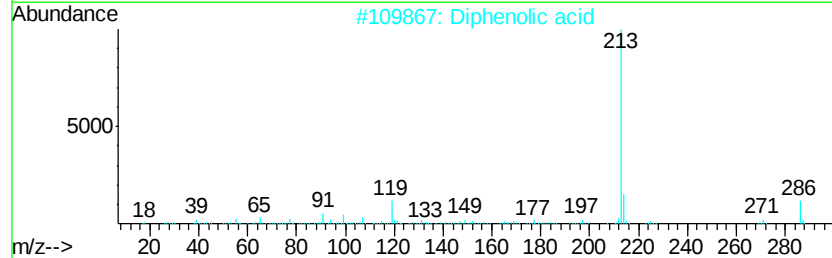
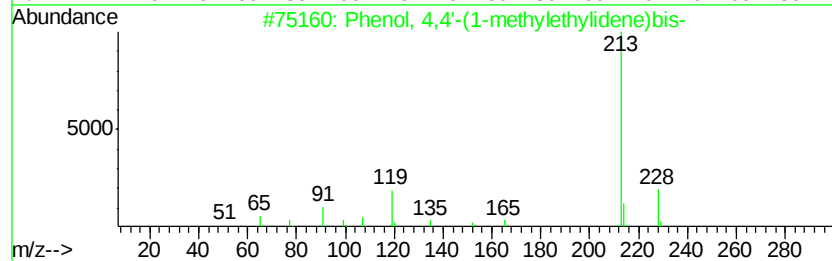
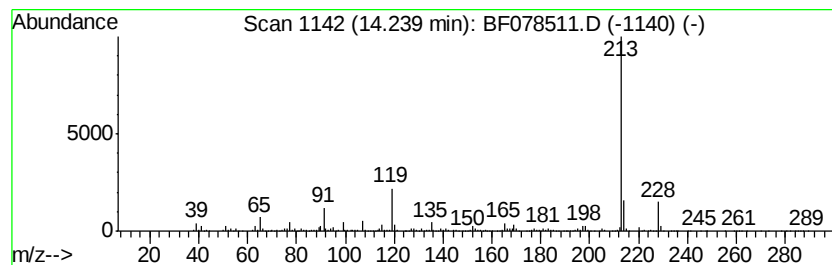
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Phenol, 4,4'-(1-methylethyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.		
14.24	12.84 ng	404732	Chrysene-d12	15.63		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97	
2	Diphenolic acid	286	C17H18O4	000126-00-1	64	
3	2,2-Bis-(4-hydroxyphenyl)-butane	242	C16H18O2	000077-40-7	59	
4	Carbanilic acid, p-phenyl-	213	C13H11NO2	004474-53-7	50	
5	As-Indacen-1(2H)-one, 3,6,7,8-te...	228	C16H20O	055591-18-9	43	



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078511.D
Acq On : 14 Apr 2015 20:57
Operator : TP/IZ
Sample : G1828-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

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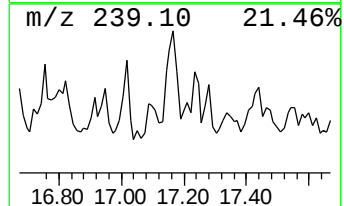
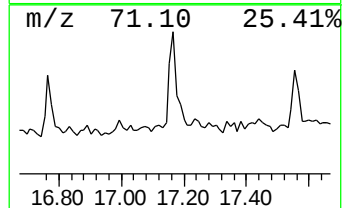
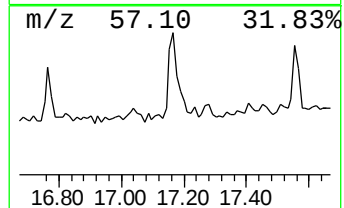
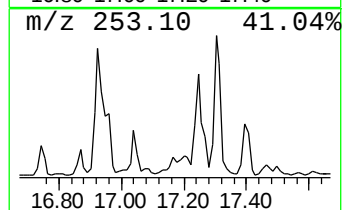
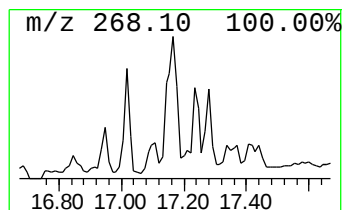
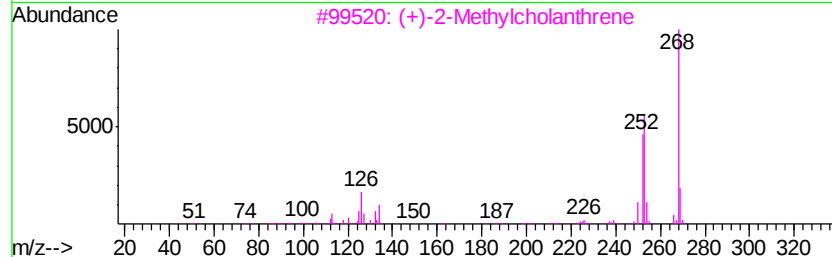
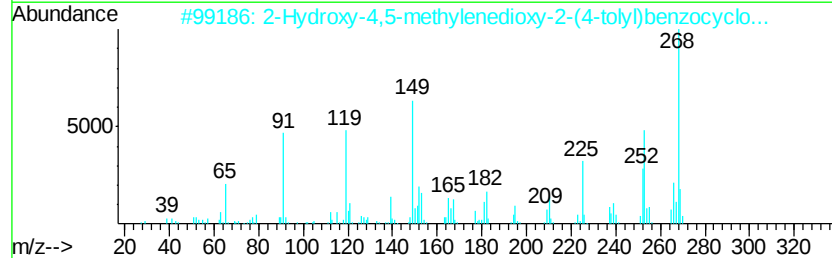
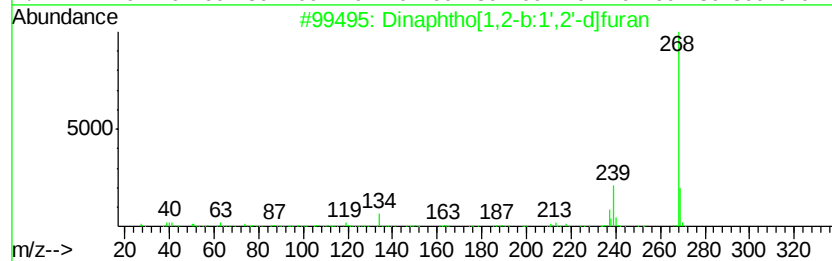
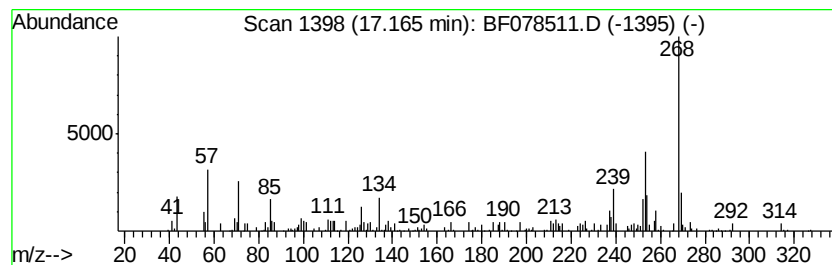
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Dinaphtho[1,2-b:1',2'-d]furan Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.17	6.16 ng	94716	Perylene-d12	17.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dinaphtho[1,2-b:1',2'-d]furan	268	C20H12O	000207-93-2	60
2			2-Hydroxy-4,5-methylenedioxy-2-(...	268	C16H12O4	1000284-51-7	59
3			(+)-2-Methylcholanthrene	268	C21H16	1000126-56-1	58
4			Dinaphtho[2,1-b:1',2'-d]furan	268	C20H12O	000194-63-8	49
5			Naphthalene, 1-(2-naphthalenyl)me...	268	C21H16	000611-48-3	49



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078511.D
Acq On : 14 Apr 2015 20:57
Operator : TP/IZ
Sample : G1828-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HS-SB02B

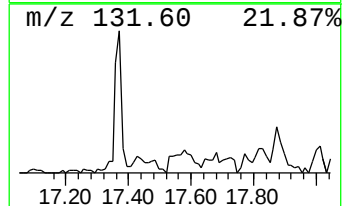
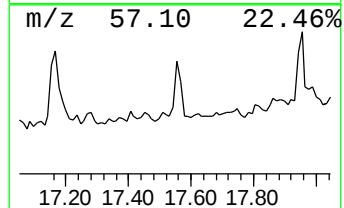
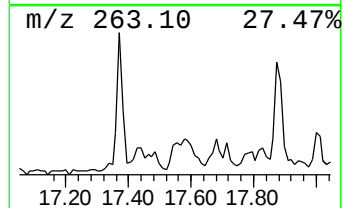
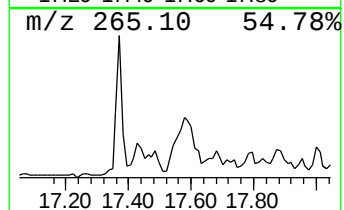
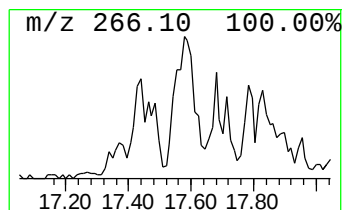
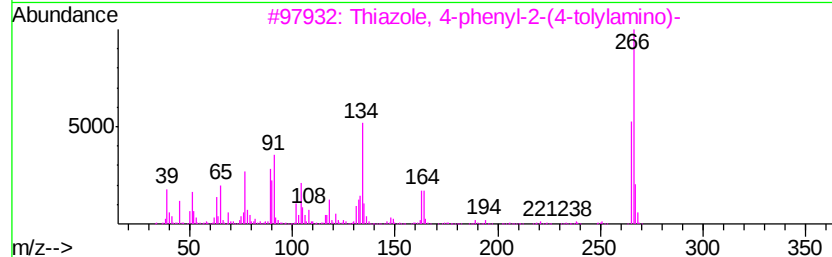
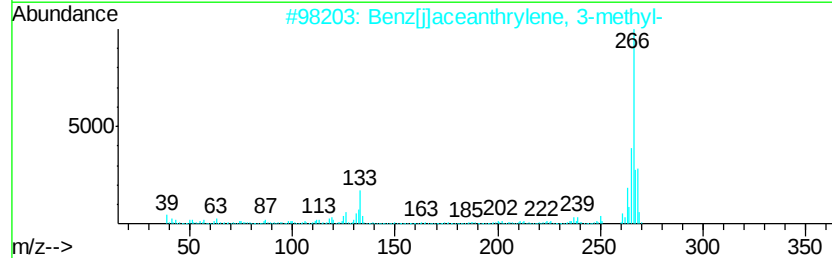
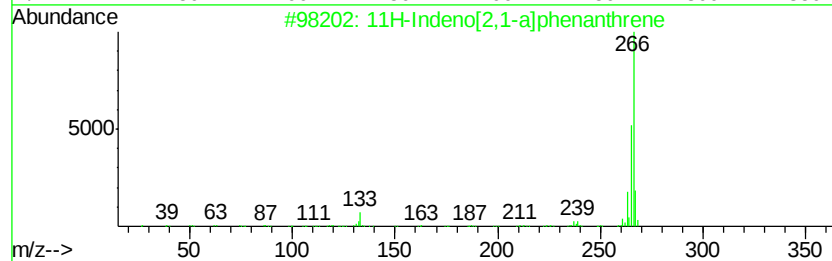
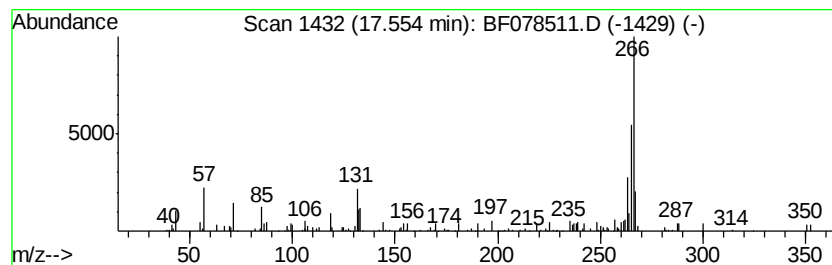
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 11H-Indeno[2,1-a]phenanthrene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.55	9.61 ng	147818	Perylene-d12	17.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Indeno[2,1-a]phenanthrene	266	C21H14	000220-97-3	62
2		Benz[j]aceanthrylene, 3-methyl-	266	C21H14	003343-10-0	60
3		Thiazole, 4-phenyl-2-(4-tolylami...	266	C16H14N2S	016098-04-7	58
4		13H-Dibenzo[a,h]fluorene	266	C21H14	000239-85-0	58
5		4H-1-Benzopyran-4-one, 3-hydroxy...	266	C17H14O3	078396-37-9	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078511.D
Acq On : 14 Apr 2015 20:57
Operator : TP/IZ
Sample : G1828-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HS-SB02B

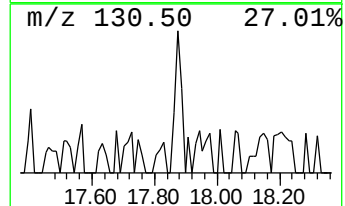
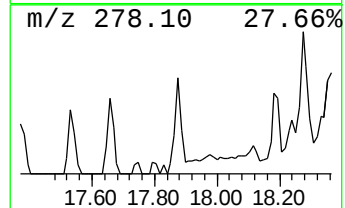
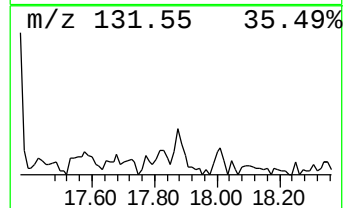
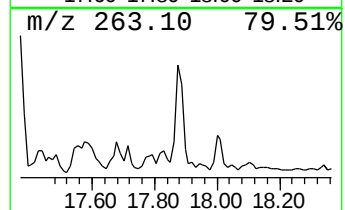
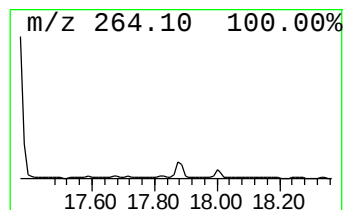
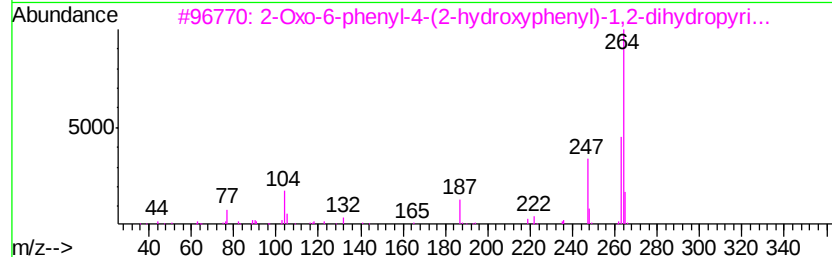
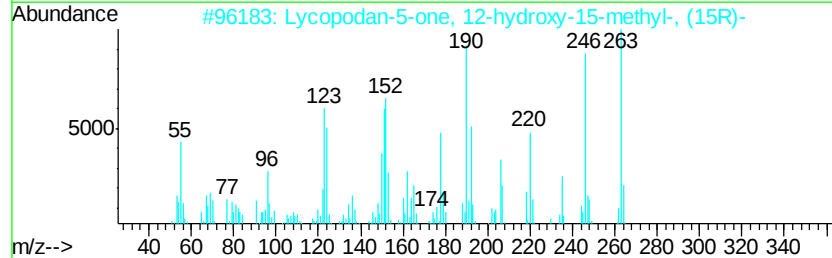
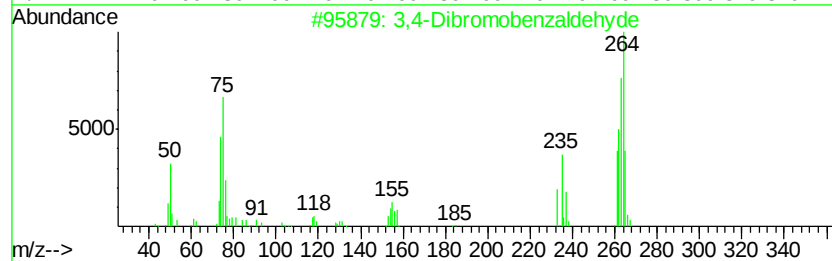
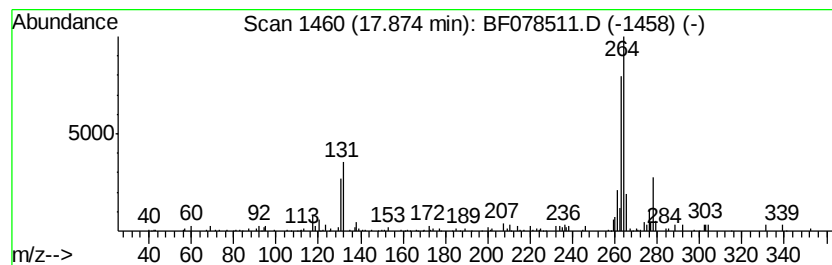
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 3,4-Dibromobenzaldehyde Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.87	5.52 ng	84921	Perylene-d12	17.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3,4-Dibromobenzaldehyde	262	C7H4Br2O	074003-55-7	58
2		Lycopodan-5-one, 12-hydroxy-15-m...	263	C16H25NO2	006900-92-1	38
3		2-Oxo-6-phenyl-4-(2-hydroxypheny...	264	C16H12N2O2	017432-82-5	32
4		6-Oxo-5-phenyl-2,3,5,6-tetrahydr...	263	C16H13N3O	087365-22-8	27
5		3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078511.D
Acq On : 14 Apr 2015 20:57
Operator : TP/IZ
Sample : G1828-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

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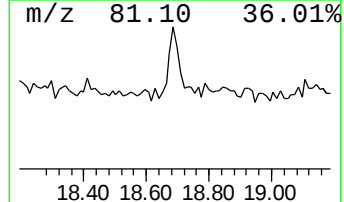
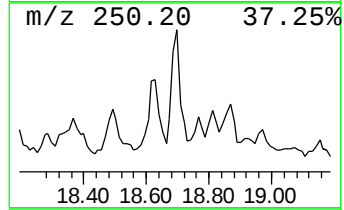
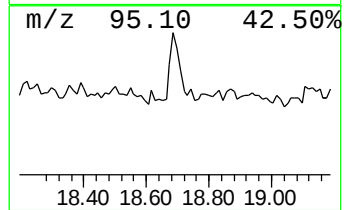
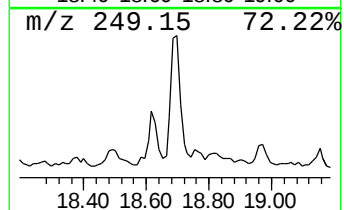
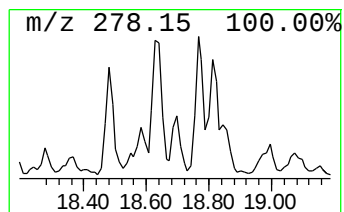
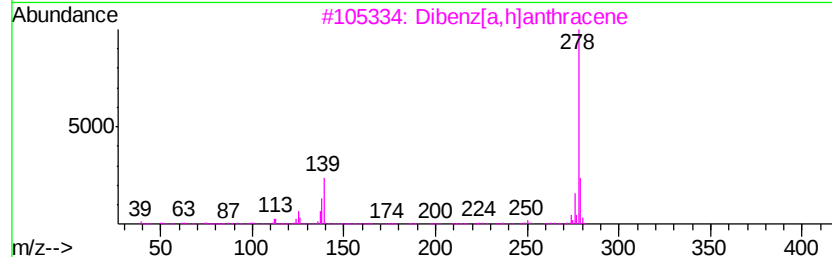
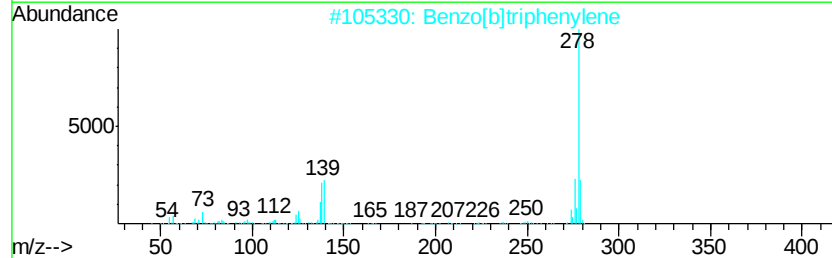
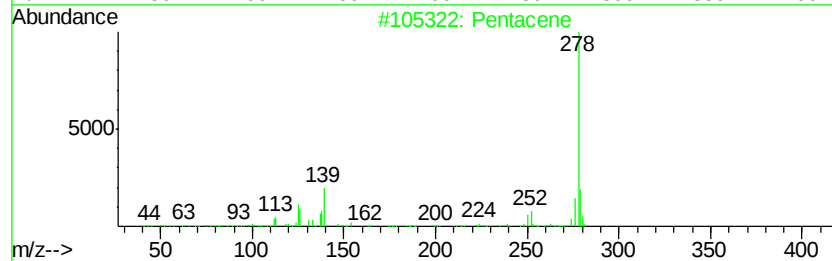
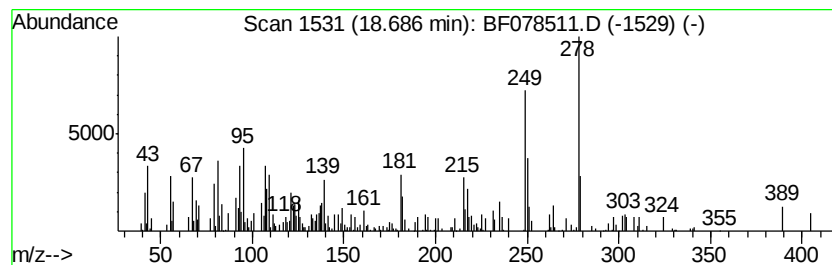
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown18.69 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.69	9.99 ng	153589	Perylene-d12	17.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentacene	278	C22H14	000135-48-8	30
2		Benzo[b]triphenylene	278	C22H14	000215-58-7	30
3		Dibenz[a,h]anthracene	278	C22H14	000053-70-3	30
4		3H-Pyrazol-3-one, 1,2-dihydro-5-...	278	C16H14N4O	053847-70-4	30
5		1,2:7,8-Dibenzophenanthrene	278	C22H14	000213-46-7	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078511.D
Acq On : 14 Apr 2015 20:57
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Misc :
ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
HS-SB02B

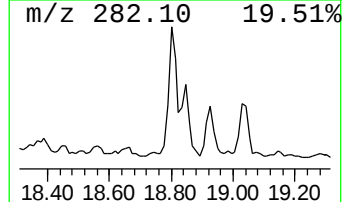
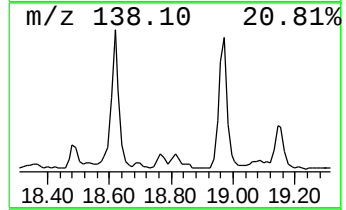
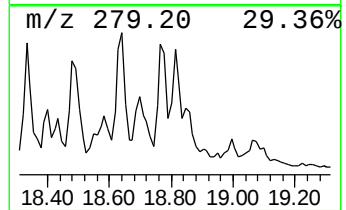
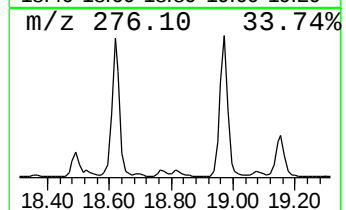
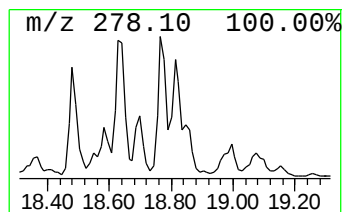
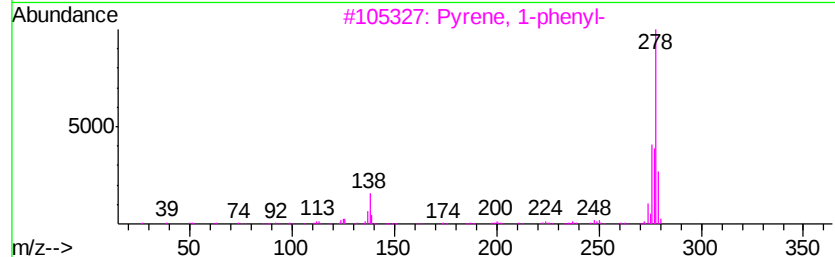
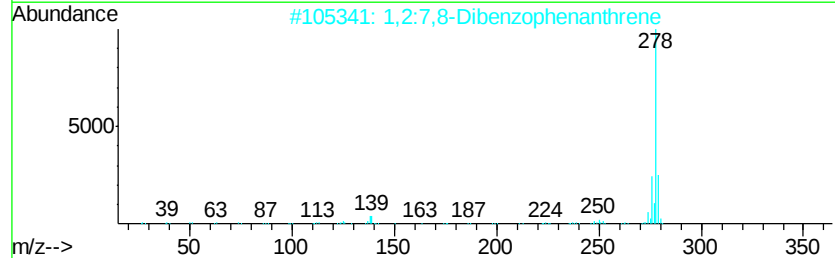
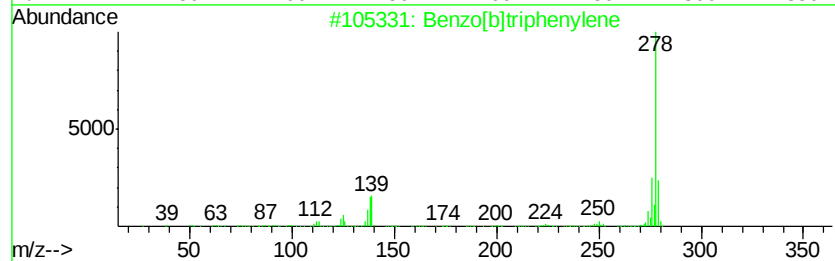
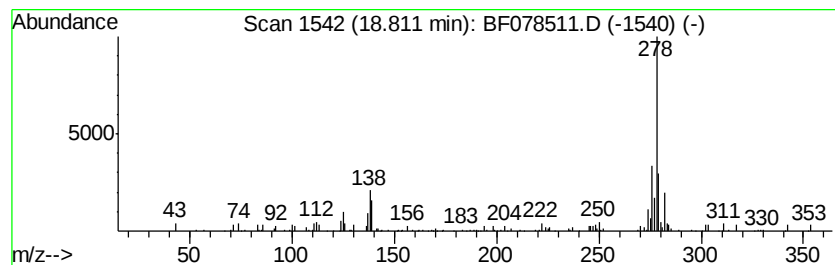
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Benzo[b]triphenylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.81	15.24 ng	234295	Perylene-d12	17.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]triphenylene	278	C22H14	000215-58-7	93
2		1,2:7,8-Dibenzophenanthrene	278	C22H14	000213-46-7	86
3		Pyrene, 1-phenyl-	278	C22H14	005101-27-9	83
4		Pentaphene	278	C22H14	000222-93-5	64
5		Benzo[a]naphthacene	278	C22H14	000226-88-0	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
Data File : BF078511.D
Acq On : 14 Apr 2015 20:57
Operator : TP/IZ
Sample : G1828-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
HS-SB02B

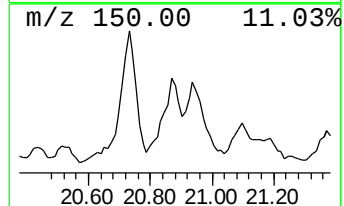
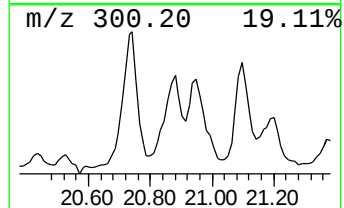
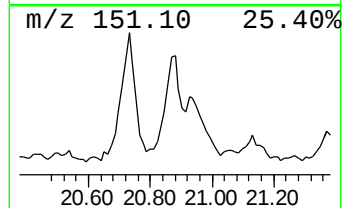
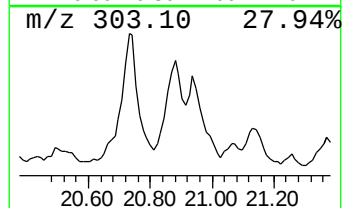
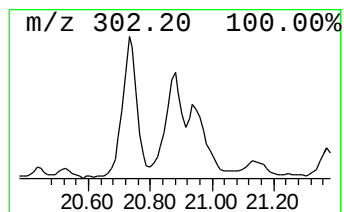
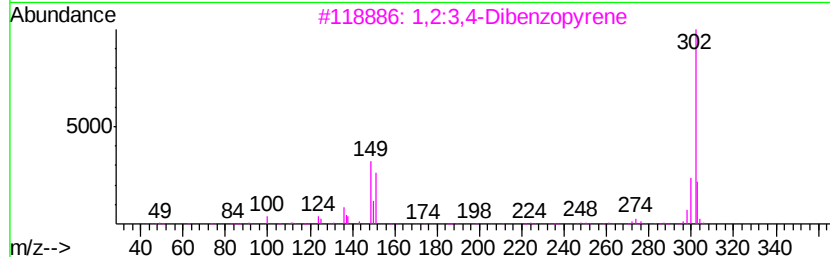
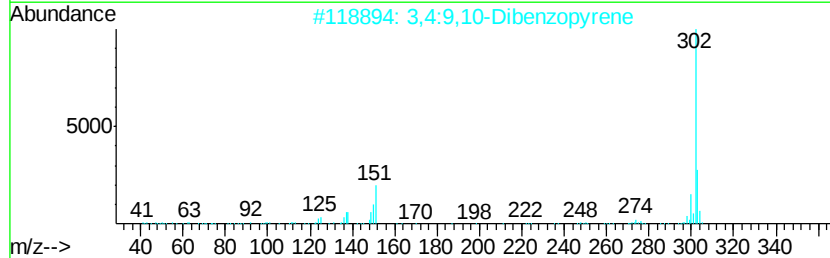
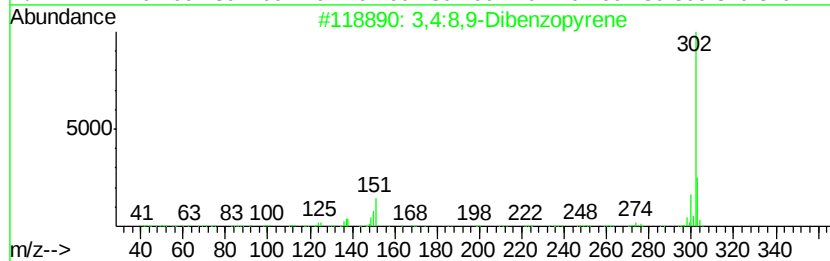
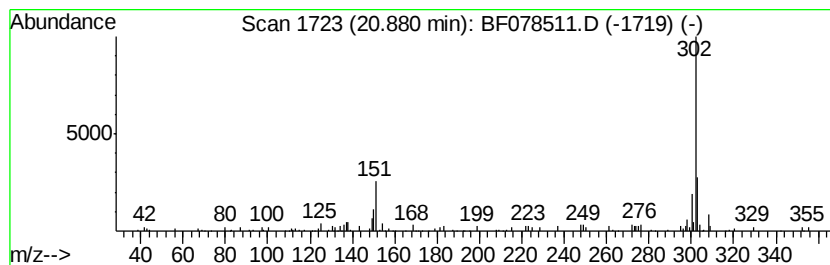
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 3,4:8,9-Dibenzopyrene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.88	6.36 ng	97742	Perylene-d12	17.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,4:8,9-Dibenzopyrene	302	C24H14	000189-64-0	94
2			3,4:9,10-Dibenzopyrene	302	C24H14	000189-55-9	90
3			1,2:3,4-Dibenzopyrene	302	C24H14	000191-30-0	76
4			Dibenz(a,e)aceanthrylene	302	C24H14	005385-75-1	62
5			1,2:4,5-Dibenzopyrene	302	C24H14	000192-65-4	58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041415\
 Data File : BF078511.D
 Acq On : 14 Apr 2015 20:57
 Operator : TP/IZ
 Sample : G1828-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB02B

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
unknown1.50	1.50	10.2	ng	54927	1	6.80	107731	20.0
2-Pentanone, 4-hy...	4.44	13.5	ng	72843	1	6.80	107731	20.0
unknown6.52	6.52	24.1	ng	129689	1	6.80	107731	20.0
Decahydro-4,4,8,9...	9.80	6.3	ng	90534	3	10.54	286769	20.0
unknown10.96	10.96	9.1	ng	130918	3	10.54	286769	20.0
Phenol, 4,4'-(1-m...	14.24	12.8	ng	404732	5	15.63	630362	20.0
Dinaphtho[1,2-b:1...	17.17	6.2	ng	94716	6	17.37	307543	20.0
11H-Indeno[2,1-a]...	17.55	9.6	ng	147818	6	17.37	307543	20.0
3,4-Dibromobenzal...	17.87	5.5	ng	84921	6	17.37	307543	20.0
unknown18.69	18.69	10.0	ng	153589	6	17.37	307543	20.0
Benzo[b]triphenylene	18.81	15.2	ng	234295	6	17.37	307543	20.0
3,4:8,9-Dibenzopy...	20.88	6.4	ng	97742	6	17.37	307543	20.0