

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111516\
 Data File : BF091178.D
 Acq On : 15 Nov 2016 14:59
 Operator : UM/SJ
 Sample : H5636-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C19013041

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

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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110316.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	1.744	3	5	61	rVB	3569231	11043010	100.00%	20.426%
2	4.773	266	270	273	rBV	2166009	3544587	32.10%	6.556%
3	5.219	307	309	319	rBV	1163650	1388199	12.57%	2.568%
4	6.282	400	402	410	rBV	725720	1014364	9.19%	1.876%
5	6.396	410	412	415	rBV	2133068	2872438	26.01%	5.313%
6	6.636	431	433	444	rVB	873147	1001624	9.07%	1.853%
7	6.784	444	446	449	rBV	2484828	2914326	26.39%	5.391%
8	6.979	460	463	466	rBV2	57307	116495	1.05%	0.215%
9	7.207	481	483	485	rBV	2306563	2306414	20.89%	4.266%
10	7.413	498	501	505	rVB3	69377	124606	1.13%	0.230%
11	7.665	521	523	526	rBV	283729	314754	2.85%	0.582%
12	7.916	543	545	555	rVB3	1083743	2159270	19.55%	3.994%
13	8.202	568	570	573	rBV4	73562	117852	1.07%	0.218%
14	8.316	577	580	582	rVB	87177	131604	1.19%	0.243%
15	8.396	582	587	588	rBV	179755	272812	2.47%	0.505%
16	8.579	599	603	605	rBV3	76970	151694	1.37%	0.281%
17	8.636	606	608	611	rVB	352991	326773	2.96%	0.604%
18	8.727	614	616	618	rBV	548796	754868	6.84%	1.396%
19	8.773	618	620	622	rVB	91581	139492	1.26%	0.258%
20	8.830	622	625	627	rBV3	59292	130537	1.18%	0.241%
21	9.002	637	640	642	rBV	4749241	4484524	40.61%	8.295%
22	9.139	650	652	654	rVB	78570	122110	1.11%	0.226%
23	9.207	654	658	660	rBV	219141	381161	3.45%	0.705%
24	9.265	660	663	667	rVV	313838	481367	4.36%	0.890%
25	9.333	667	669	670	rVV	507010	506413	4.59%	0.937%
26	9.368	670	672	674	rVV2	192191	320057	2.90%	0.592%
27	9.402	674	675	677	rVB	126701	113311	1.03%	0.210%
28	9.448	677	679	684	rVB2	288154	479418	4.34%	0.887%
29	9.528	684	686	689	rBV	327836	400441	3.63%	0.741%
30	9.665	696	698	700	rBV	1333143	1704912	15.44%	3.154%
31	9.699	700	701	707	rVV2	141689	290195	2.63%	0.537%
32	9.790	707	709	710	rVV	119222	155750	1.41%	0.288%
33	9.825	710	712	714	rVV2	103226	151226	1.37%	0.280%
34	9.882	714	717	718	rVV2	150697	225025	2.04%	0.416%

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35	9.916	718	720	722 rVV	145976	194243	1.76%	0.359%
36	9.985	724	726	728 rVV	166340	217717	1.97%	0.403%
37	10.076	732	734	735 rVV	92624	170268	1.54%	0.315%
38	10.099	735	736	739 rVV2	115881	139470	1.26%	0.258%
39	10.213	743	746	749 rVV2	215319	302190	2.74%	0.559%
40	10.282	749	752	755 rVV	282035	434248	3.93%	0.803%
41	10.328	755	756	759 rVV2	77256	142803	1.29%	0.264%
42	10.453	765	767	770 rVV2	1036151	1289298	11.68%	2.385%
43	10.796	795	797	800 rVB2	128315	182098	1.65%	0.337%
44	10.911	803	807	809 rVB4	52861	110651	1.00%	0.205%
45	11.048	813	819	823 rBV3	103637	229507	2.08%	0.425%
46	11.151	825	828	830 rBV	1893615	1797371	16.28%	3.325%
47	12.739	964	967	969 rBV	4880083	5402120	48.92%	9.992%
48	13.779	1055	1058	1060 rBV	1612353	1553080	14.06%	2.873%
49	15.151	1175	1178	1187 rBV	881605	1255906	11.37%	2.323%

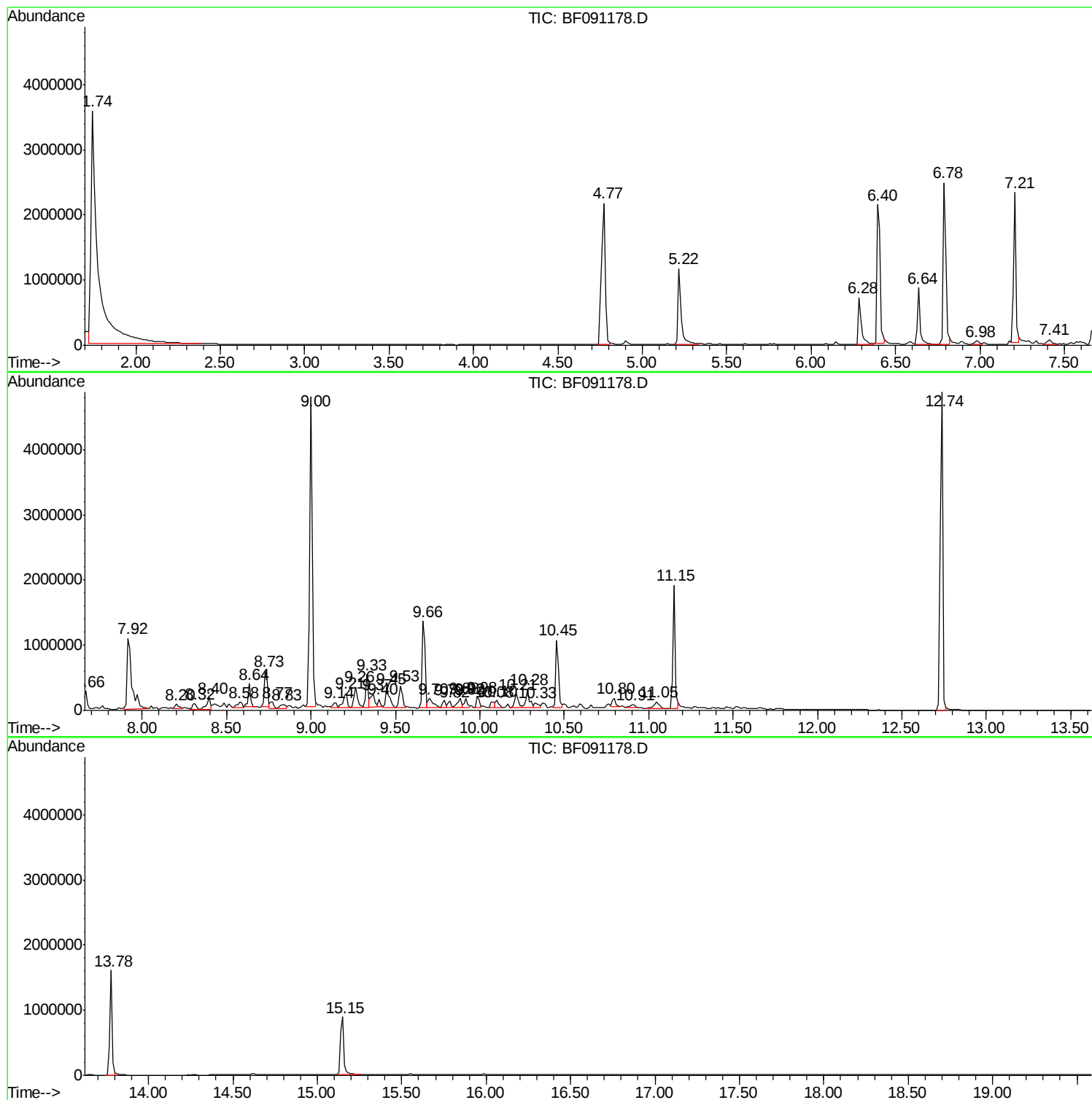
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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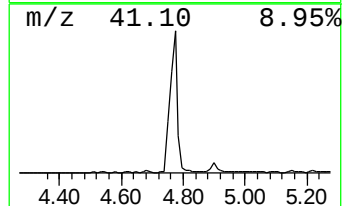
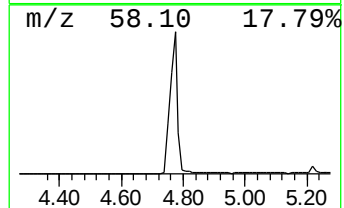
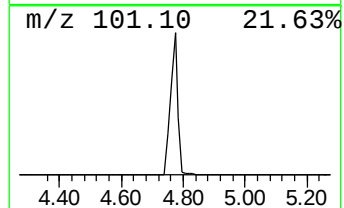
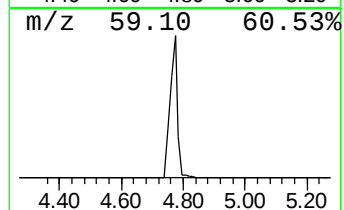
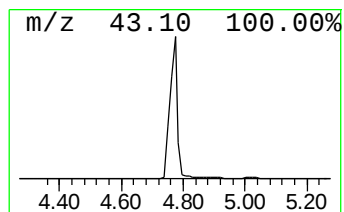
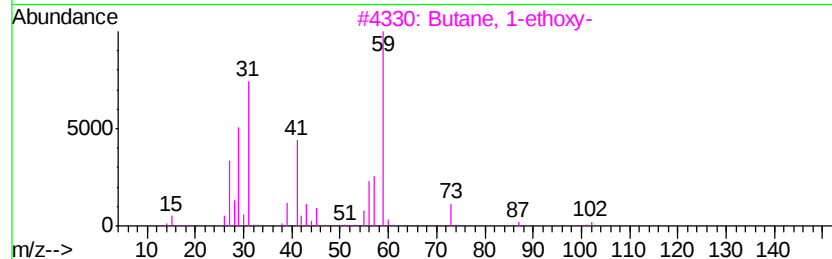
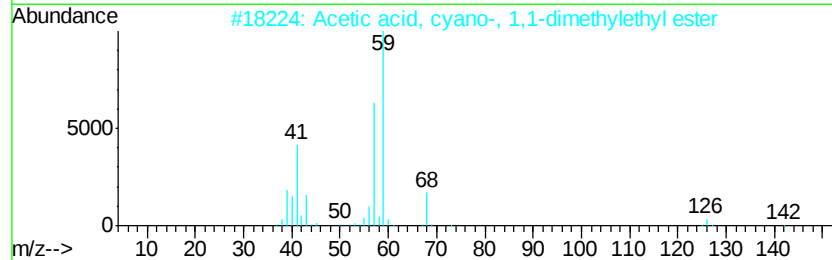
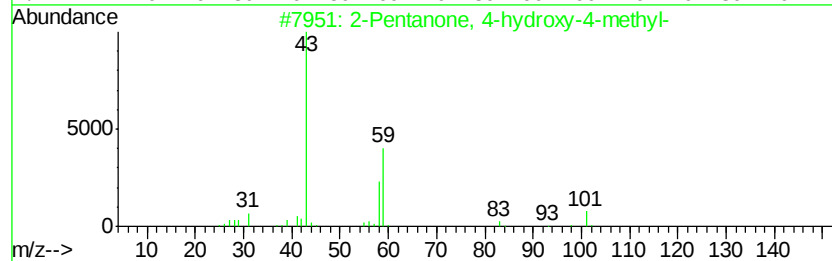
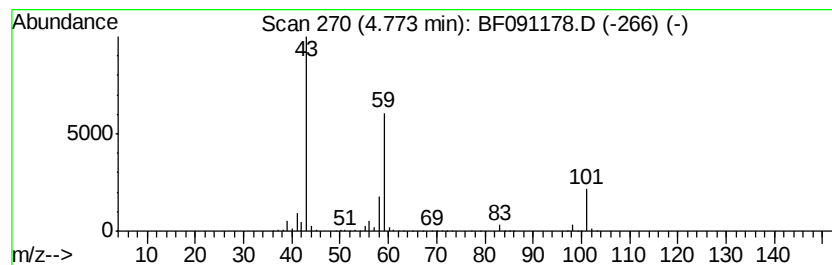
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.77	70.78 ng	3544590	1,4-Dichlorobenzene-d4	6.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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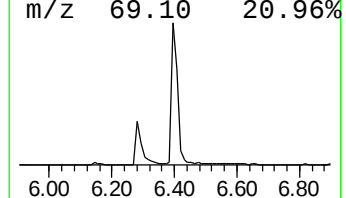
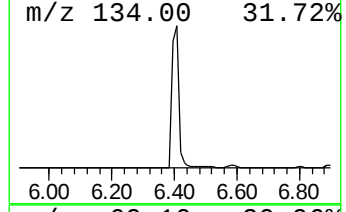
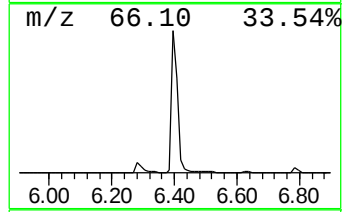
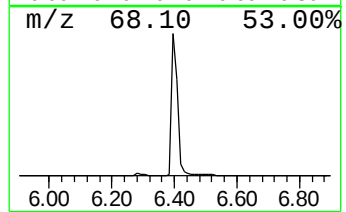
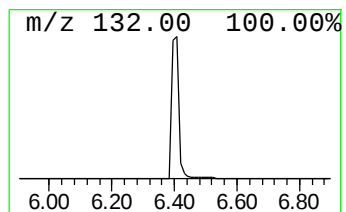
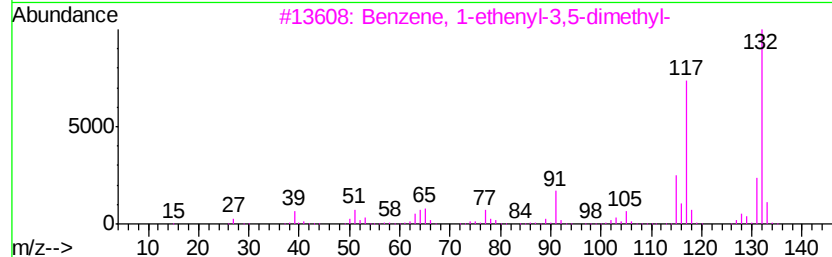
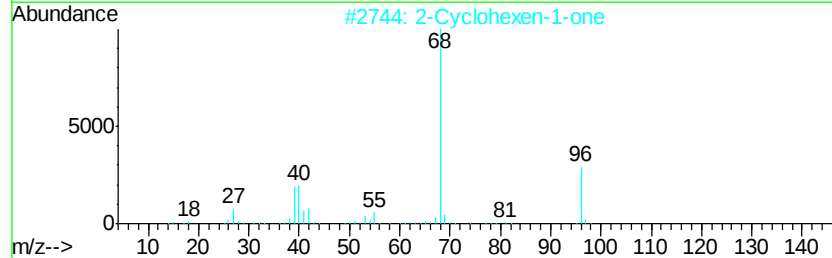
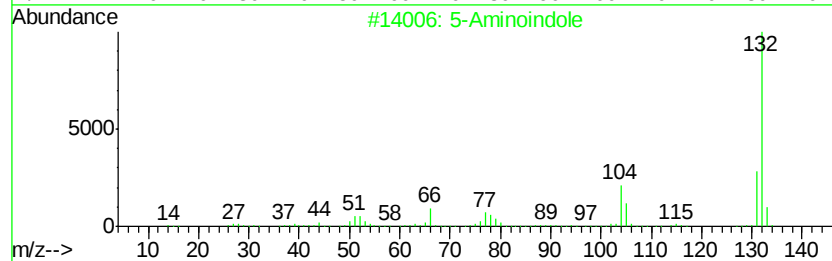
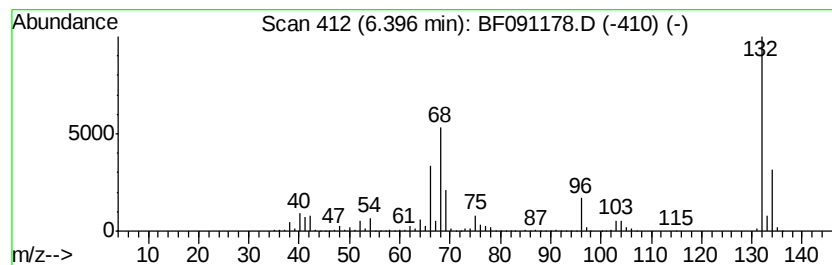
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 unknown6.40 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.40	57.36 ng	2872440	1,4-Dichlorobenzene-d4	6.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	22
2		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	11
3		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
4		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9
5		Bicyclo[4.2.0]octa-1,3,5-triene, ...	132	C10H12	028749-81-7	9



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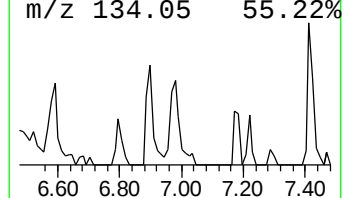
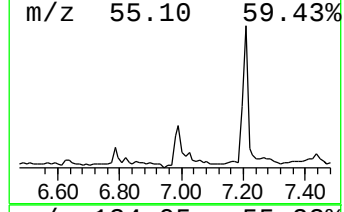
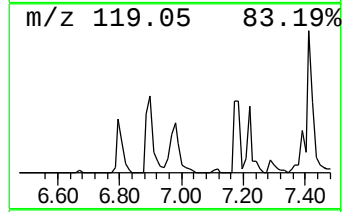
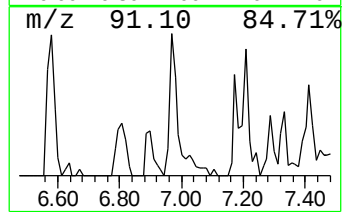
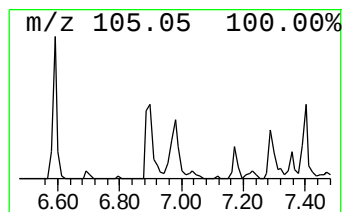
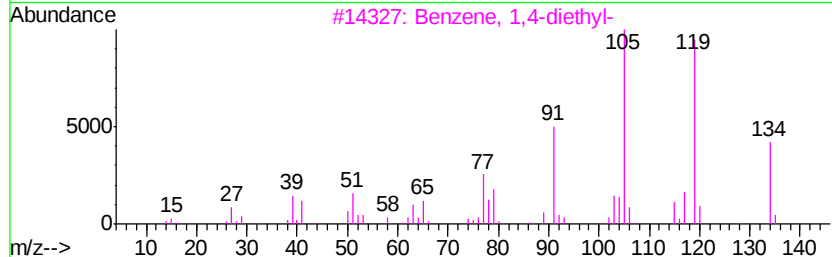
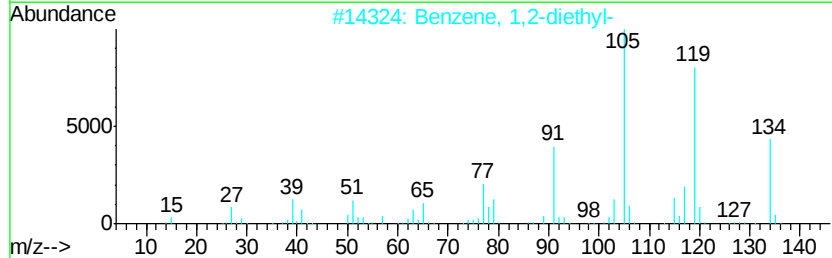
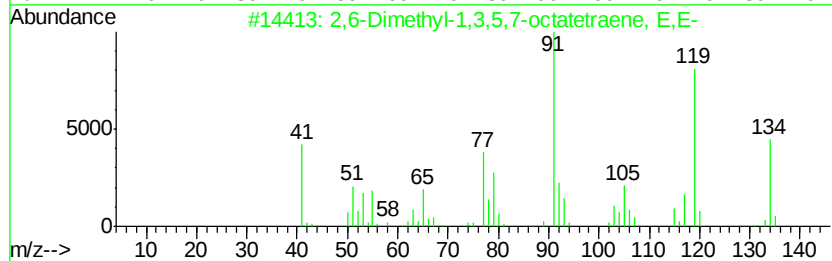
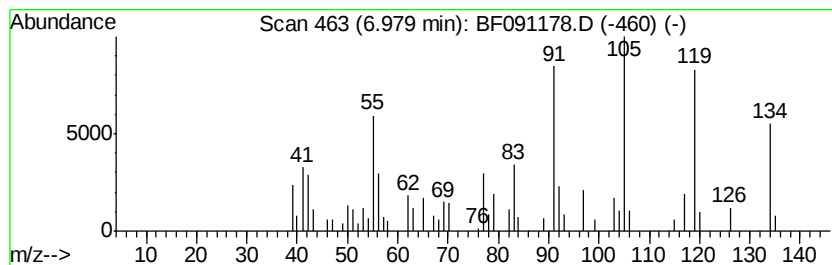
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 2,6-Dimethyl-1,3,5,7-octate... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.98	2.33 ng	116495	1,4-Dichlorobenzene-d4	6.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,6-Dimethyl-1,3,5,7-octatetraen...	134	C10H14	000460-01-5	93
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	64
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	64
4		Cyclobutene, bis(1-methylethylid...	134	C10H14	003642-14-6	64
5		1,3,8-p-Menthatriene	134	C10H14	021195-59-5	62



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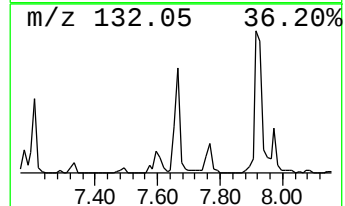
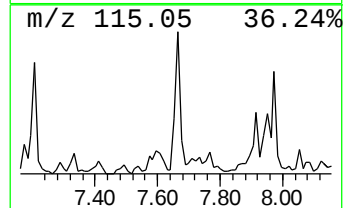
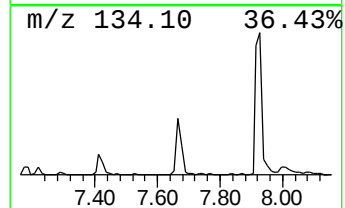
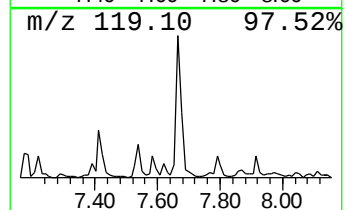
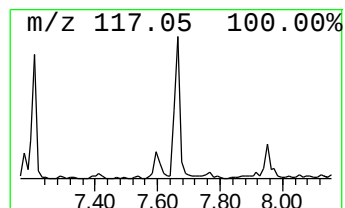
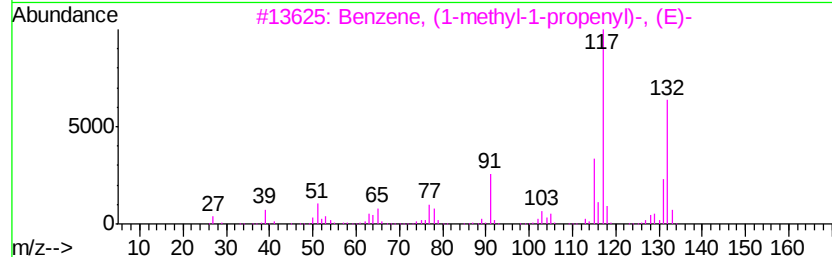
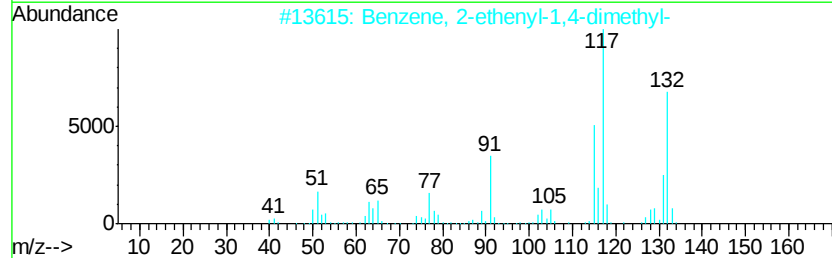
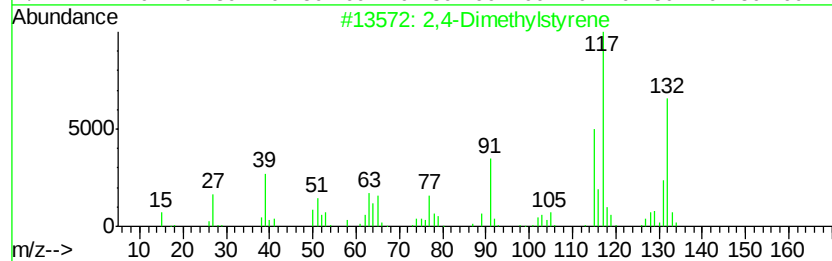
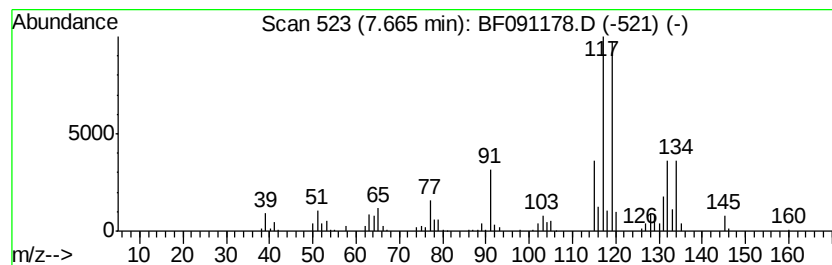
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 6 2,4-Dimethylstyrene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.66	2.92 ng	314754	Naphthalene-d8	7.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,4-Dimethylstvrene	132	C10H12	002234-20-0	74
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	70
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	50
4		1-Phenyl-1-butene	132	C10H12	000824-90-8	50
5		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	50



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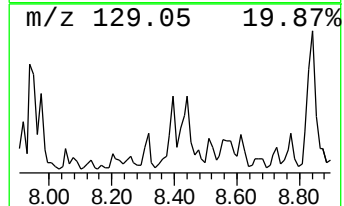
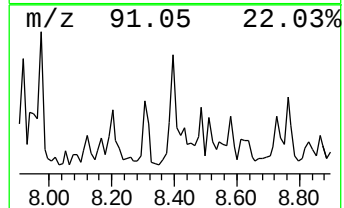
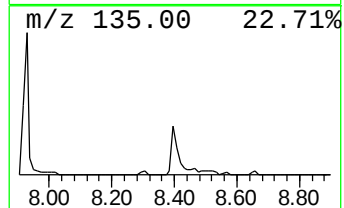
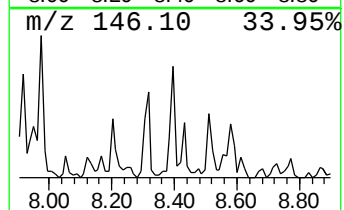
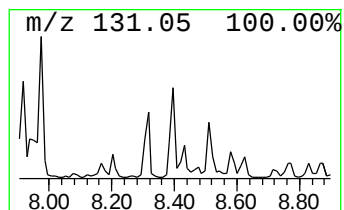
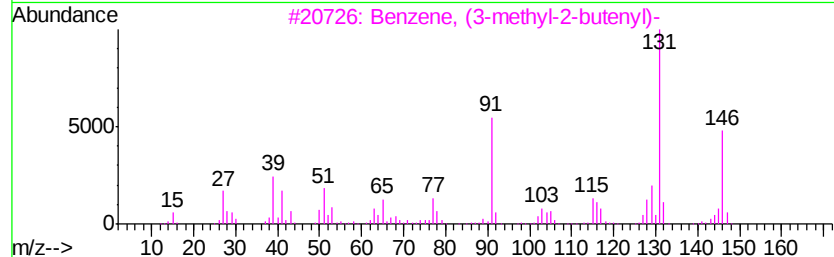
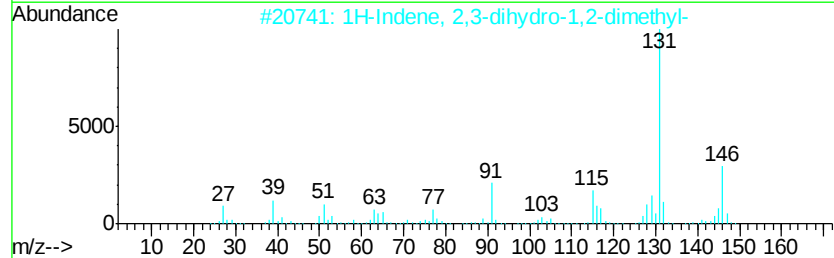
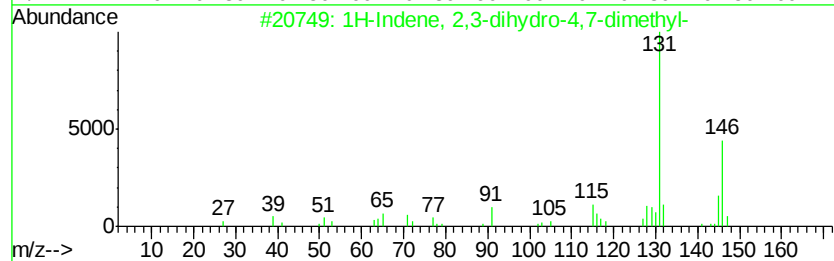
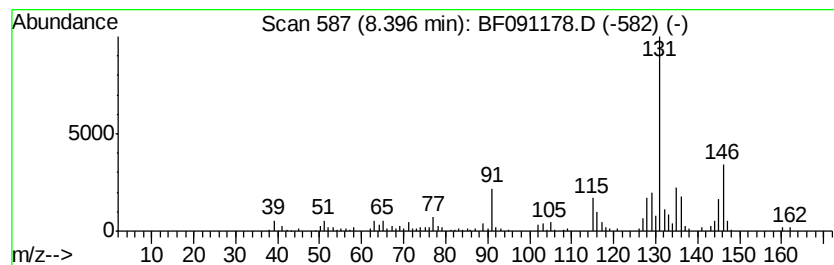
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.40	2.53 ng	272812	Naphthalene-d8	7.93

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
2		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
3		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	81
4		.alpha.,.beta.,.beta.-Trimethyls...	146	C11H14	000769-57-3	76
5		1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	70



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111516\
Data File : BF091178.D
Acq On : 15 Nov 2016 14:59
Operator : UM/SJ
Sample : H5636-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

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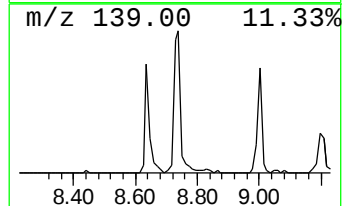
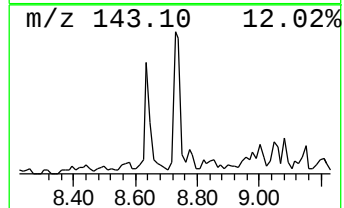
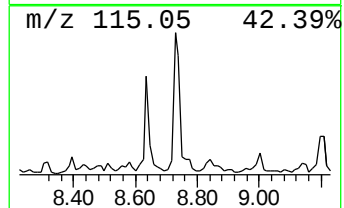
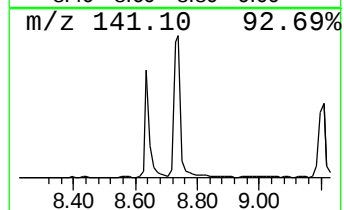
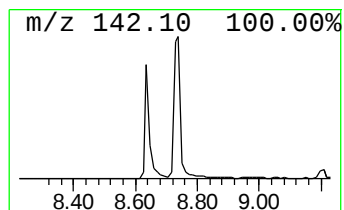
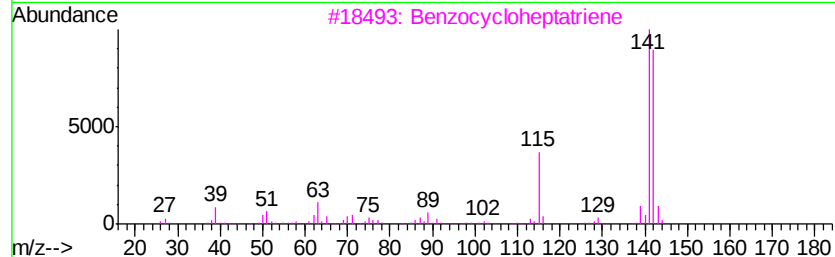
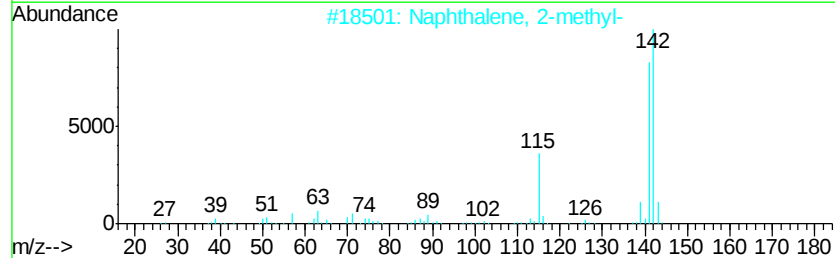
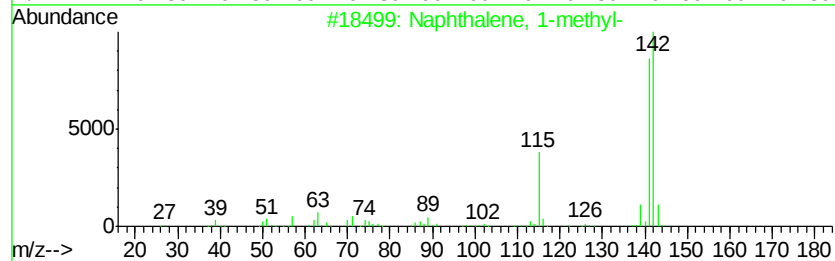
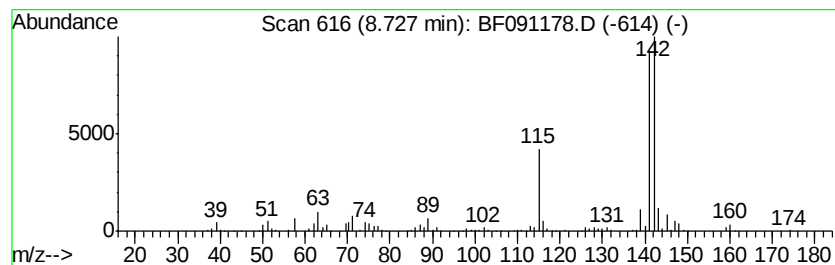
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.73	6.99 ng	754868	Naphthalene-d8	7.93

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	93
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
5		Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	83



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111516\
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ClientSampleId :
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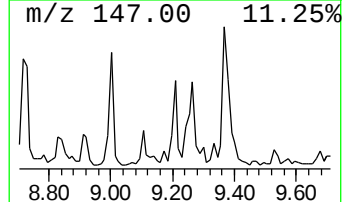
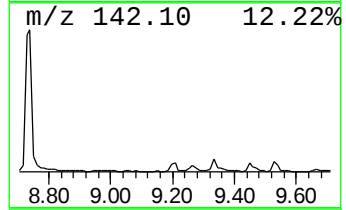
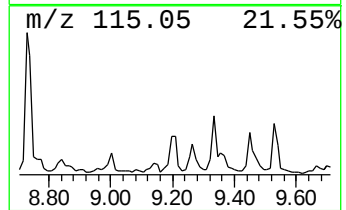
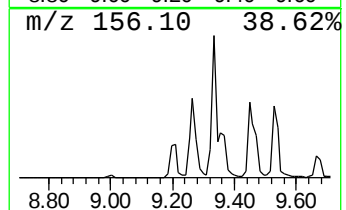
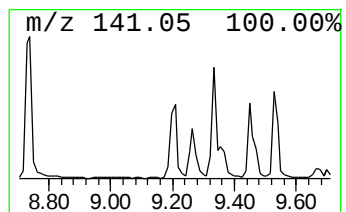
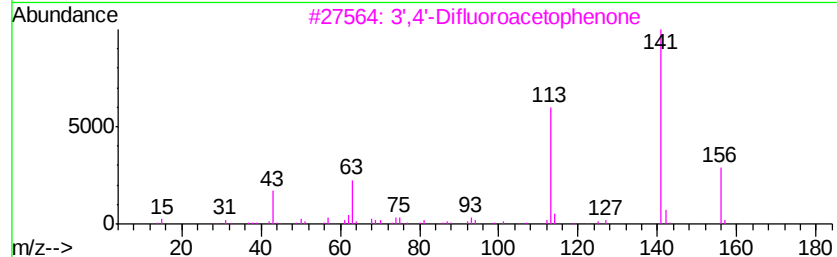
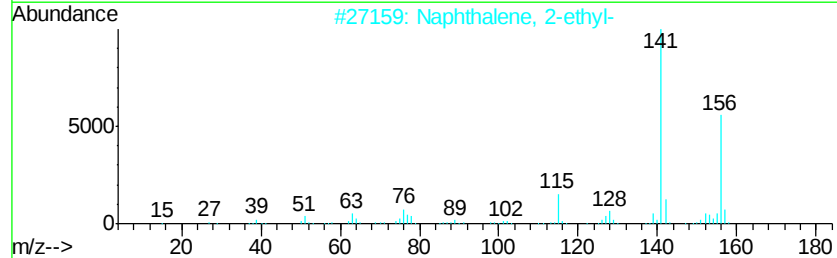
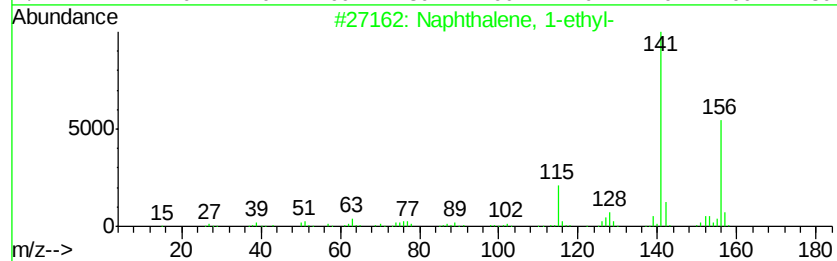
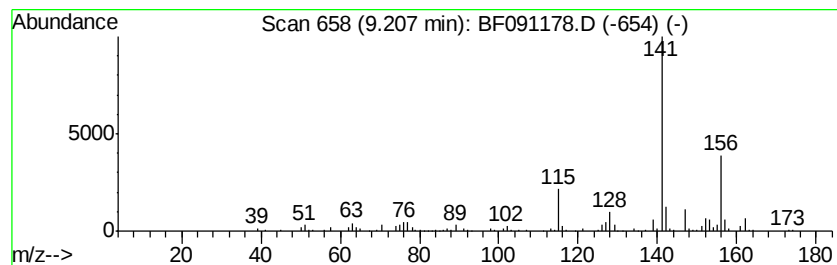
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, 1-ethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.21	4.47 ng	381161	Acenaphthene-d10	9.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	96
2		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	95
3		3',4'-Difluoroacetophenone	156	C8H6F2O	000369-33-5	53
4		2-Thiophenecarboxylic acid, 5-et...	156	C7H8O2S	023229-72-3	50
5		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	49



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111516\
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Acq On : 15 Nov 2016 14:59
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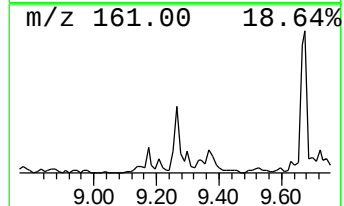
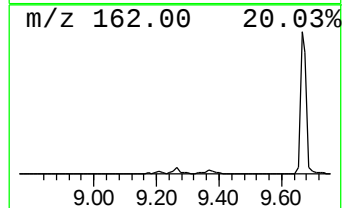
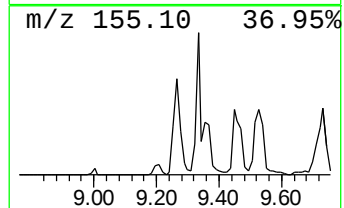
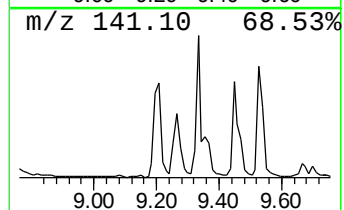
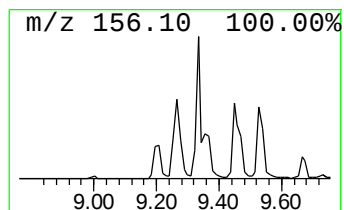
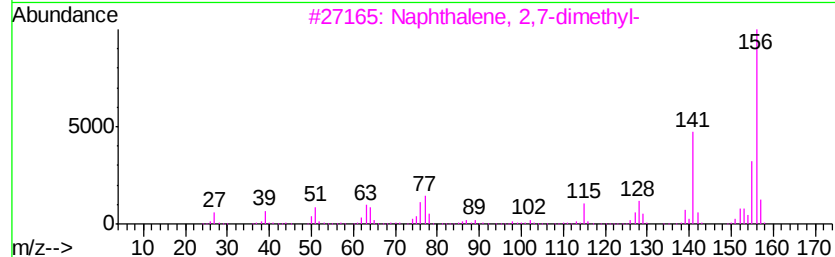
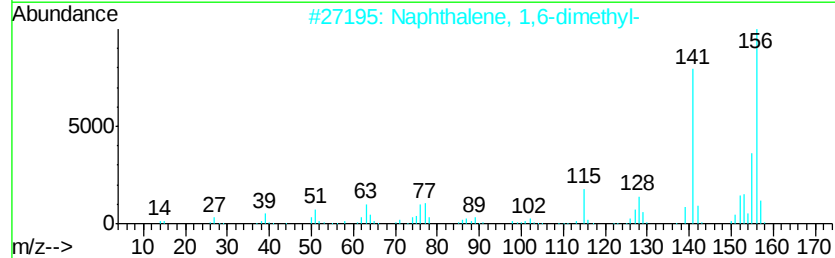
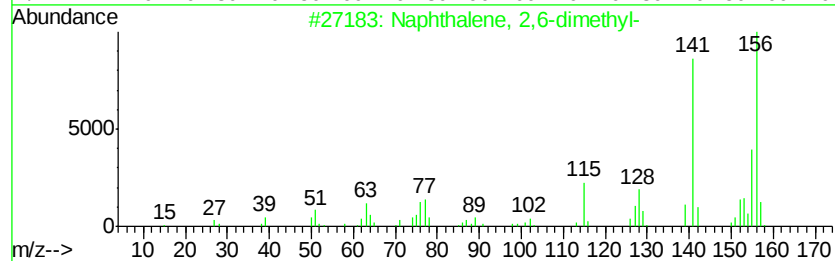
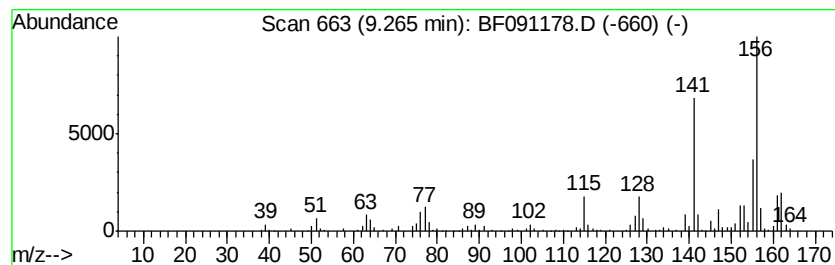
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Naphthalene, 2,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.26	5.65 ng	481367	Acenaphthene-d10	9.66	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111516\
Data File : BF091178.D
Acq On : 15 Nov 2016 14:59
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ALS Vial : 13 Sample Multiplier: 1

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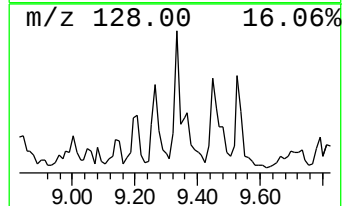
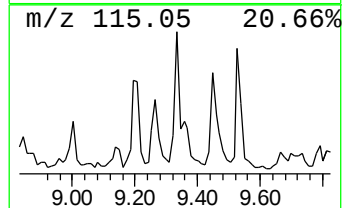
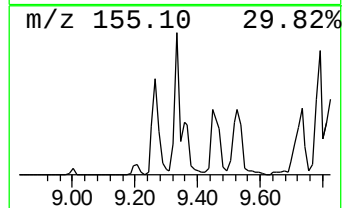
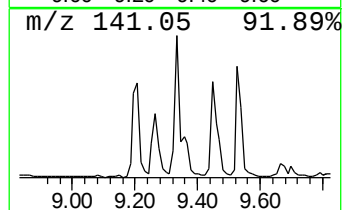
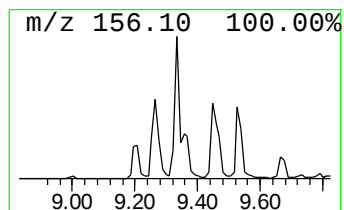
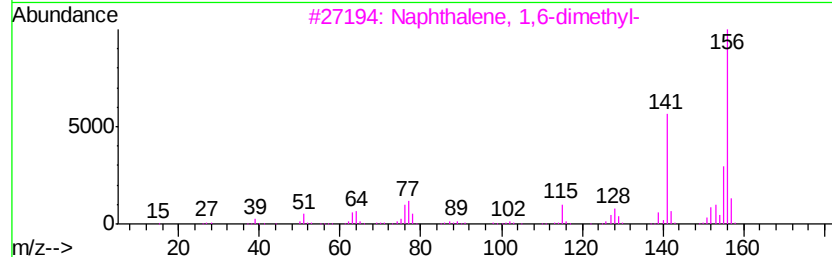
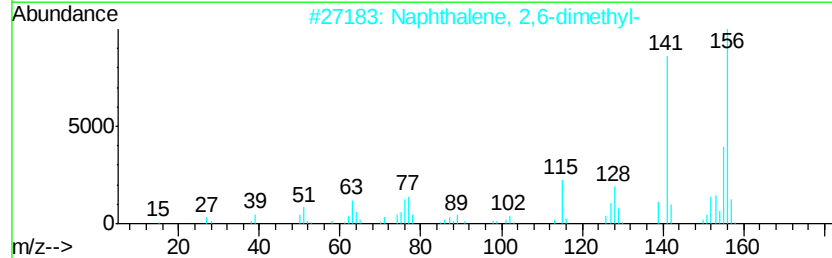
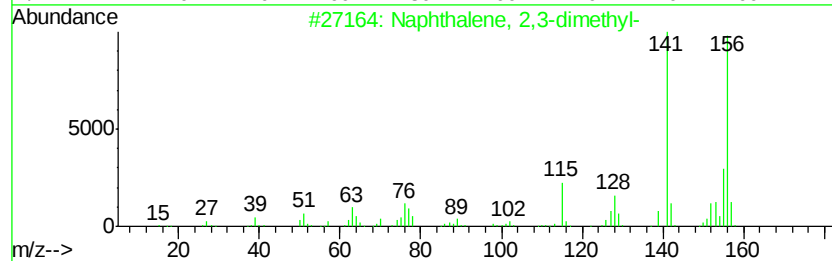
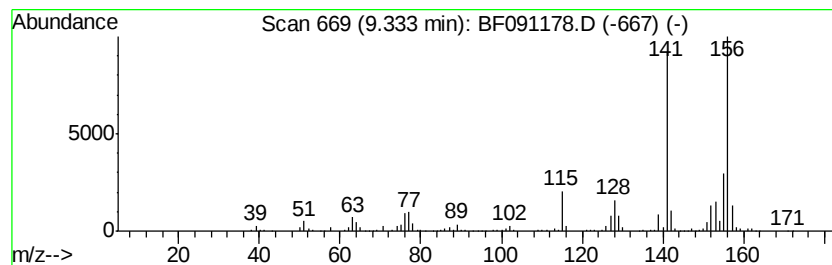
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Naphthalene, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.33	5.94 ng	506413	Acenaphthene-d10	9.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111516\
Data File : BF091178.D
Acq On : 15 Nov 2016 14:59
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ClientSampleId :
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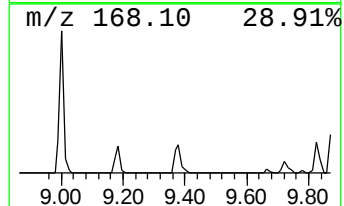
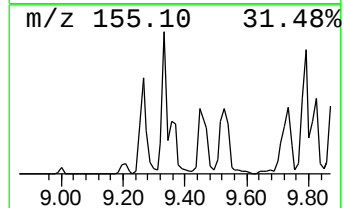
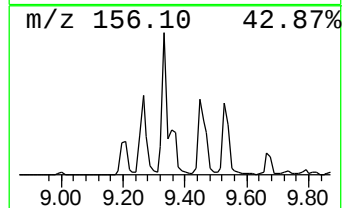
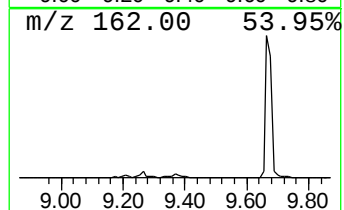
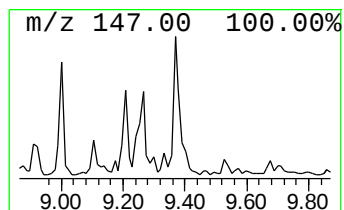
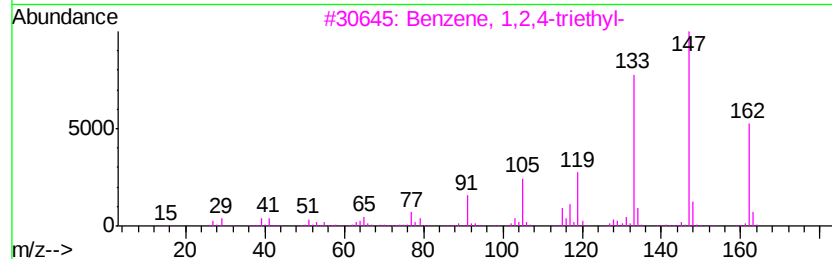
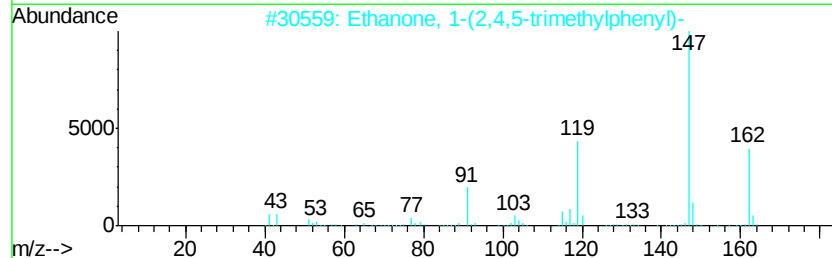
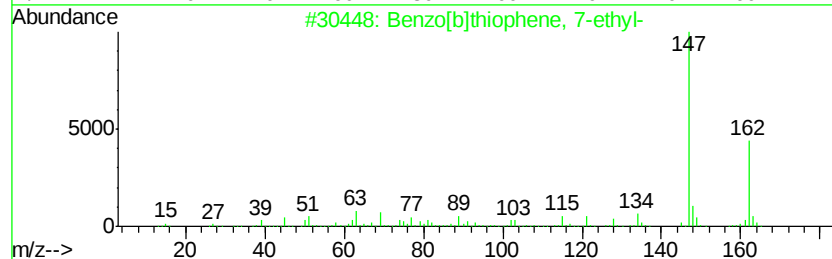
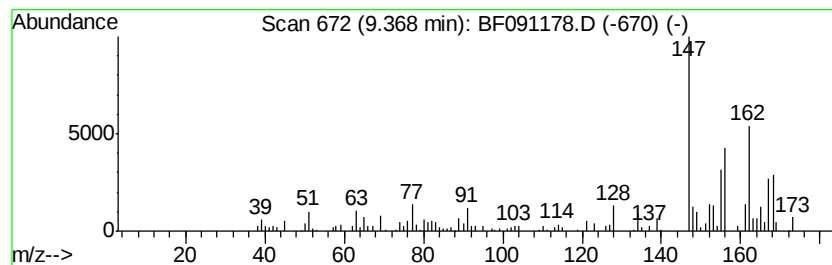
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown9.37 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.37	3.75 ng	320057	Acenaphthene-d10	9.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]thiophene, 7-ethyl-	162	C10H10S	016587-42-1	38
2		Ethanone, 1-(2,4,5-trimethylphen...	162	C11H14O	002040-07-5	38
3		Benzene, 1,2,4-triethyl-	162	C12H18	000877-44-1	38
4		Ethanone, 1-(2,3,4-trimethylphen...	162	C11H14O	001467-36-3	38
5		Benzene, hexamethyl-	162	C12H18	000087-85-4	38



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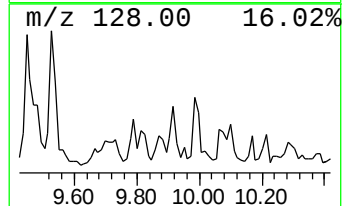
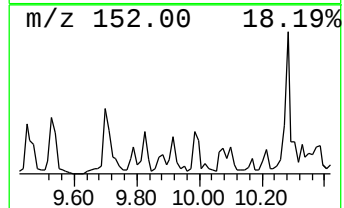
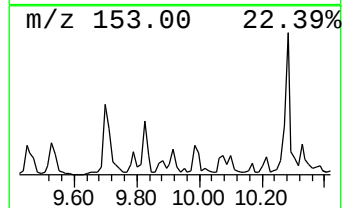
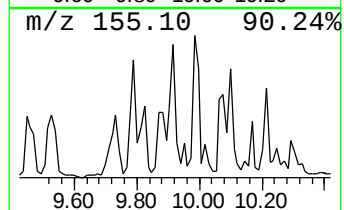
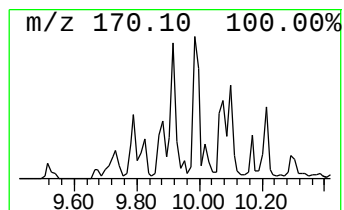
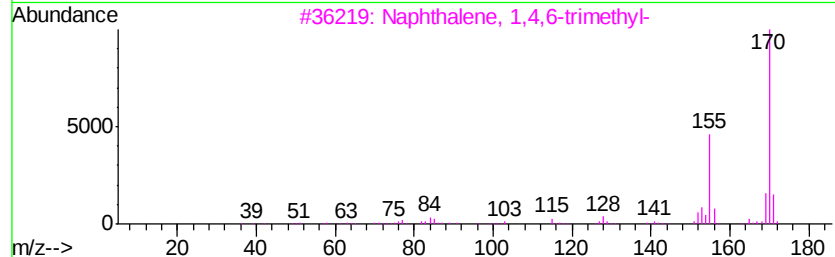
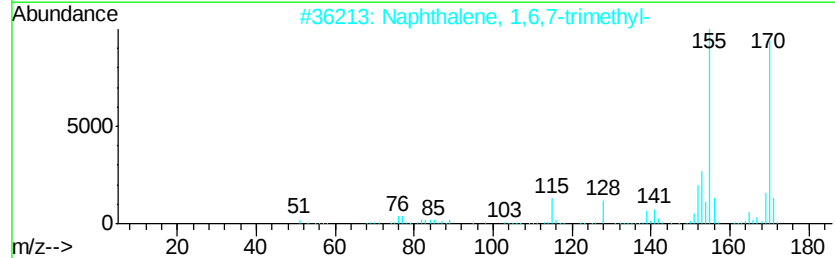
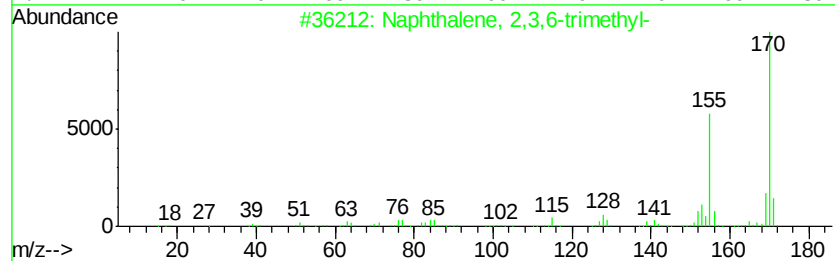
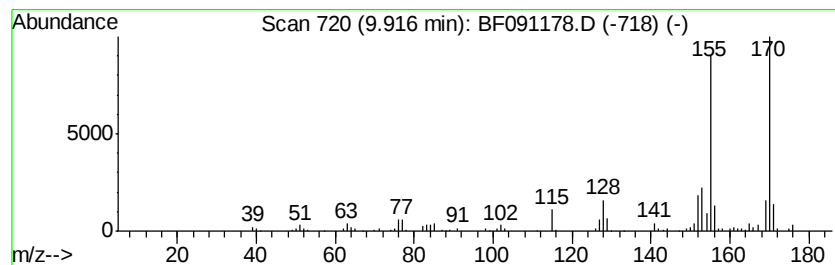
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Naphthalene, 2,3,6-trimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.92	2.28 ng	194243	Acenaphthene-d10	9.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	94
4			Azulene, 4,6,8-trimethyl-	170	C13H14	000941-81-1	94
5			Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	94



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111516\
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ALS Vial : 13 Sample Multiplier: 1

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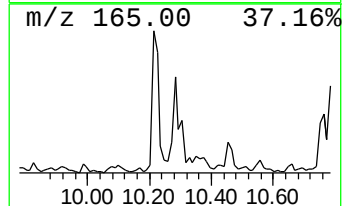
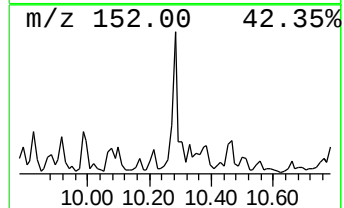
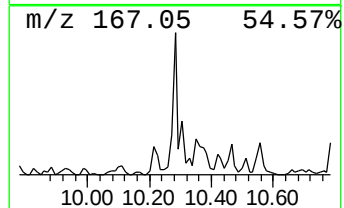
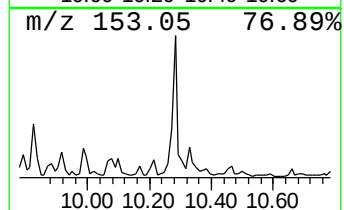
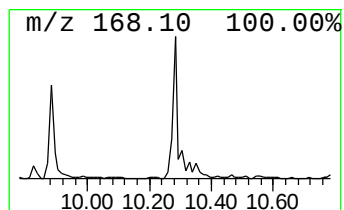
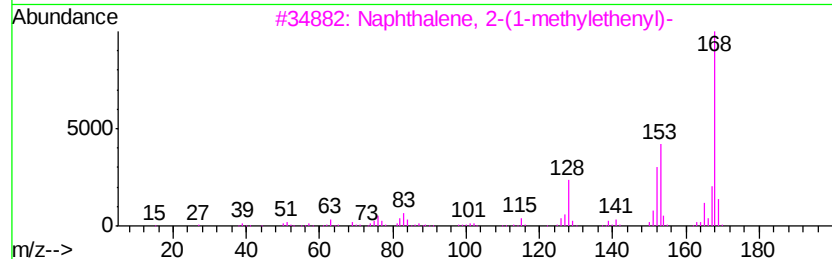
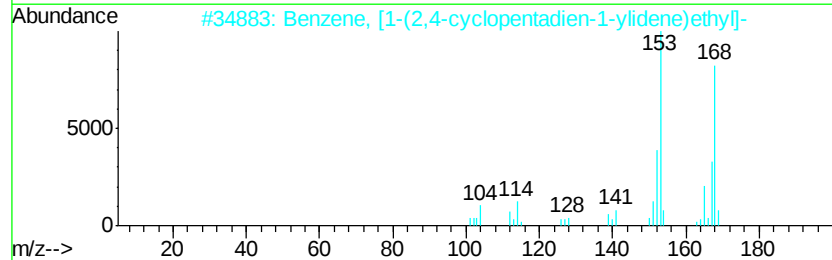
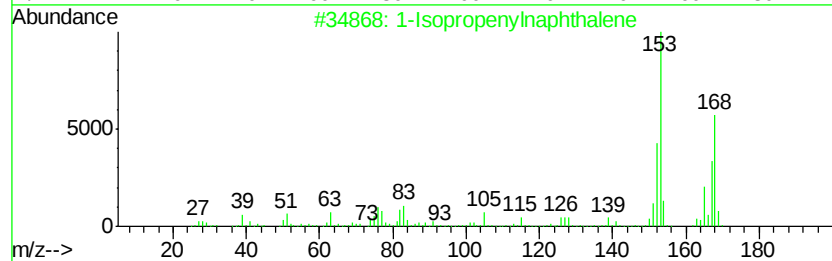
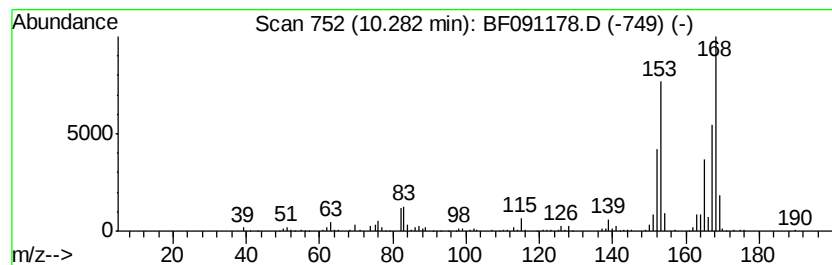
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 1-Isopropenylnaphthalene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.28	5.09 ng	434248	Acenaphthene-d10	9.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Isopropenylnaphthalene	168	C13H12	001855-47-6	87
2			Benzene, [1-(2,4-cyclopentadien-...	168	C13H12	002320-32-3	78
3			Naphthalene, 2-(1-methylethenyl)-	168	C13H12	003710-23-4	64
4			1,1'-Biphenyl, 3-methyl-	168	C13H12	000643-93-6	58
5			1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	50



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111516\
Data File : BF091178.D
Acq On : 15 Nov 2016 14:59
Operator : UM/SJ
Sample : H5636-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
C19013041

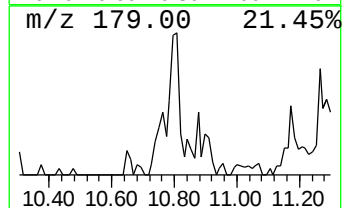
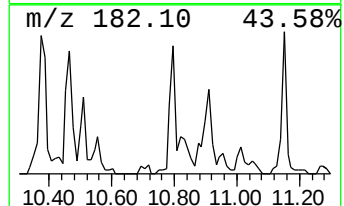
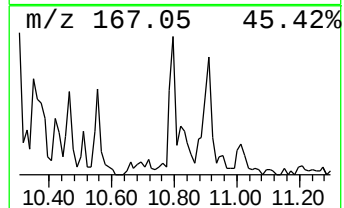
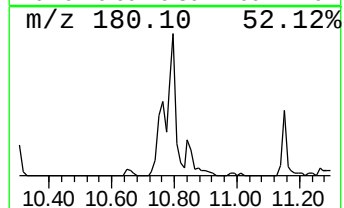
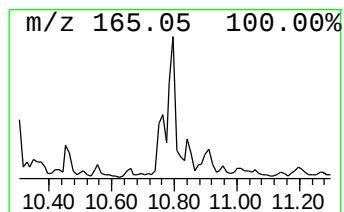
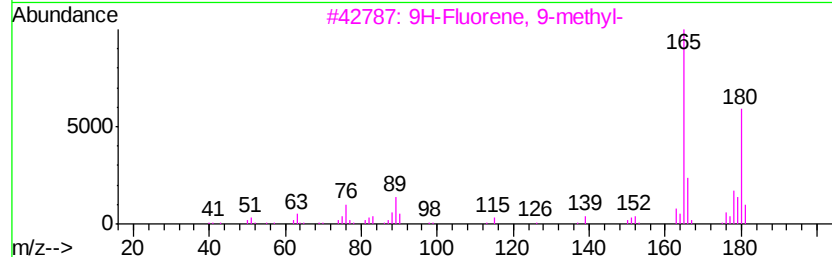
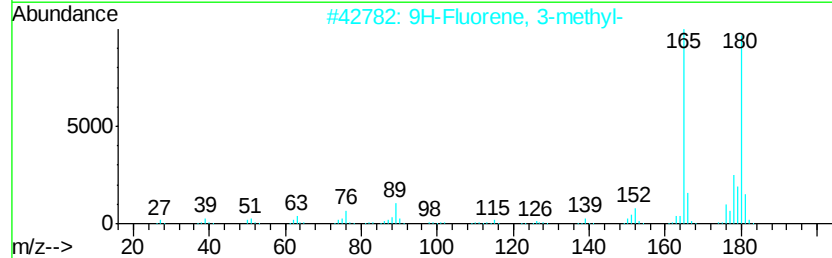
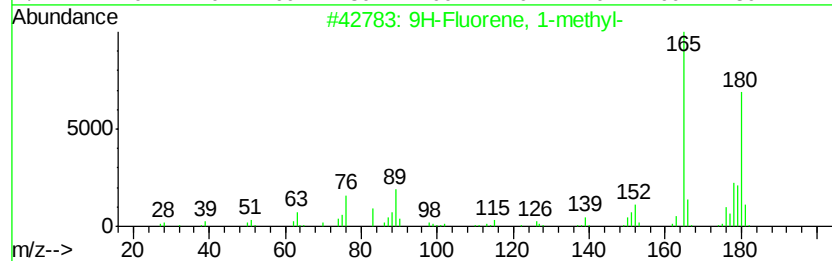
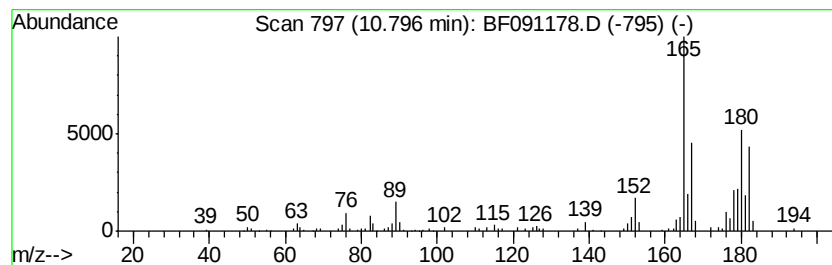
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 9H-Fluorene, 1-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.80	2.03 ng	182098	Phenanthrene-d10	11.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	70
2		9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	49
3		9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	49
4		9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	49
5		3-Buten-2-one, 4-(4-chlorophenyl)-	180	C10H9ClO	003160-40-5	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111516\
Data File : BF091178.D
Acq On : 15 Nov 2016 14:59
Operator : UM/SJ
Sample : H5636-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
C19013041

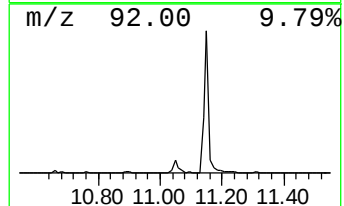
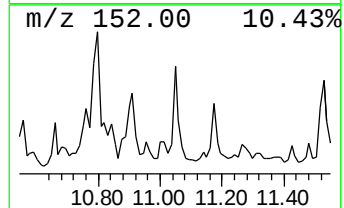
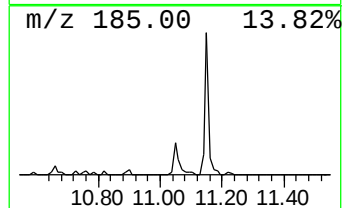
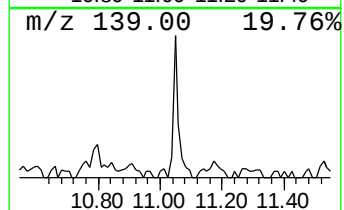
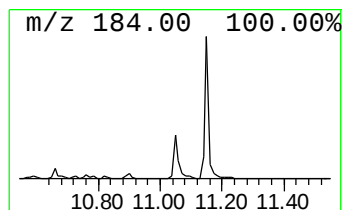
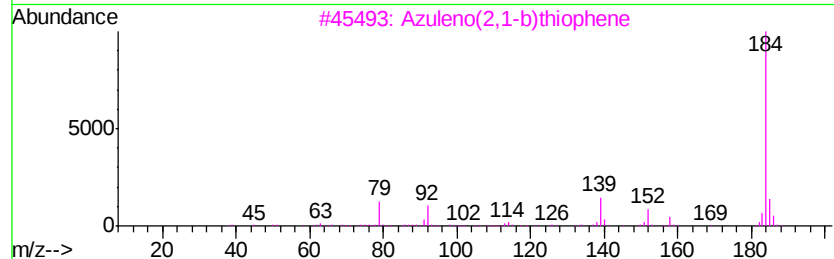
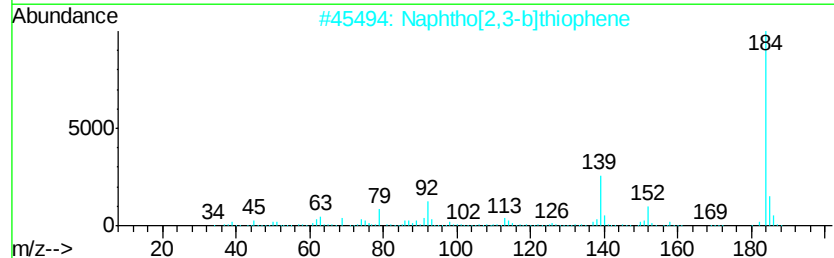
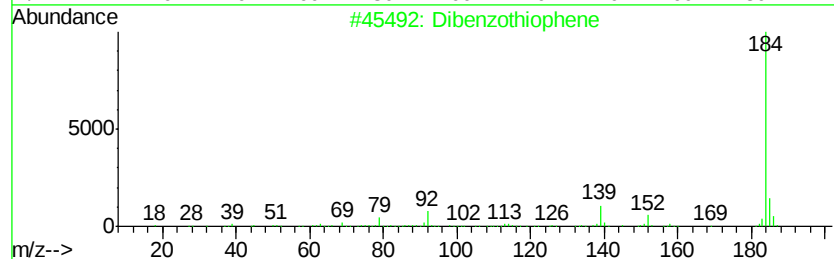
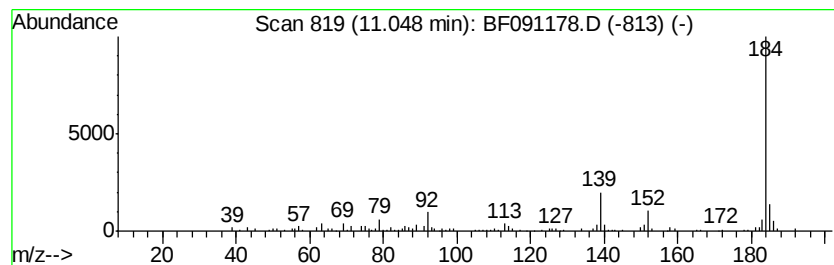
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 Dibenzothiophene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.05	2.55 ng	229507	Phenanthrene-d10	11.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dibenzothiophene	184	C12H8S	000132-65-0	97
2		Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	95
3		Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	91
4		2,5-Cyclohexadiene-1,4-dione, 2,...	184	C8H8O5	000085-23-4	52
5		2,2'-Bipyridine, 4,4'-dimethyl-	184	C12H12N2	001134-35-6	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111516\
 Data File : BF091178.D
 Acq On : 15 Nov 2016 14:59
 Operator : UM/SJ
 Sample : H5636-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C19013041

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110316.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-Pentanone, 4-hy...	4.77	70.8	ng	3544590	1	6.64	1001620	20.0
unknown6.40	6.40	57.4	ng	2872440	1	6.64	1001620	20.0
2,6-Dimethyl-1,3,...	6.98	2.3	ng	116495	1	6.64	1001620	20.0
2,4-Dimethylstyrene	7.66	2.9	ng	314754	2	7.93	2159270	20.0
1H-Indene, 2,3-di...	8.40	2.5	ng	272812	2	7.93	2159270	20.0
Naphthalene, 1-me...	8.73	7.0	ng	754868	2	7.93	2159270	20.0
Naphthalene, 1-et...	9.21	4.5	ng	381161	3	9.66	1704910	20.0
Naphthalene, 2,6-...	9.26	5.7	ng	481367	3	9.66	1704910	20.0
Naphthalene, 2,3-...	9.33	5.9	ng	506413	3	9.66	1704910	20.0
unknown9.37	9.37	3.8	ng	320057	3	9.66	1704910	20.0
Naphthalene, 2,3,...	9.92	2.3	ng	194243	3	9.66	1704910	20.0
1-Isopropenylnaph...	10.28	5.1	ng	434248	3	9.66	1704910	20.0
9H-Fluorene, 1-me...	10.80	2.0	ng	182098	4	11.15	1797370	20.0
Dibenzothiophene	11.05	2.5	ng	229507	4	11.15	1797370	20.0