

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111414\
 Data File : BF075528.D
 Acq On : 15 Nov 2014 8:47
 Operator : TP / JJ
 Sample : F4695-08
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FB111014

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-BF111214.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	1.761	56	59	62	rVB	360535	512622	21.98%	1.920%
2	1.830	64	65	73	rVB	96818	143072	6.14%	0.536%
3	1.967	73	77	84	rBV2	1166355	2331730	100.00%	8.733%
4	5.339	370	372	378	rBV	179950	201698	8.65%	0.755%
5	5.830	413	415	418	rBV	565673	756626	32.45%	2.834%
6	7.087	523	525	528	rBV	575118	487178	20.89%	1.825%
7	7.259	537	540	542	rBV	1471900	1417899	60.81%	5.310%
8	7.545	563	565	568	rBV	450988	540600	23.18%	2.025%
9	7.739	579	582	584	rVB	1405066	1301202	55.80%	4.873%
10	8.242	623	626	628	rBV	1147633	1087938	46.66%	4.075%
11	9.133	701	704	706	rVB	730903	772861	33.15%	2.895%
12	9.282	714	717	719	rBV2	27834	23652	1.01%	0.089%
13	9.339	720	722	724	rVV	104919	86225	3.70%	0.323%
14	9.431	727	730	732	rBV	30303	33935	1.46%	0.127%
15	9.705	752	754	757	rVB	191064	220612	9.46%	0.826%
16	10.151	792	793	795	rBV	27214	26263	1.13%	0.098%
17	10.219	795	799	801	rVV	43488	61190	2.62%	0.229%
18	10.276	802	804	807	rVB	202107	259354	11.12%	0.971%
19	10.368	809	812	816	rBV3	24683	47691	2.05%	0.179%
20	10.471	819	821	824	rVB	1634140	1939812	83.19%	7.265%
21	10.699	839	841	843	rBV	1961327	1841634	78.98%	6.897%
22	10.745	843	845	856	rVB	248648	436227	18.71%	1.634%
23	11.054	870	872	874	rBV	1074973	980621	42.06%	3.673%
24	11.088	874	875	880	rVB3	28193	29433	1.26%	0.110%
25	11.305	892	894	896	rBV	668642	852825	36.57%	3.194%
26	11.339	896	897	899	rVB	254014	195521	8.39%	0.732%
27	11.408	899	903	906	rBV2	629079	920293	39.47%	3.447%
28	11.465	906	908	913	rVB	598209	862544	36.99%	3.230%
29	11.739	929	932	936	rVB2	23549	41176	1.77%	0.154%
30	12.025	955	957	959	rVV	72321	79830	3.42%	0.299%
31	12.094	961	963	966	rVB	992292	1173660	50.33%	4.396%
32	12.197	970	972	975	rBV2	44943	70477	3.02%	0.264%
33	12.254	975	977	978	rBV	27676	28952	1.24%	0.108%
34	12.288	978	980	983	rVV	1146169	1205398	51.70%	4.514%

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Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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35	12.357	983	986	987	rVV	42864	53388	2.29%	0.200%
36	12.425	987	992	1001	rVV3	44997	180316	7.73%	0.675%
37	13.157	1054	1056	1059	rBV	873239	849420	36.43%	3.181%
38	13.694	1101	1103	1106	rBV	98659	86206	3.70%	0.323%
39	13.854	1115	1117	1120	rBV	29321	29808	1.28%	0.112%
40	13.922	1120	1123	1124	rVV2	29031	48070	2.06%	0.180%
41	13.945	1124	1125	1133	rVB2	18030	33244	1.43%	0.125%
42	14.837	1201	1203	1209	rBV3	25713	29100	1.25%	0.109%
43	15.157	1228	1231	1233	rBV	1847522	2095314	89.86%	7.847%
44	15.248	1236	1239	1248	rBV4	25522	78082	3.35%	0.292%
45	16.414	1335	1341	1342	rBV	40160	90626	3.89%	0.339%
46	16.448	1342	1344	1347	rVB	859092	949756	40.73%	3.557%
47	17.511	1433	1437	1442	rBV2	32105	87163	3.74%	0.326%
48	18.197	1494	1497	1500	rBV	736968	977262	41.91%	3.660%
49	18.654	1533	1537	1542	rBV5	21183	82802	3.55%	0.310%
50	20.277	1674	1679	1688	rVB5	10737	59288	2.54%	0.222%

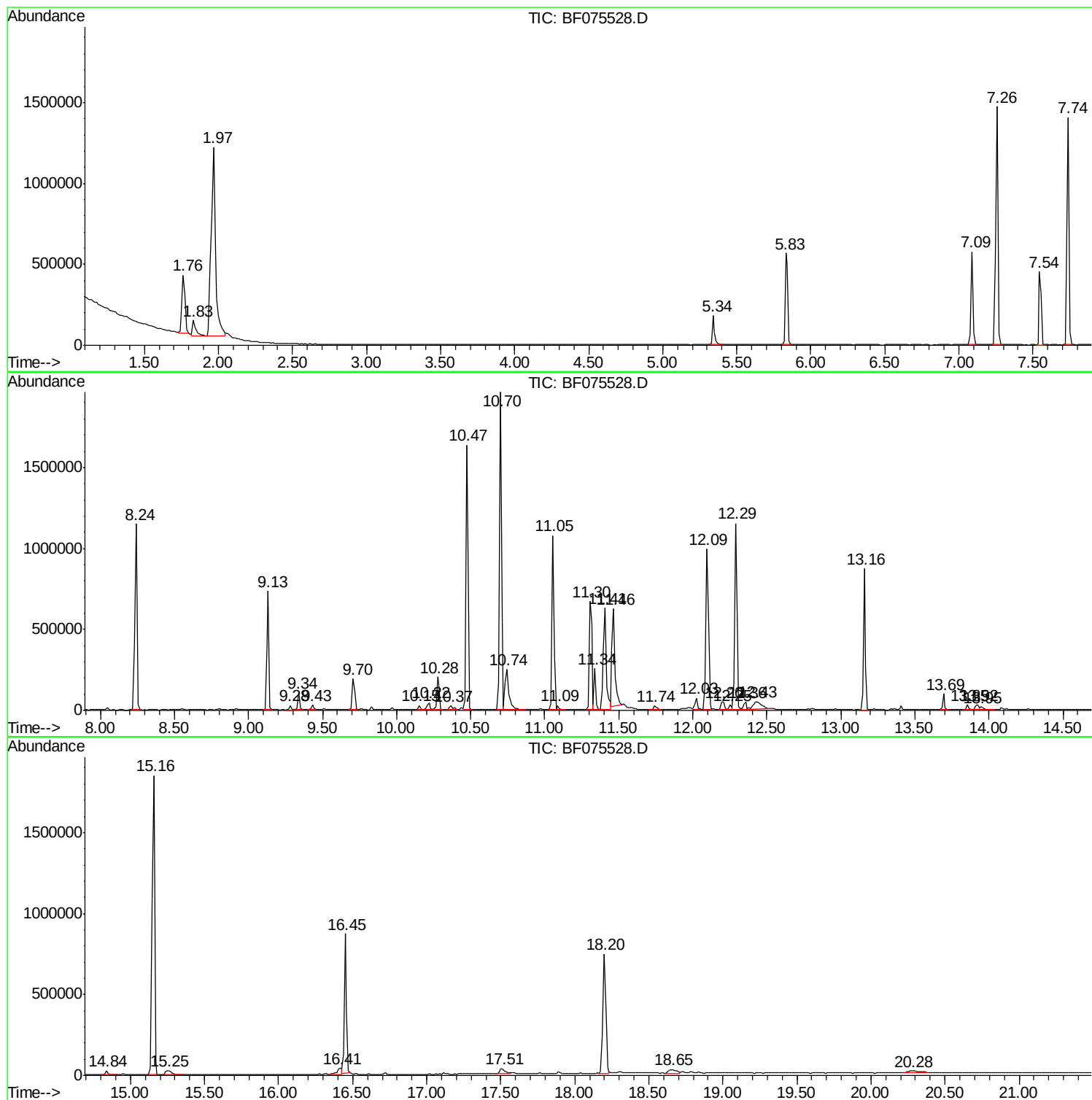
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF111214.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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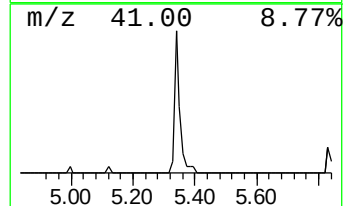
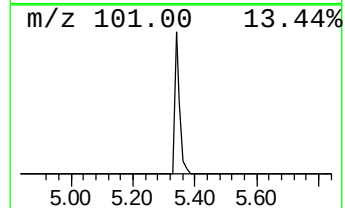
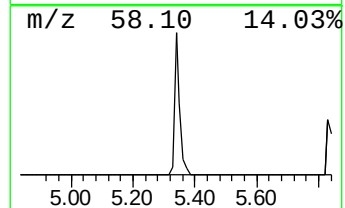
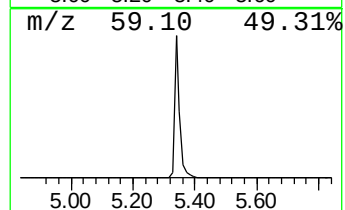
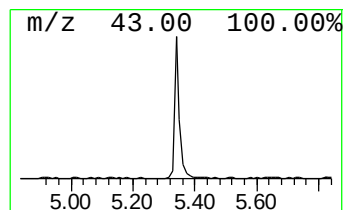
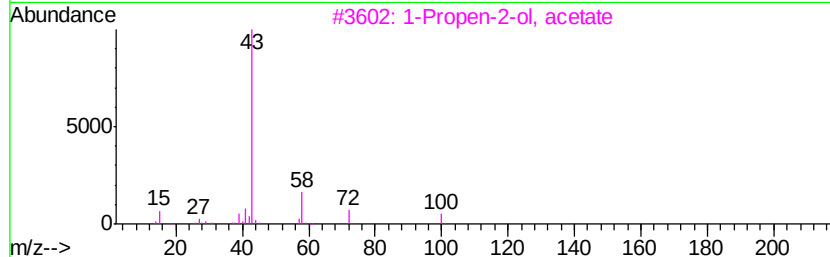
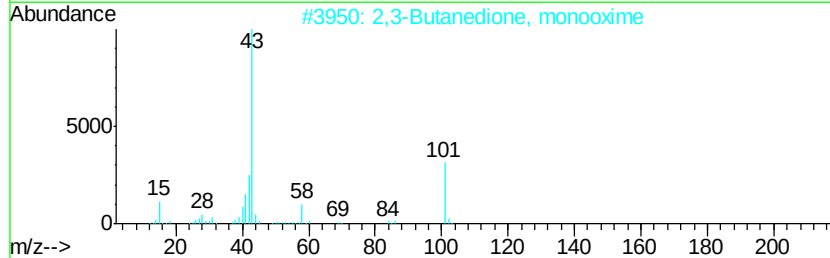
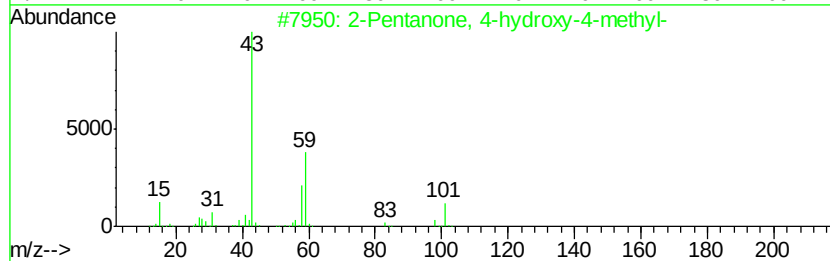
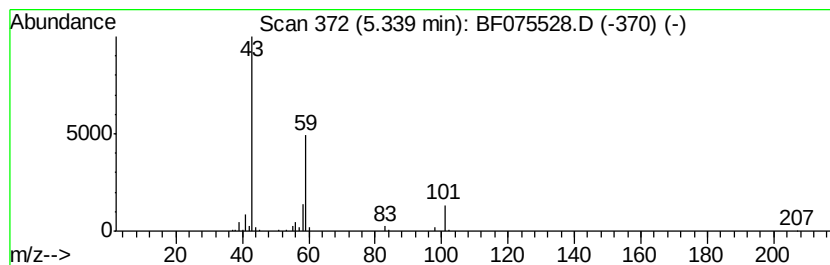
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF111214.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.34	7.46 ng	201698	1,4-Dichlorobenzene-d4	7.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	8
4		Diacetamide	101	C4H7NO2	000625-77-4	7
5		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	7



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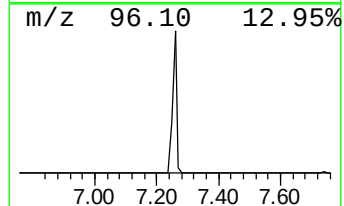
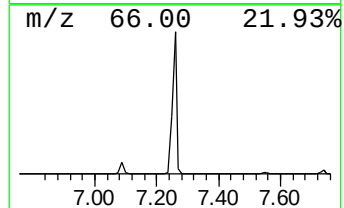
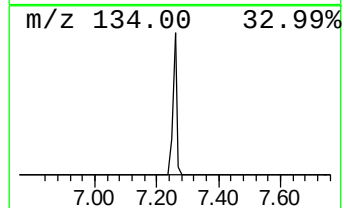
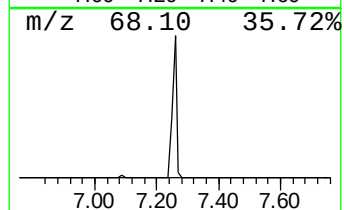
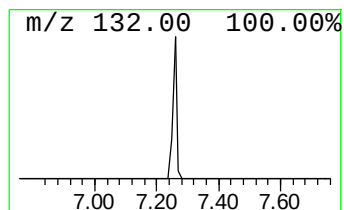
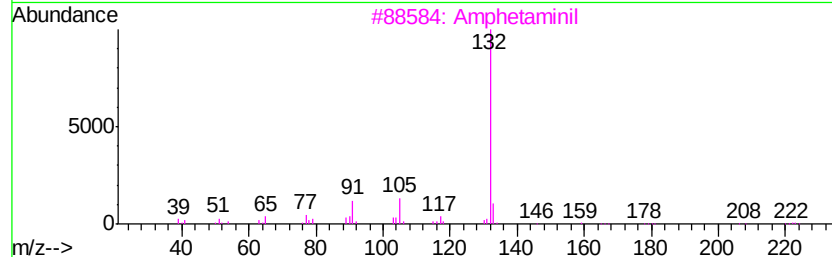
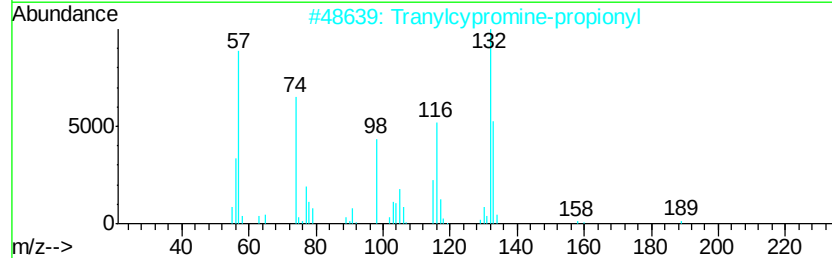
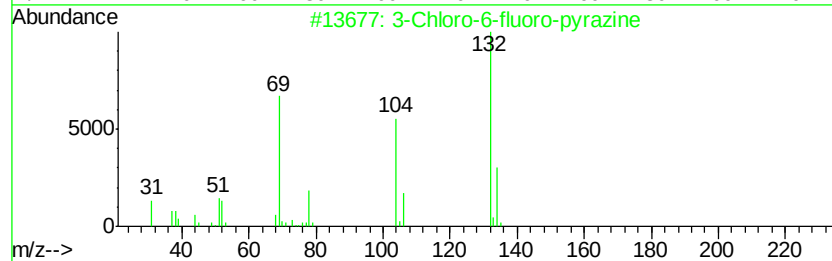
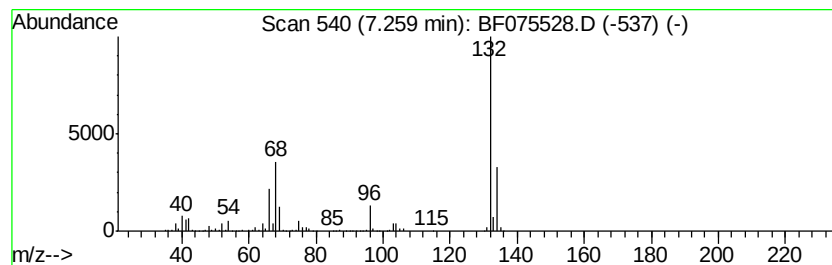
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown7.26 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.26	52.46 ng	1417900	1,4-Dichlorobenzene-d4	7.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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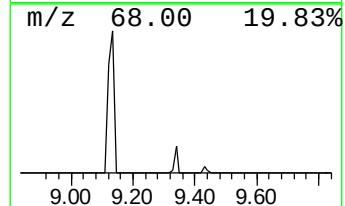
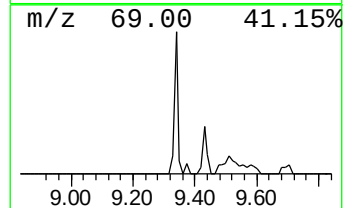
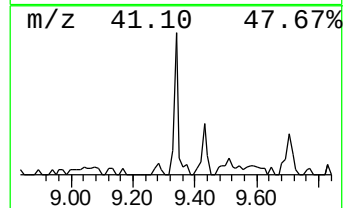
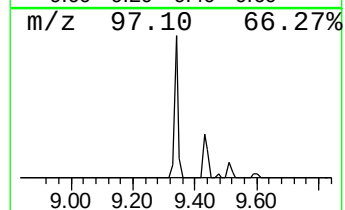
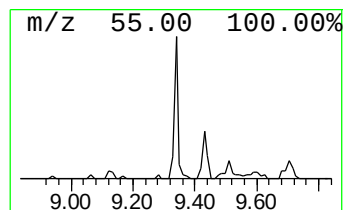
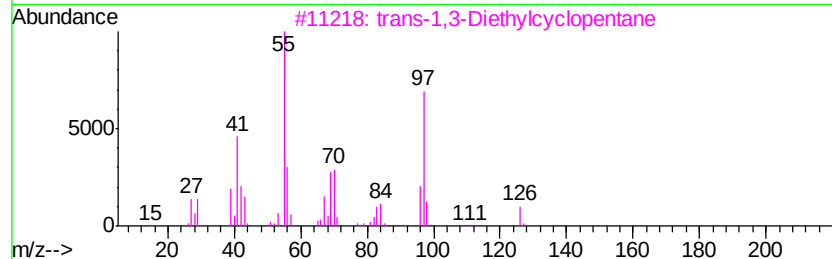
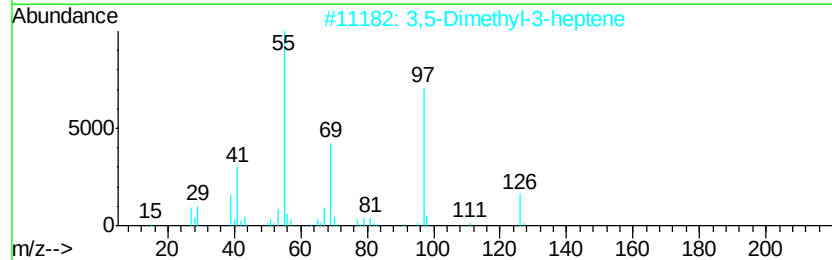
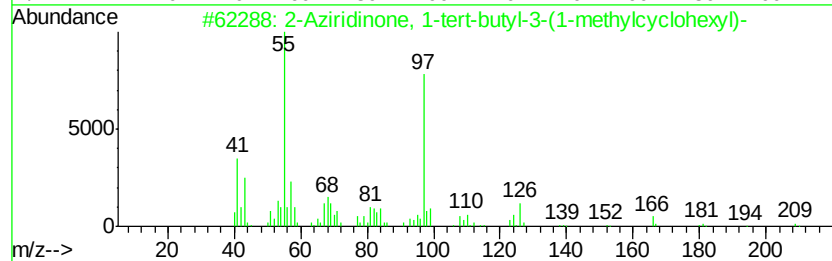
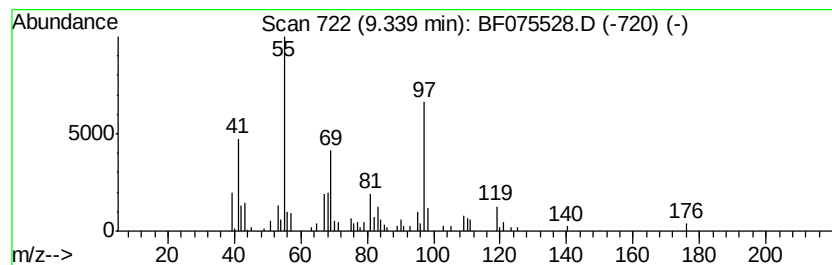
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 2-Aziridinone, 1-tert-butyl... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.34	2.23 ng	86225	Naphthalene-d8	9.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Aziridinone, 1-tert-butyl-3-(1...	209	C13H23NO	024161-49-7	59
2		3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	47
3		trans-1,3-Diethylcyclopentane	126	C9H18	1000113-87-1	47
4		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	43
5		Cyclododecanemethanol	198	C13H26O	001892-12-2	35



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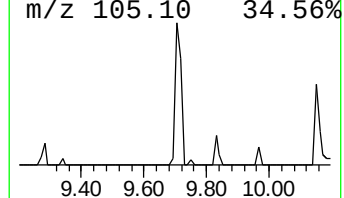
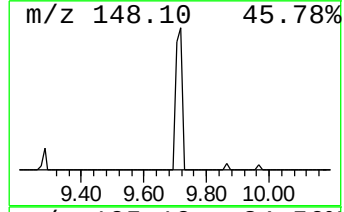
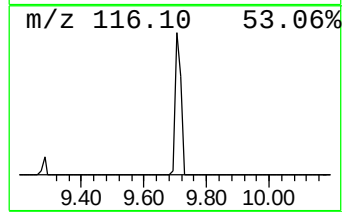
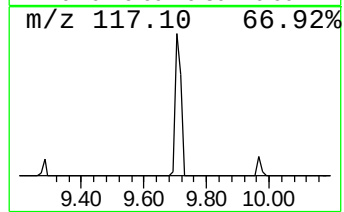
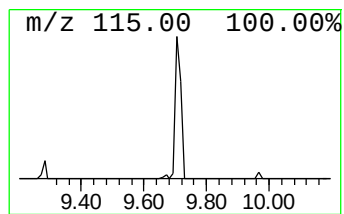
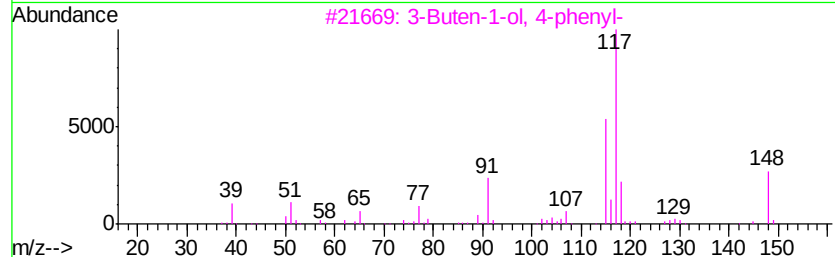
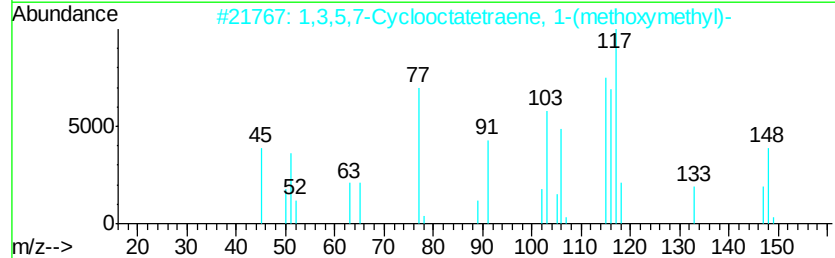
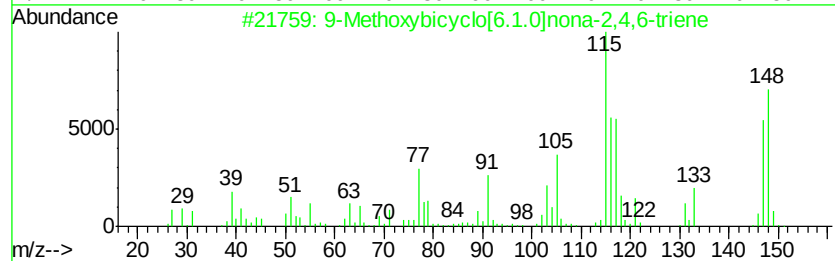
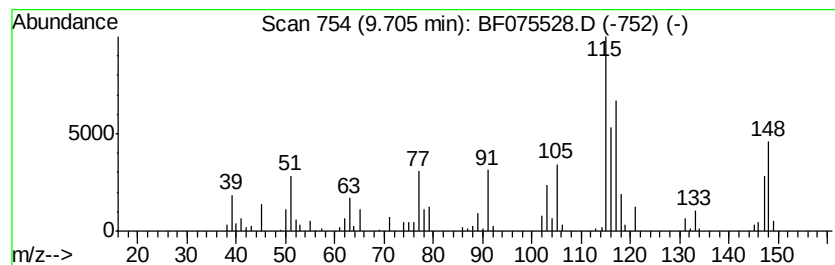
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 9-Methoxybicyclo[6.1.0]nona... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.70	5.71 ng	220612	Naphthalene-d8	9.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Methoxybicyclo[6.1.0]nona-2,4,...	148	C10H12O	000826-14-2	58
2		1,3,5,7-Cyclooctatetraene, 1-(me...	148	C10H12O	030844-13-4	52
3		3-Buten-1-ol, 4-phenyl-	148	C10H12O	000937-58-6	38
4		Benzene, 1-propynyl-	116	C9H8	000673-32-5	38
5		Benzeneacetaldehyde, .alpha.-eth...	146	C10H10O	004411-89-6	27



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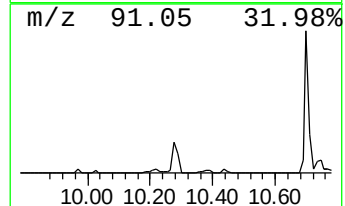
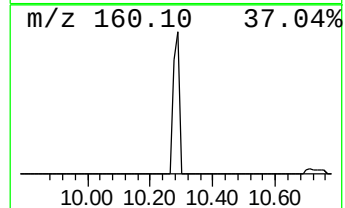
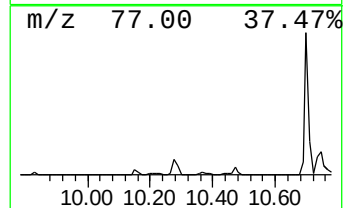
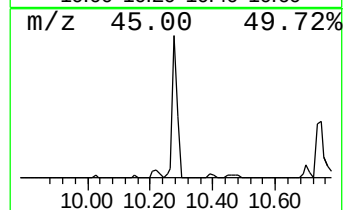
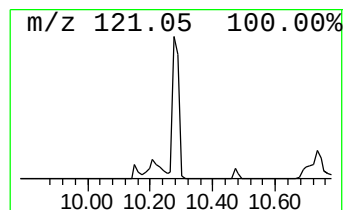
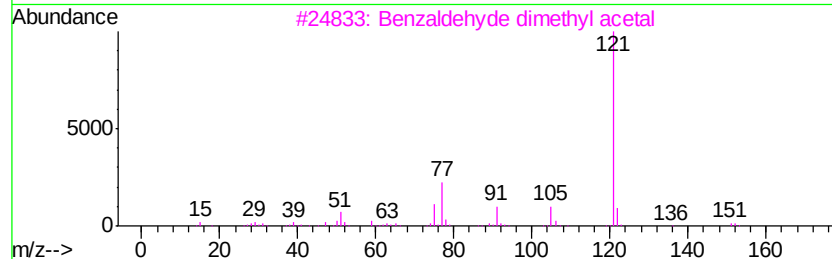
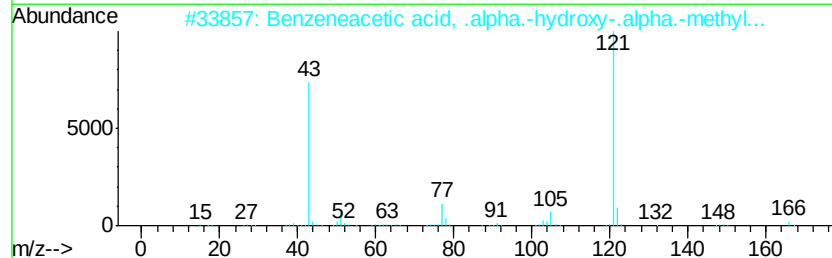
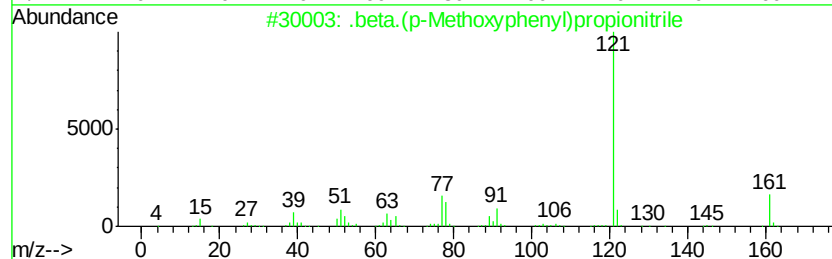
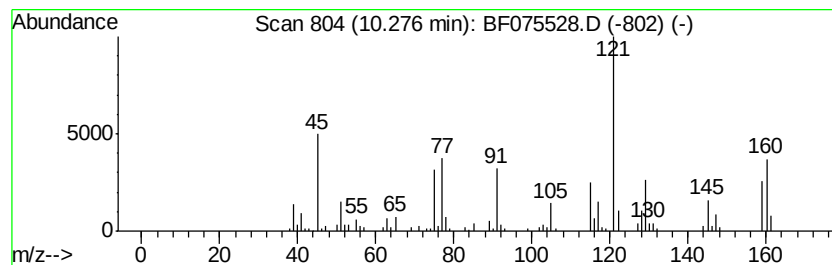
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown10.28 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.28	6.08 ng	259354	Acenaphthene-d10	11.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.(p-Methoxyphenyl)propionit...	161	C10H11NO	022442-48-4	22
2		Benzeneacetic acid, .alpha.-hydr...	166	C9H10O3	004607-38-9	22
3		Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	22
4		Phenol, 2-(1-methylpropyl)-, met...	207	C12H17NO2	003766-81-2	22
5		Pyridine, 2,4,6-trimethyl-	121	C8H11N	000108-75-8	22



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111414\
Data File : BF075528.D
Acq On : 15 Nov 2014 8:47
Operator : TP / JJ
Sample : F4695-08
Misc :
ALS Vial : 29 Sample Multiplier: 1

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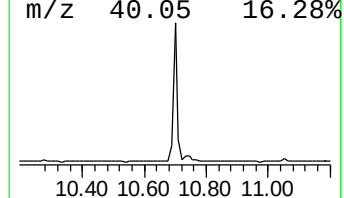
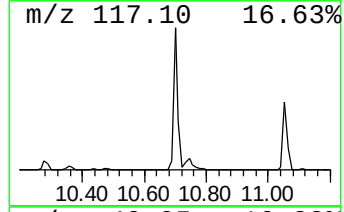
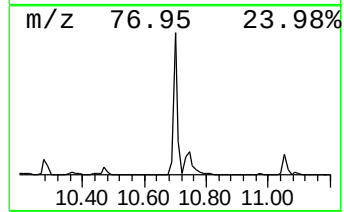
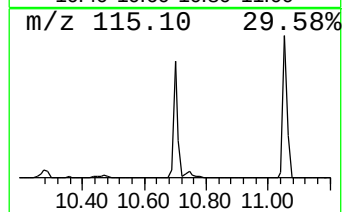
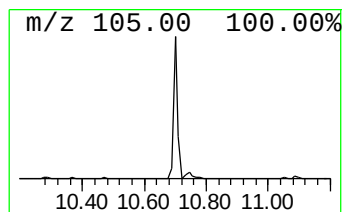
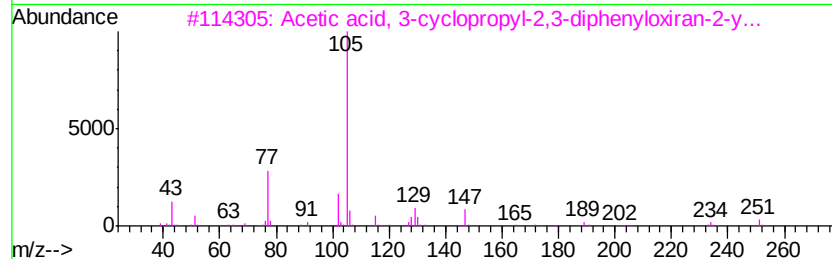
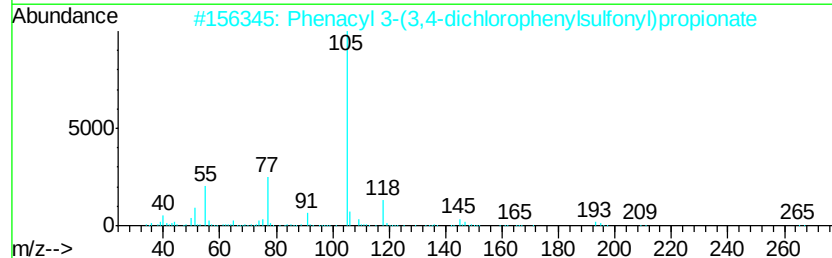
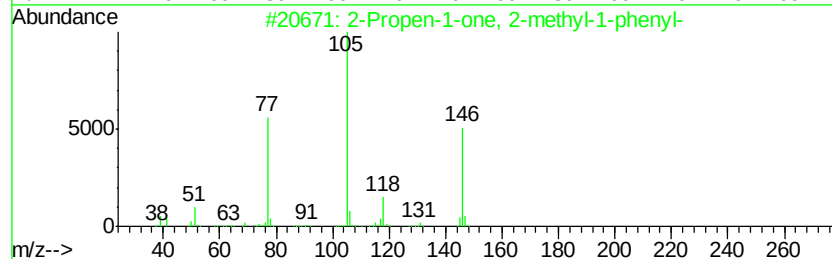
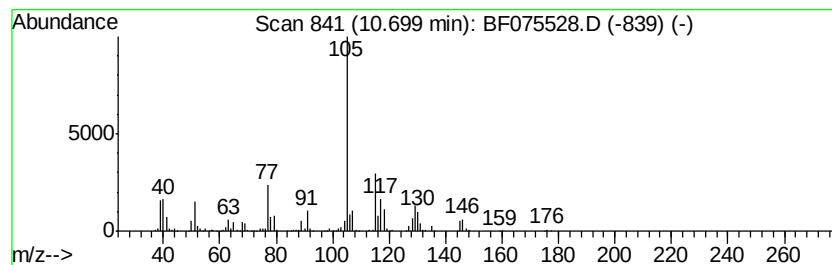
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown10.70 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.70	43.19 ng	1841630	Acenaphthene-d10	11.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propen-1-one, 2-methyl-1-phenyl-	146	C10H10O	000769-60-8	38
2		Phenacyl 3-(3,4-dichlorophenyl)sulfonylpropionate	400	C17H14Cl2O5S	1000225-01-1	35
3		Acetic acid, 3-cyclopropyl-2,3-diphenyloxiran-2-yl...	294	C19H18O3	128103-54-8	32
4		2-Phenylbutenolide	160	C10H8O2	001955-39-1	22
5		5-Amino-4-methyl-3-phenylisoxazole	174	C10H10N2O	004320-84-7	16



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111414\
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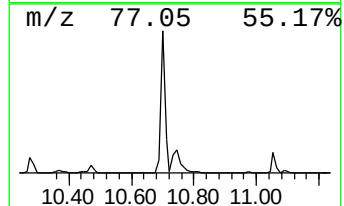
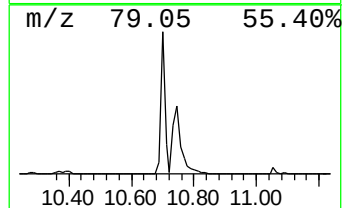
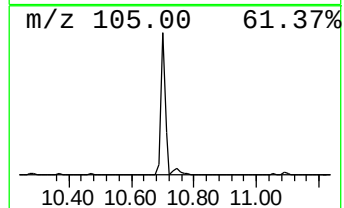
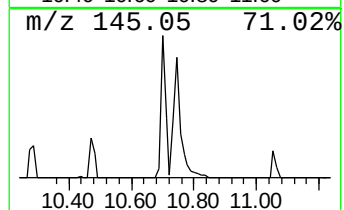
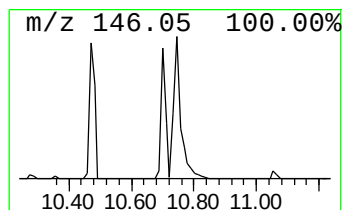
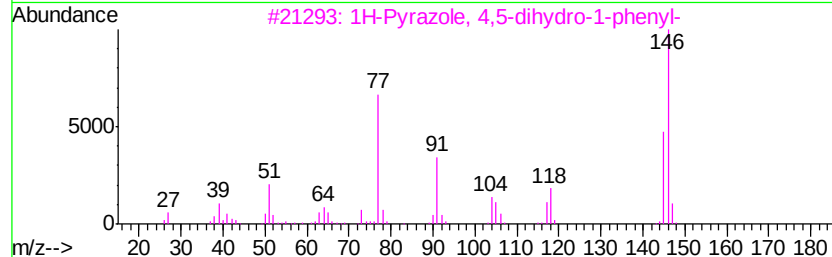
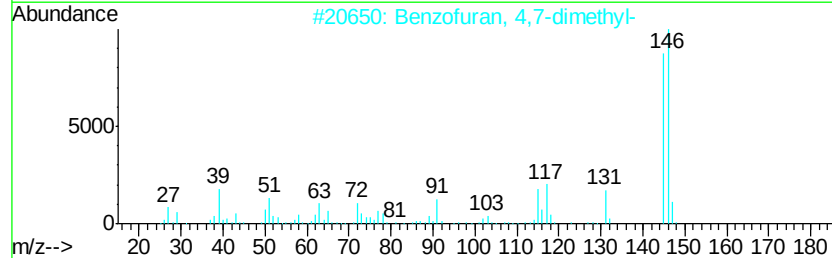
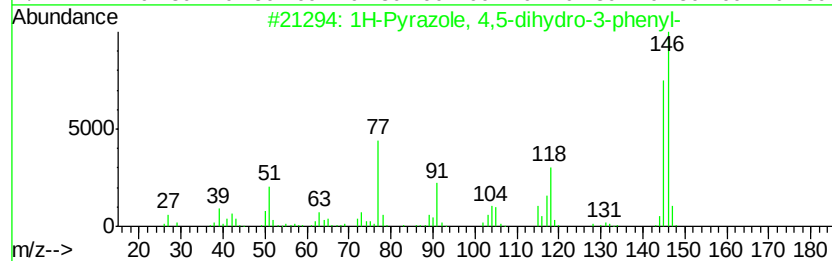
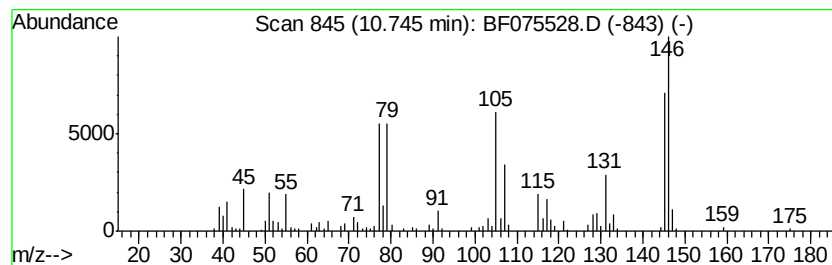
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF111214.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1H-Pyrazole, 4,5-dihydro-3-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.74	10.23 ng	436227	Acenaphthene-d10	11.31
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Pyrazole, 4,5-dihydro-3-phenyl-	146 C9H10N2	000936-48-1 64
2		Benzofuran, 4,7-dimethyl-	146 C10H10O	028715-26-6 55
3		1H-Pyrazole, 4,5-dihydro-1-phenyl-	146 C9H10N2	000936-53-8 47
4		2-Propen-1-one, 2-methyl-1-phenyl-	146 C10H10O	000769-60-8 43
5		p-Propargyloxytoluene	146 C10H10O	005651-90-1 43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111414\
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Sample : F4695-08
Misc :
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Instrument :
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ClientSampleId :
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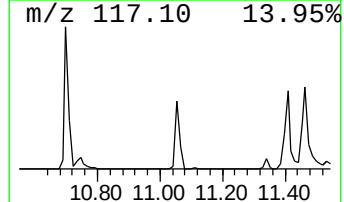
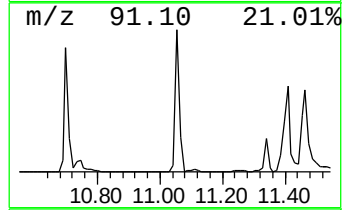
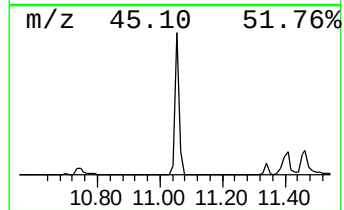
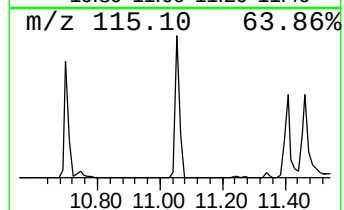
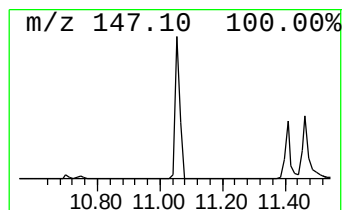
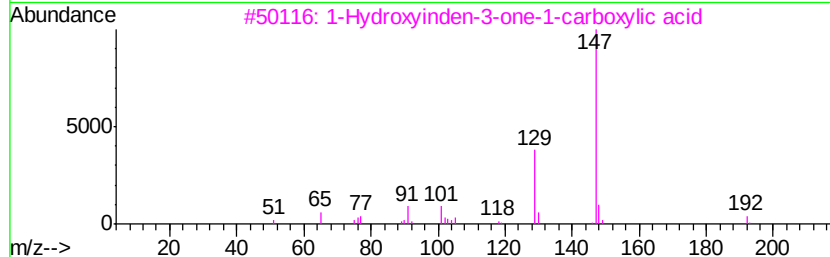
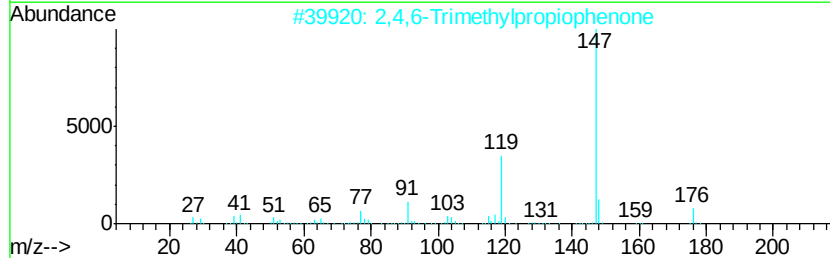
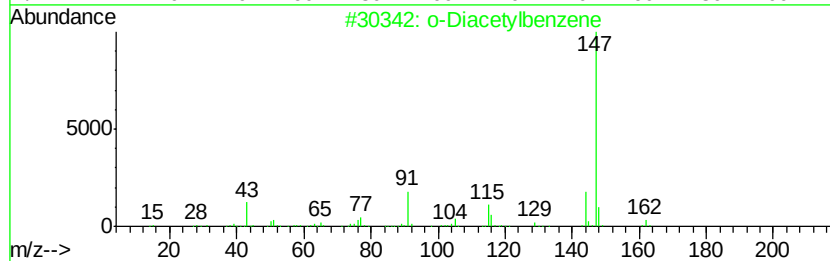
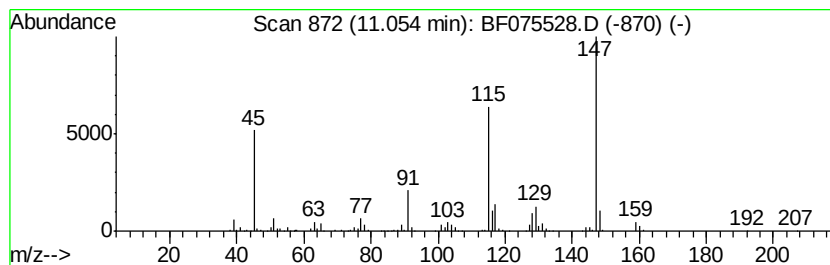
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown11.05 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.05	23.00 ng	980621	Acenaphthene-d10	11.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Diacetylbenzene	162	C10H10O2	000704-00-7	37
2		2,4,6-Trimethylpropiophenone	176	C12H16O	002040-15-5	35
3		1-Hydroxyinden-3-one-1-carboxyli...	192	C10H8O4	1000215-33-8	32
4		1(3H)-Isobenzofuranone, 3,3-dime...	162	C10H10O2	001689-09-4	27
5		Ethanone, 1-(2,4,6-trimethylphen...	162	C11H14O	001667-01-2	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111414\
Data File : BF075528.D
Acq On : 15 Nov 2014 8:47
Operator : TP / JJ
Sample : F4695-08
Misc :
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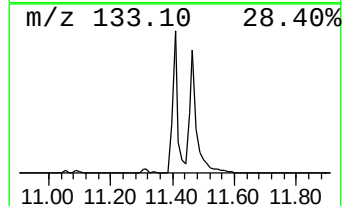
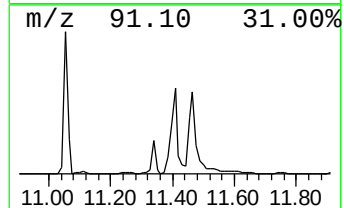
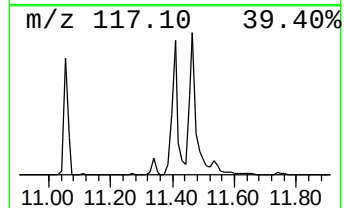
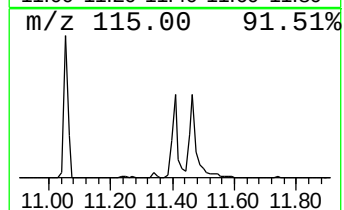
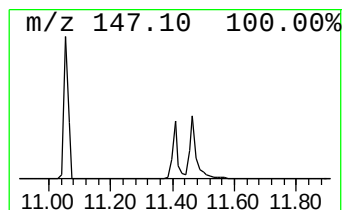
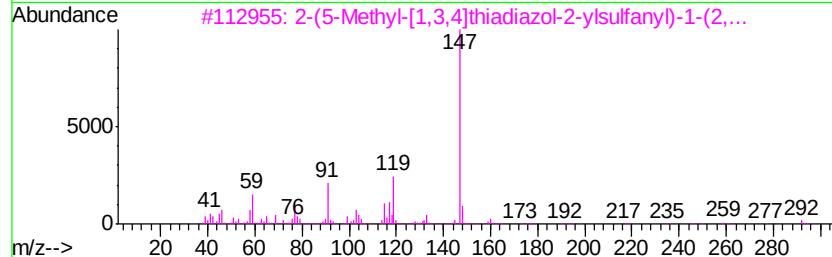
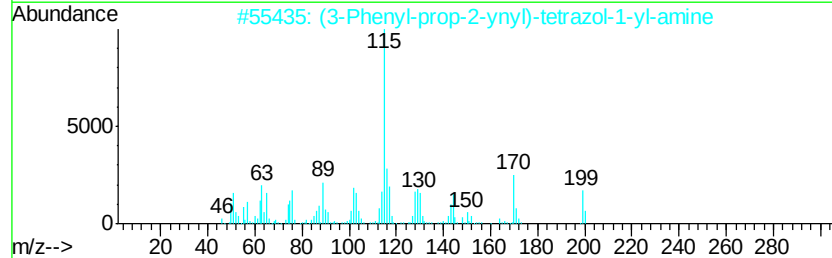
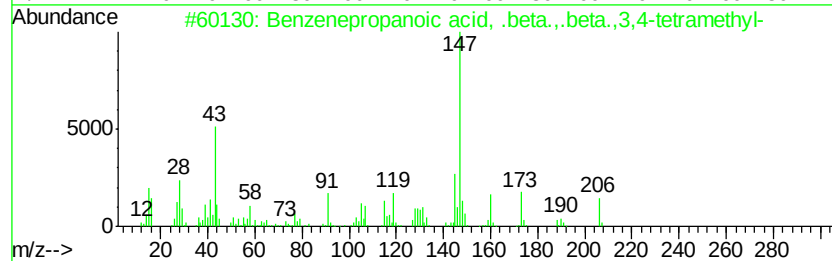
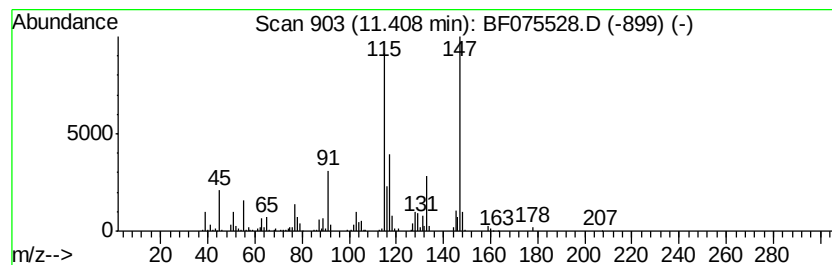
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown11.41 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.41	21.58 ng	920293	Acenaphthene-d10	11.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzenepropanoic acid, .beta.,.b...	206	C13H18O2	055683-10-8	27
2		(3-Phenyl-prop-2-ynyl)-tetrazol-...	199	C10H9N5	1000296-21-0	25
3		2-(5-Methyl-[1,3,4]thiadiazol-2-...	292	C14H16N2OS2	1000296-87-8	25
4		2H-Indazol-3-amine, 2-methyl-	147	C8H9N3	097990-19-7	22
5		Indano[2,1-d]1,3-dioxane,	176	C11H12O2	102688-70-0	16



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111414\
Data File : BF075528.D
Acq On : 15 Nov 2014 8:47
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Sample : F4695-08
Misc :
ALS Vial : 29 Sample Multiplier: 1

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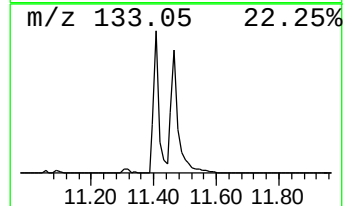
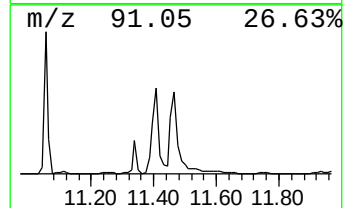
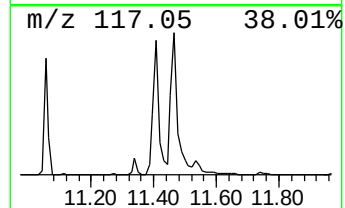
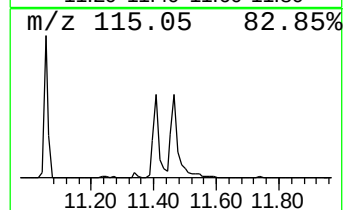
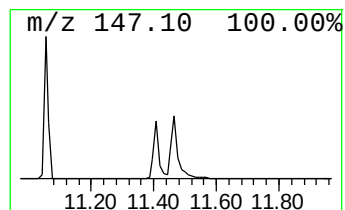
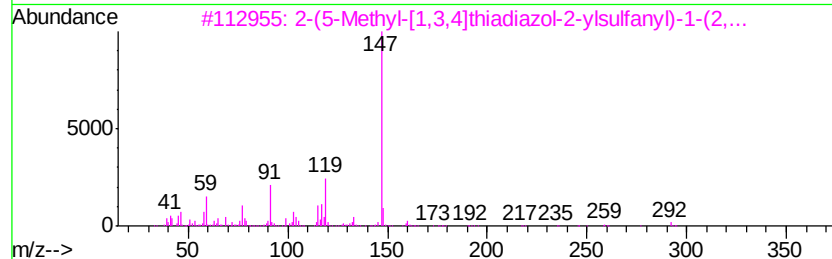
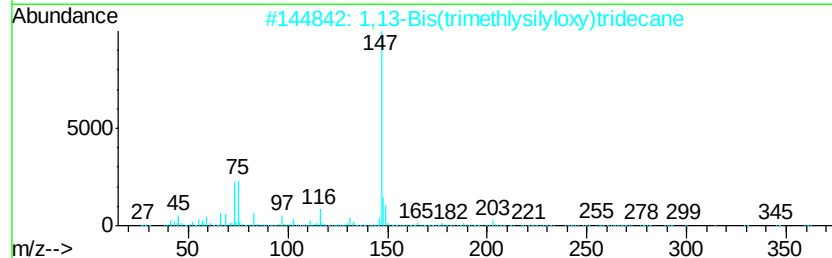
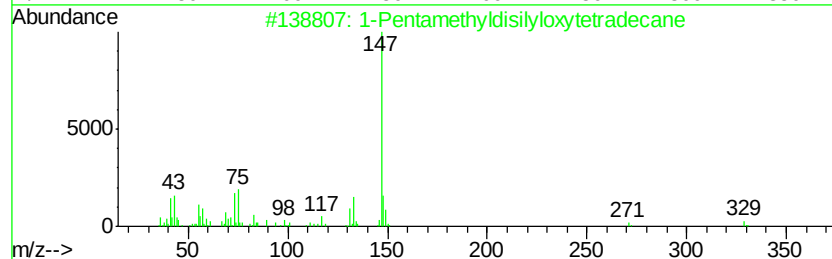
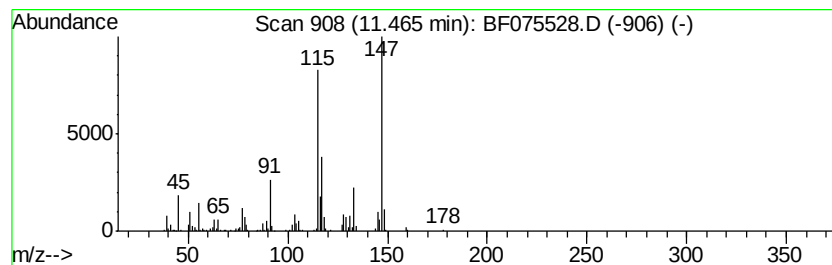
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown11.47 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.47	20.23 ng	862544	Acenaphthene-d10	11.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Pentamethylidisilvloxvtetradecane	344	C19H44OSi2	1000216-85-5	35
2		1,13-Bis(trimethylsilyloxy)tride...	360	C19H44O2Si2	1000196-23-0	35
3		2-(5-Methyl-[1,3,4]thiadiazol-2-...	292	C14H16N2OS2	1000296-87-8	32
4		Ethanone, 1-(2,3,4-trimethylphen...	162	C11H14O	001467-36-3	27
5		1-Naphthalenamine, 5,6,7,8-tetra...	147	C10H13N	002217-41-6	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111414\
Data File : BF075528.D
Acq On : 15 Nov 2014 8:47
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Sample : F4695-08
Misc :
ALS Vial : 29 Sample Multiplier: 1

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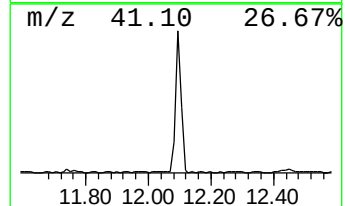
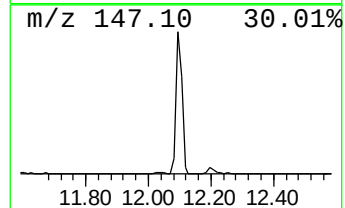
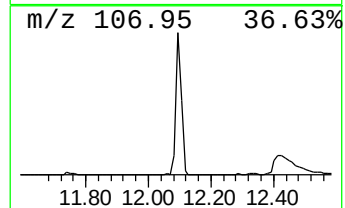
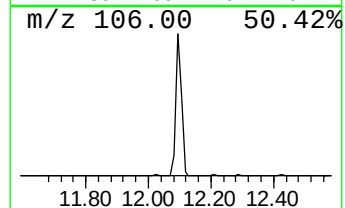
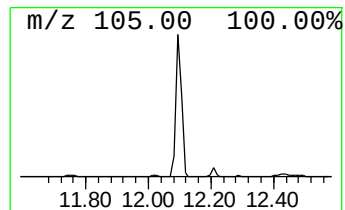
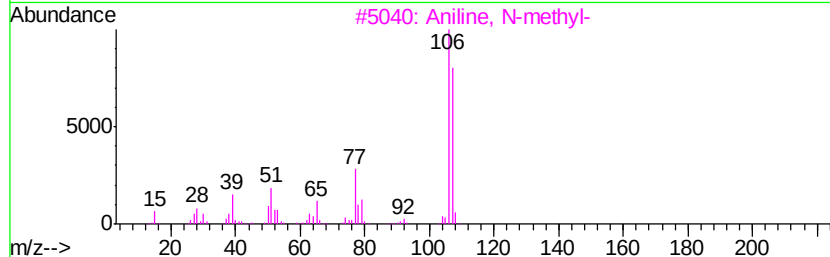
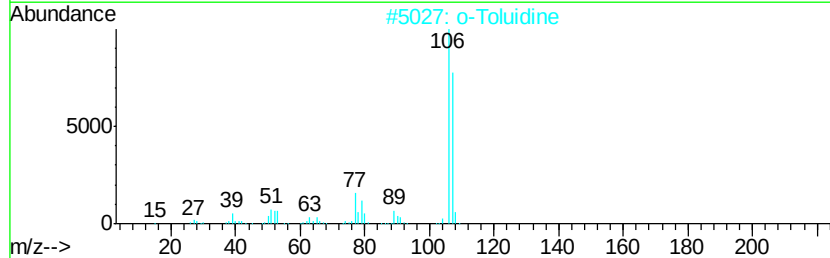
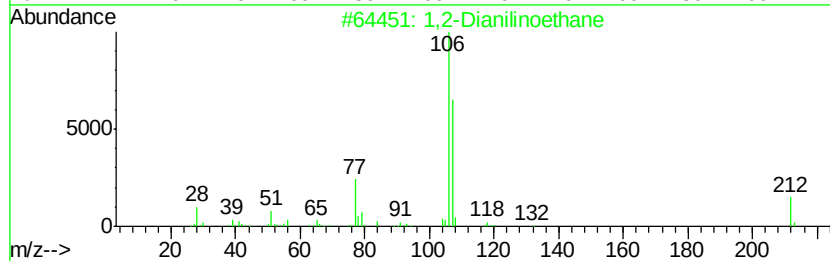
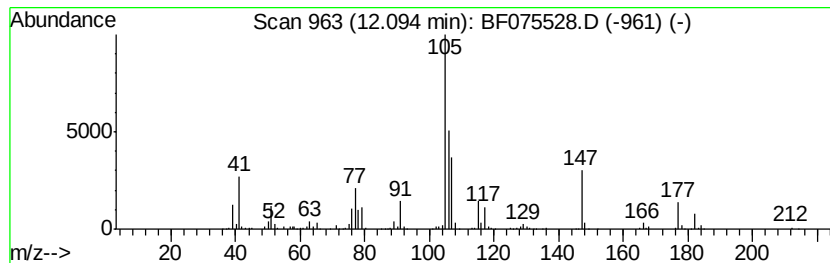
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown12.09 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.09	27.52 ng	1173660	Acenaphthene-d10	11.31

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2-Dianilinoethane	212	C14H16N2	000150-61-8	45
2		o-Toluidine	107	C7H9N	000095-53-4	43
3		Aniline, N-methyl-	107	C7H9N	000100-61-8	43
4		Aminodiphenylmethane	183	C13H13N	000091-00-9	38
5		Benzylamine	107	C7H9N	000100-46-9	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF111414\
Data File : BF075528.D
Acq On : 15 Nov 2014 8:47
Operator : TP / JJ
Sample : F4695-08
Misc :
ALS Vial : 29 Sample Multiplier: 1

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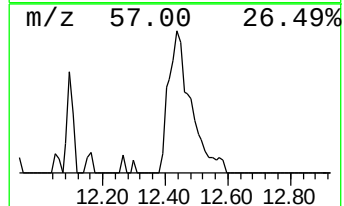
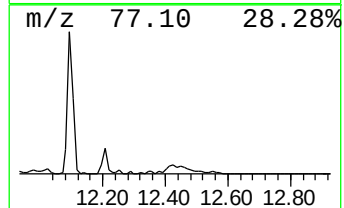
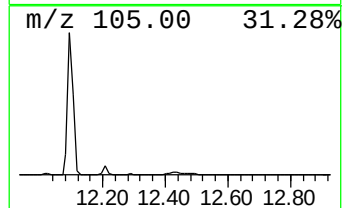
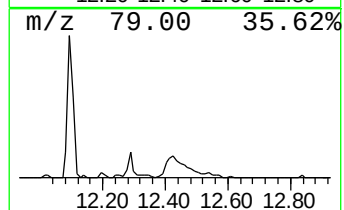
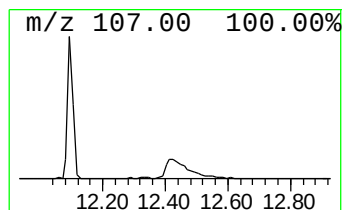
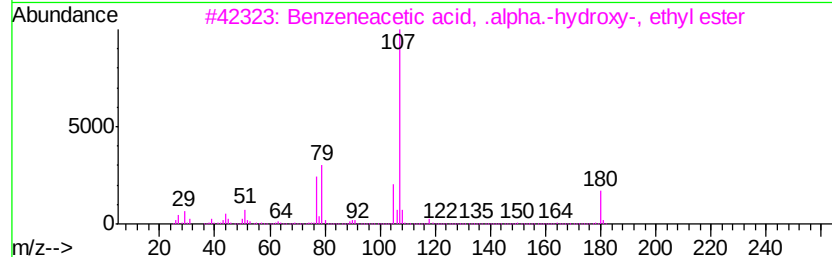
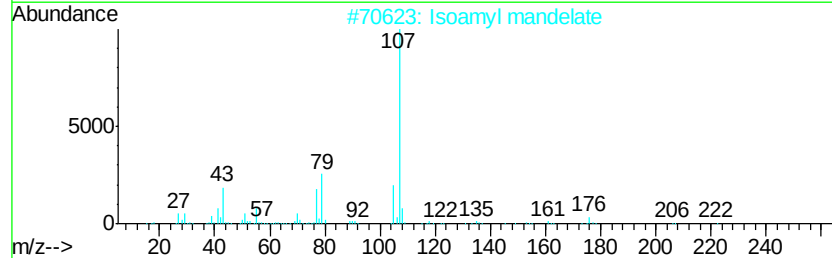
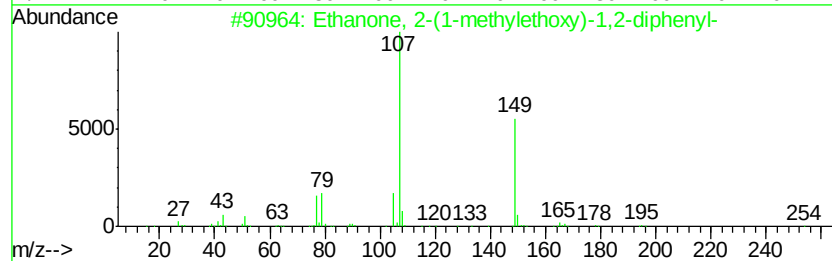
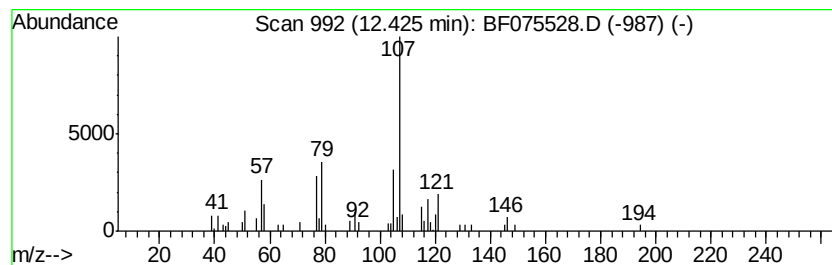
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Ethanone, 2-(1-methylethoxy... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.43	4.25 ng	180316	Phenanthrene-d10	13.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanone, 2-(1-methylethoxy)-1,2...	254	C17H18O2	006652-28-4	59
2		Isoamyl mandelate	222	C13H18O3	005421-04-5	59
3		Benzeneacetic acid, .alpha.-hydr...	180	C10H12O3	000774-40-3	59
4		Cyclohexene, 1-ethyl-6-ethylidene-	136	C10H16	061141-57-9	53
5		Benzeneacetic acid, .alpha.-hydr...	152	C8H8O3	017199-29-0	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111414\
Data File : BF075528.D
Acq On : 15 Nov 2014 8:47
Operator : TP / JJ
Sample : F4695-08
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FB111014

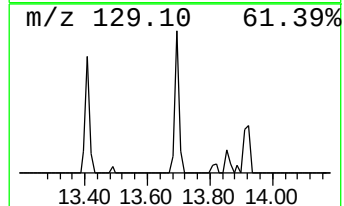
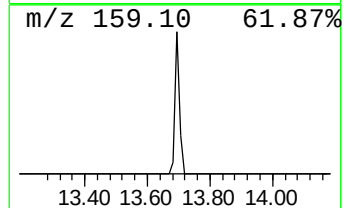
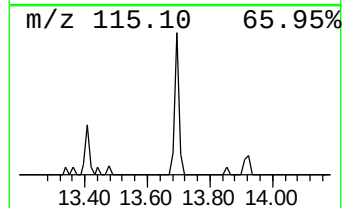
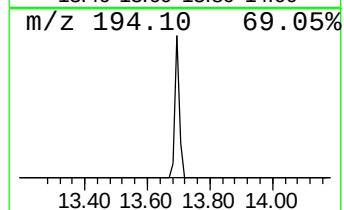
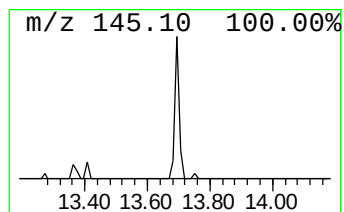
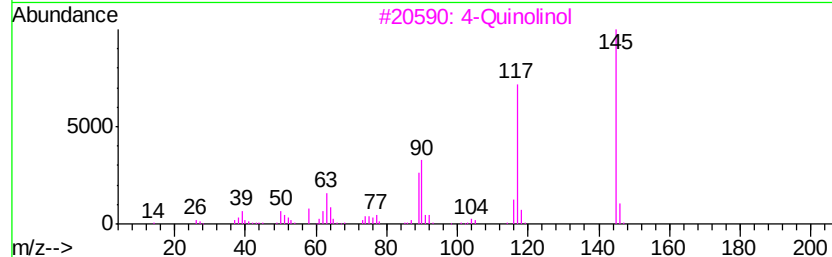
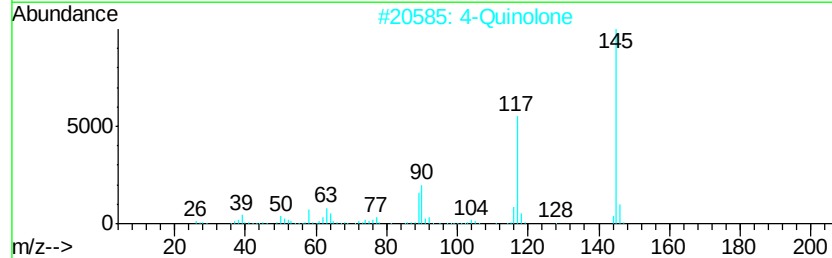
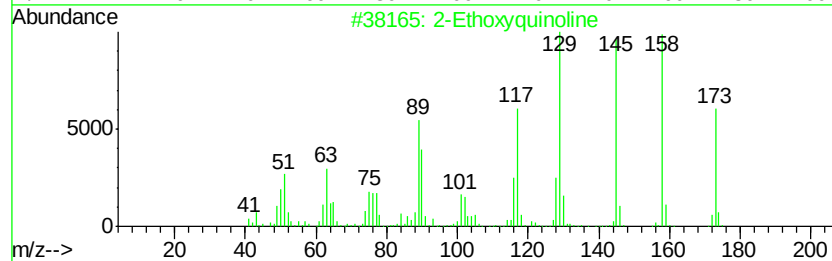
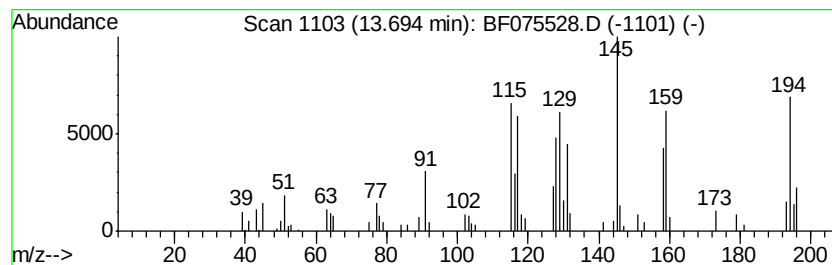
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown13.69 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.69	2.03 ng	86206	Phenanthrene-d10	13.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Ethoxyquinoline	173	C11H11NO	046185-83-5	25
2		4-Quinolone	145	C9H7NO	1000234-25-1	20
3		4-Quinololinol	145	C9H7NO	000611-36-9	18
4		2(1H)-Quinololinone	145	C9H7NO	000059-31-4	14
5		3-Benzyl-1H-1,2,4-triazole	159	C9H9N3	021117-34-0	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111414\
 Data File : BF075528.D
 Acq On : 15 Nov 2014 8:47
 Operator : TP / JJ
 Sample : F4695-08
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FB111014

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF111214.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...	5.34	7.5	ng	201698	1	7.54	540600	20.0
unknown7.26	7.26	52.5	ng	1417900	1	7.54	540600	20.0
2-Aziridinone, 1-...	9.34	2.2	ng	86225	2	9.13	772861	20.0
9-Methoxybicyclo[...	9.70	5.7	ng	220612	2	9.13	772861	20.0
unknown10.28	10.28	6.1	ng	259354	3	11.31	852825	20.0
unknown10.70	10.70	43.2	ng	1841630	3	11.31	852825	20.0
1H-Pyrazole, 4,5-...	10.74	10.2	ng	436227	3	11.31	852825	20.0
unknown11.05	11.05	23.0	ng	980621	3	11.31	852825	20.0
unknown11.41	11.41	21.6	ng	920293	3	11.31	852825	20.0
unknown11.47	11.47	20.2	ng	862544	3	11.31	852825	20.0
unknown12.09	12.09	27.5	ng	1173660	3	11.31	852825	20.0
Ethanone, 2-(1-me...	12.43	4.3	ng	180316	4	13.16	849420	20.0
unknown13.69	13.69	2.0	ng	86206	4	13.16	849420	20.0