

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051316\
 Data File : BF086999.D
 Acq On : 14 May 2016 5:07
 Operator : UM/SJ
 Sample : H2966-05
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-36

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-BF050916.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	1.735	12	13	50	rVB	4946996	10250816	100.00%	22.366%
2	5.199	314	316	328	rBV	1569565	1800158	17.56%	3.928%
3	6.273	408	410	417	rBV	1074657	1161085	11.33%	2.533%
4	6.387	417	420	422	rBV	2658534	3414521	33.31%	7.450%
5	6.604	437	439	442	rBV	651844	919541	8.97%	2.006%
6	6.764	451	453	456	rBV	3070599	3013457	29.40%	6.575%
7	7.187	488	490	491	rVV	2948057	3009114	29.35%	6.566%
8	7.210	491	492	494	rVV	514385	490921	4.79%	1.071%
9	7.370	503	506	510	rVB3	65964	130023	1.27%	0.284%
10	7.587	520	525	527	rVB2	53200	109135	1.06%	0.238%
11	7.644	527	530	532	rBV	139874	156396	1.53%	0.341%
12	7.702	532	535	540	rBV5	64084	120637	1.18%	0.263%
13	7.839	544	547	550	rBV2	61376	122410	1.19%	0.267%
14	7.896	550	552	556	rBV	1149373	1287510	12.56%	2.809%
15	8.056	563	566	567	rBV2	113109	112831	1.10%	0.246%
16	8.662	616	619	622	rBV4	130820	307617	3.00%	0.671%
17	8.982	644	647	649	rBV	4755010	4650341	45.37%	10.146%
18	9.153	661	662	665	rVB3	89985	107060	1.04%	0.234%
19	9.267	667	672	673	rBV5	36620	119107	1.16%	0.260%
20	9.347	676	679	681	rVV4	83846	170430	1.66%	0.372%
21	9.382	681	682	688	rVV3	69967	113662	1.11%	0.248%
22	9.462	688	689	691	rVB2	105217	115666	1.13%	0.252%
23	9.576	697	699	702	rVB2	265838	301262	2.94%	0.657%
24	9.645	703	705	707	rBV	1329571	1272758	12.42%	2.777%
25	9.976	731	734	736	rBV4	124124	259553	2.53%	0.566%
26	10.010	736	737	740	rVV3	130257	169856	1.66%	0.371%
27	10.090	742	744	745	rVB2	119363	128549	1.25%	0.280%
28	10.159	746	750	752	rBV2	105322	164213	1.60%	0.358%
29	10.353	765	767	769	rVB	100445	103764	1.01%	0.226%
30	10.399	769	771	772	rVV	423860	450419	4.39%	0.983%
31	10.433	772	774	777	rVV2	2340366	2891267	28.21%	6.308%
32	10.490	777	779	782	rVV2	181110	231915	2.26%	0.506%
33	10.548	782	784	787	rVV2	194750	329561	3.21%	0.719%
34	10.982	820	822	826	rBV4	159661	206098	2.01%	0.450%

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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 >
Peak separation: 5

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35	11.119	832	834	837	rBV2	971985	1168570	11.40%	2.550%
36	11.199	840	841	844	rVB	125706	168303	1.64%	0.367%
37	11.679	881	883	888	rVV2	202310	258084	2.52%	0.563%
38	11.793	891	893	899	rVB3	164456	236440	2.31%	0.516%
39	12.456	949	951	953	rBV	105378	111928	1.09%	0.244%
40	12.708	970	973	976	rBV	3342838	3899558	38.04%	8.508%
41	13.611	1050	1052	1058	rBV	78637	120783	1.18%	0.264%
42	13.748	1061	1064	1067	rBV	1045000	966212	9.43%	2.108%
43	15.108	1180	1183	1186	rBV	629679	710530	6.93%	1.550%

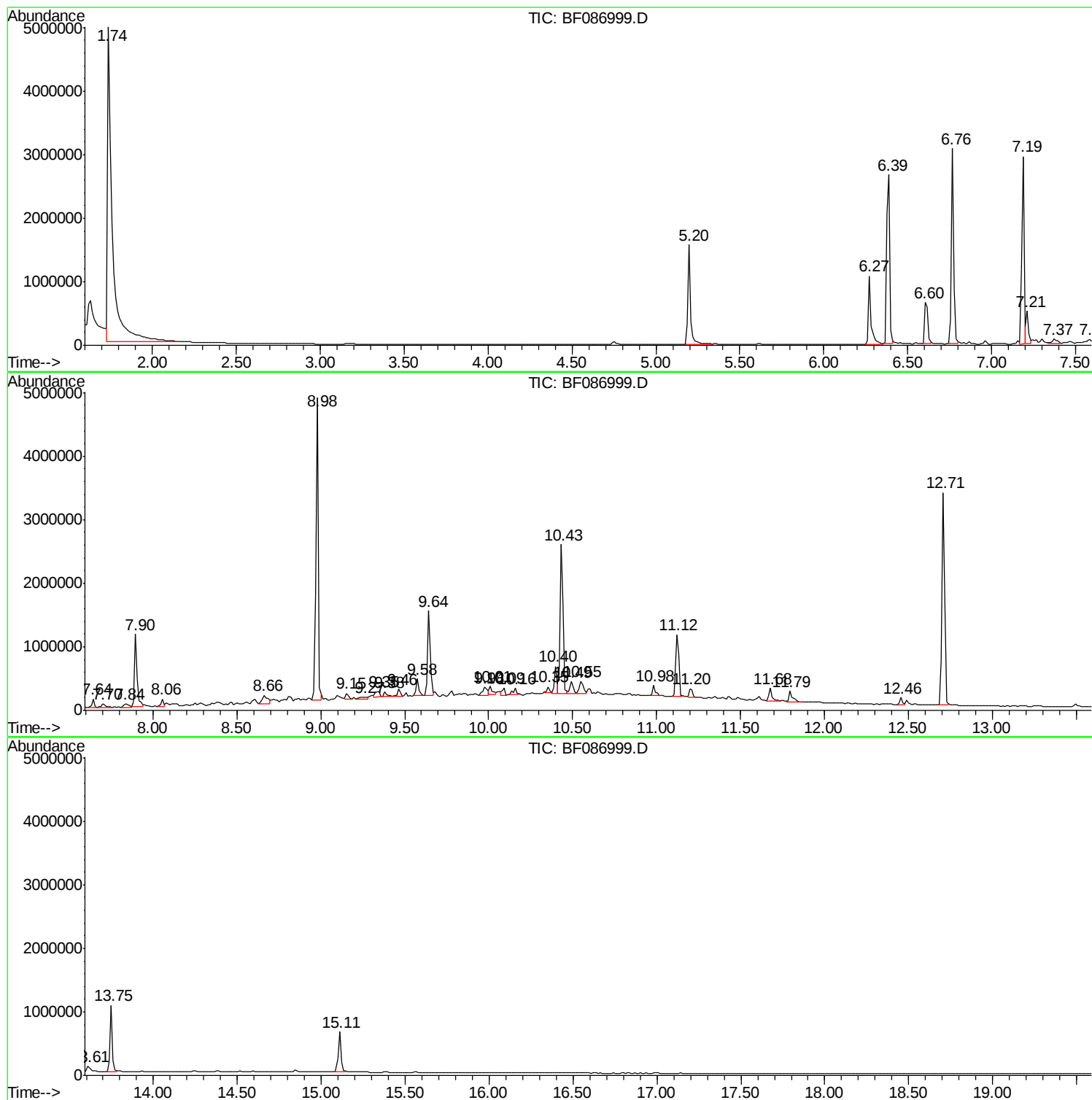
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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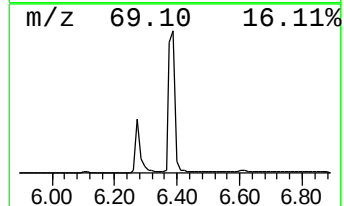
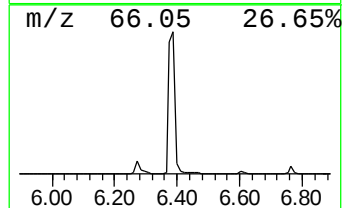
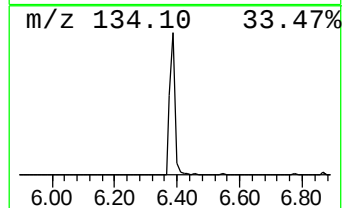
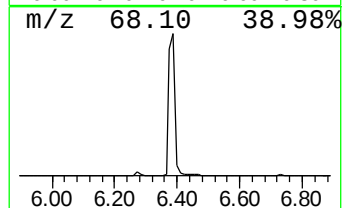
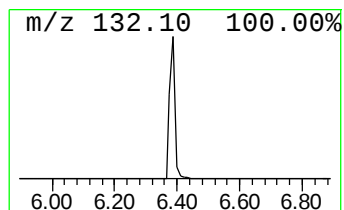
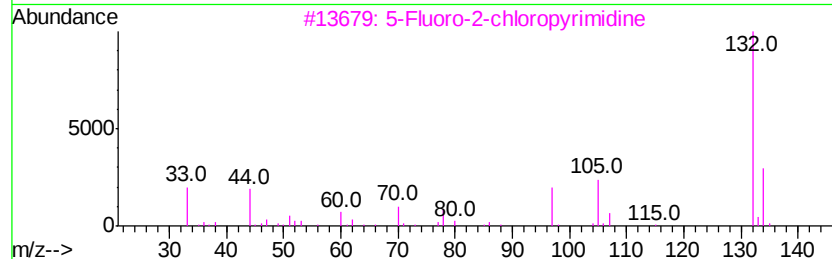
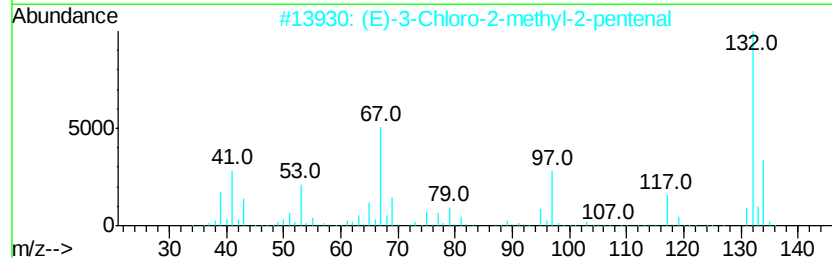
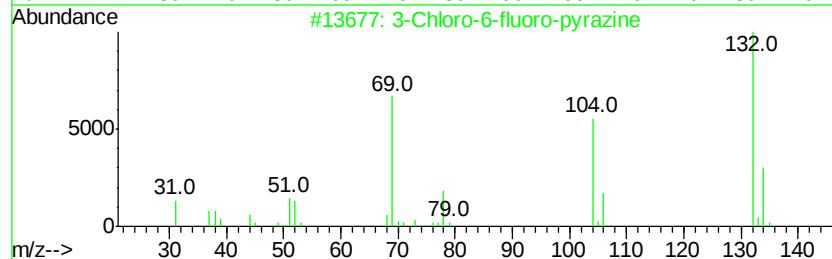
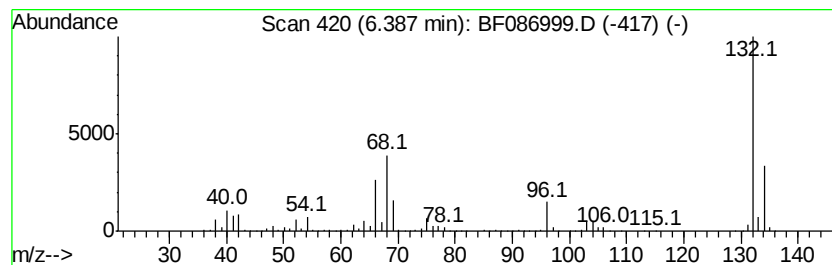
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.39 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.39	74.27 ng	3414520	1,4-Dichlorobenzene-d4	6.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9



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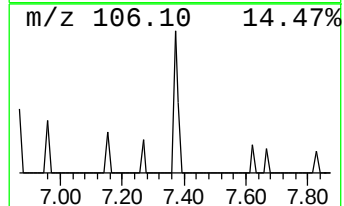
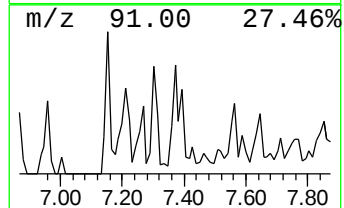
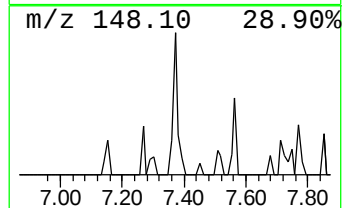
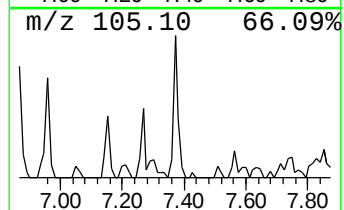
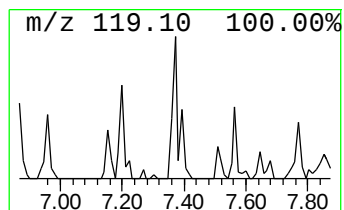
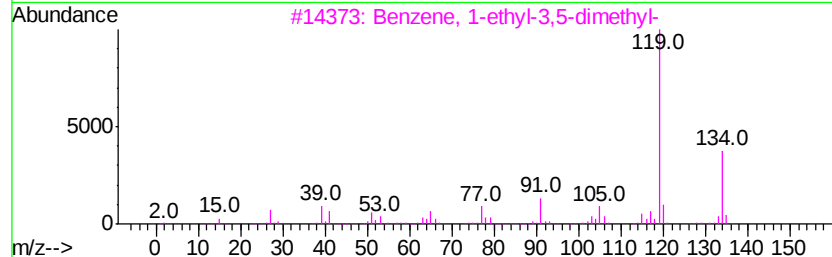
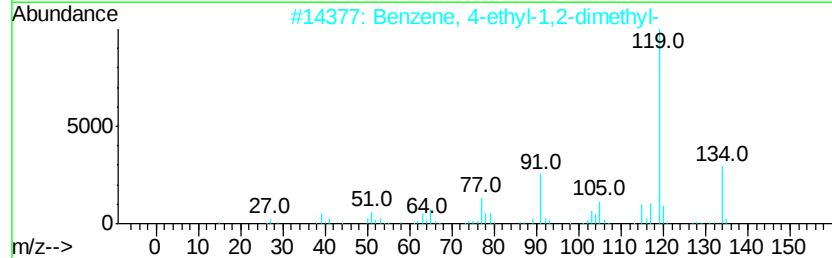
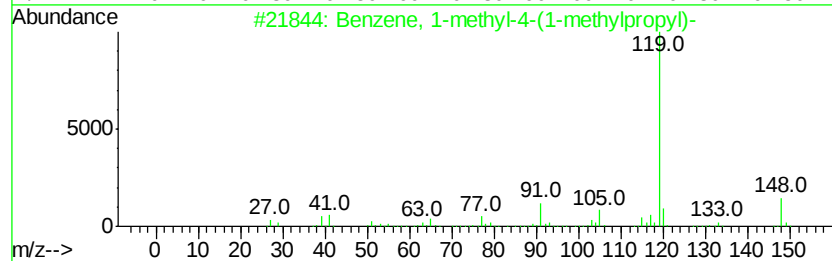
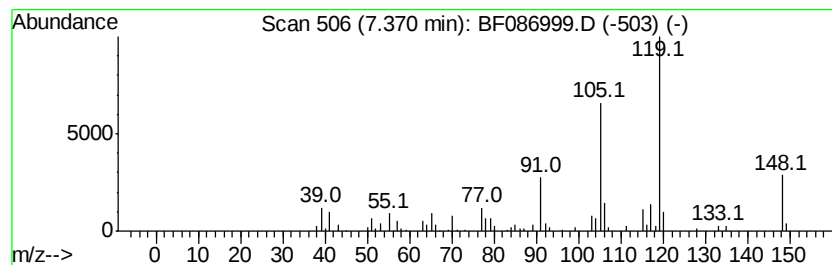
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 Benzene, 1-methyl-4-(1-meth... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.37	2.02 ng	130023	Napthalene-d8	7.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	83
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	72
3		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	72
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	72
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	72



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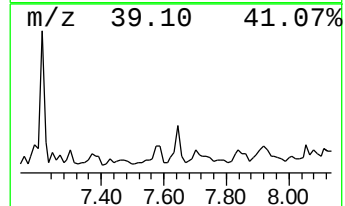
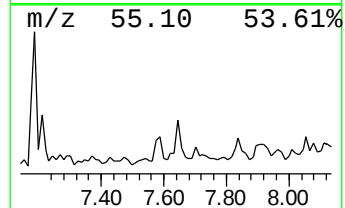
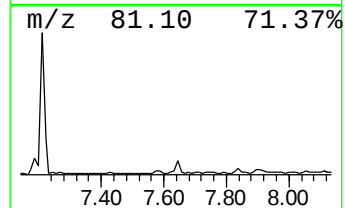
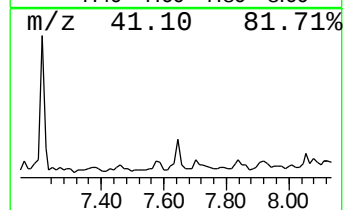
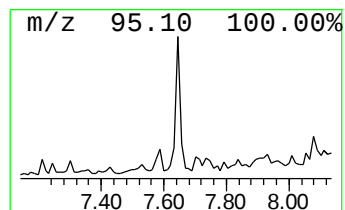
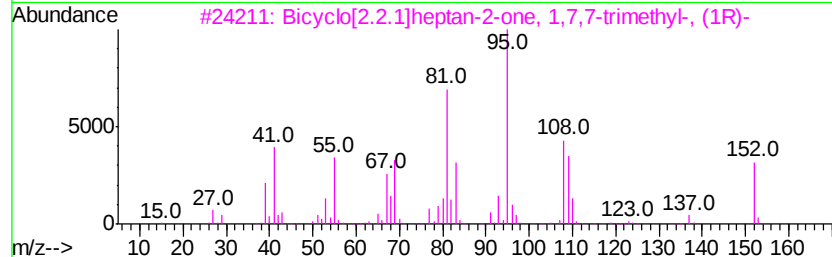
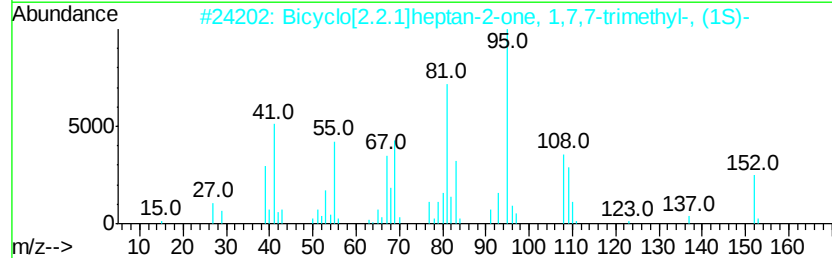
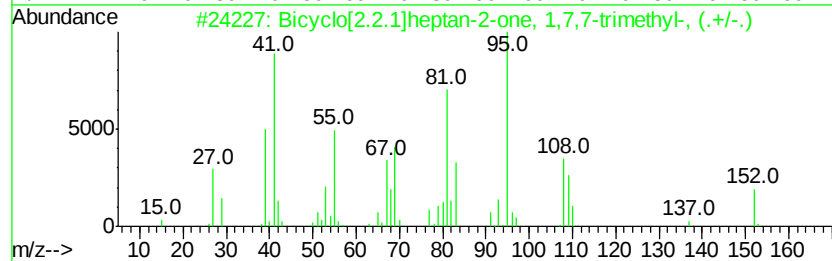
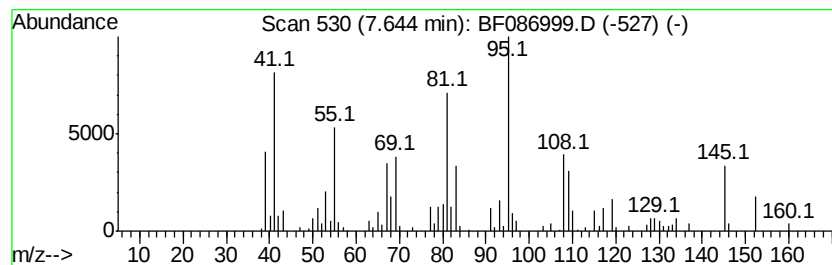
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.64	2.43 ng	156396	Naphthalene-d8	7.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	021368-68-3	98
2		Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	97
3		Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-49-3	97
4		Camphor	152	C10H16O	000076-22-2	93
5		5,7-Octadien-4-one, 2,6-dimethyl...	152	C10H16O	003588-18-9	38



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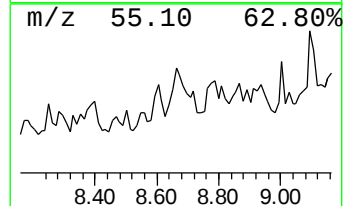
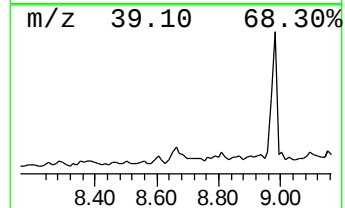
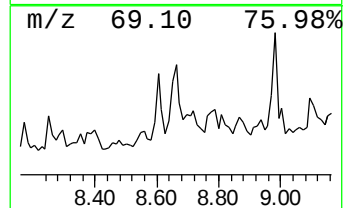
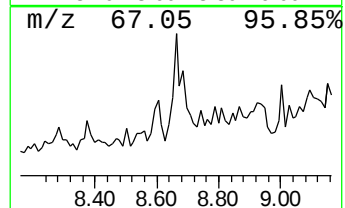
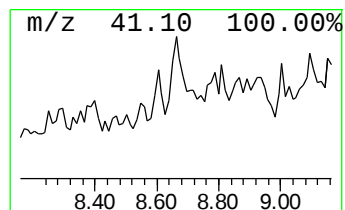
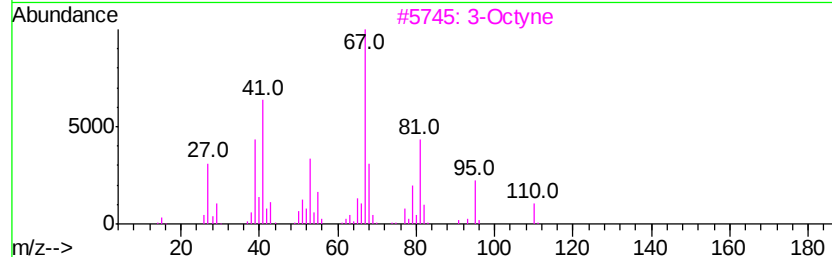
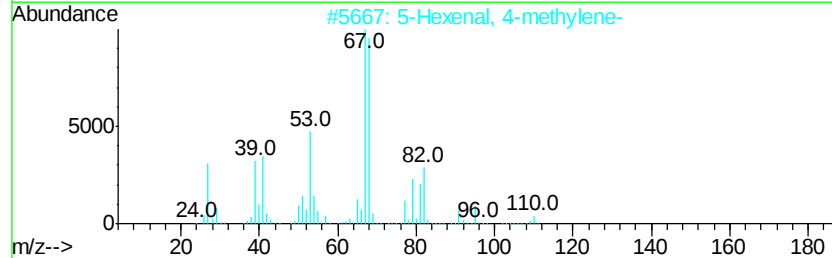
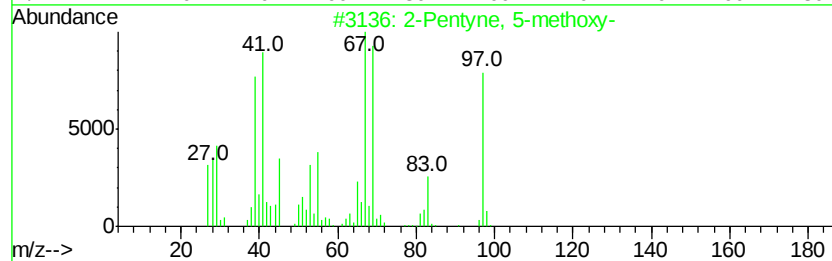
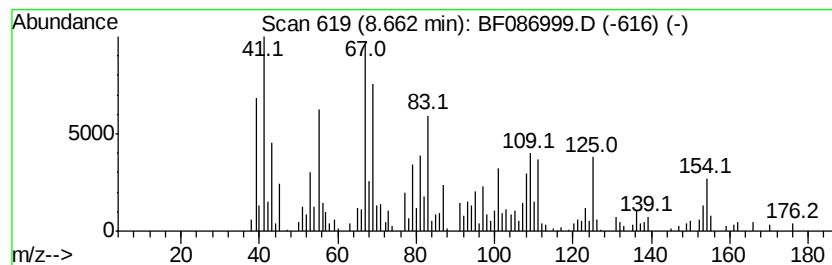
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Peak Number 6 unknown8.66 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.66	4.78 ng	307617	Naphthalene-d8	7.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentyne, 5-methoxy-	98	C6H10O	062108-18-3	41
2		5-Hexenal, 4-methylene-	110	C7H10O	017844-21-2	25
3		3-Octyne	110	C8H14	015232-76-5	25
4		4-Octene, 2,3,7-trimethyl-, [S-(...]	154	C11H22	052763-13-0	25
5		Bicyclo[3.1.1]heptan-3-one, 2,6,...	152	C10H16O	015358-88-0	25



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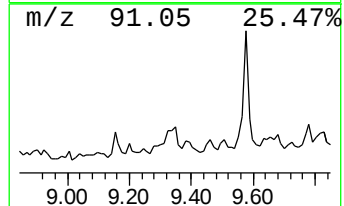
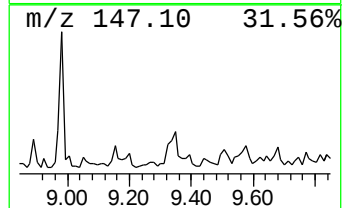
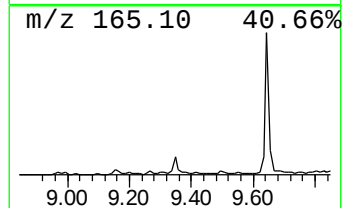
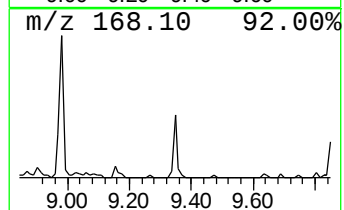
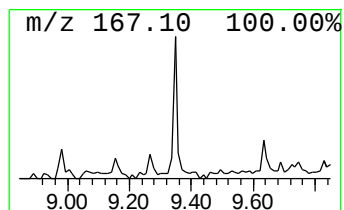
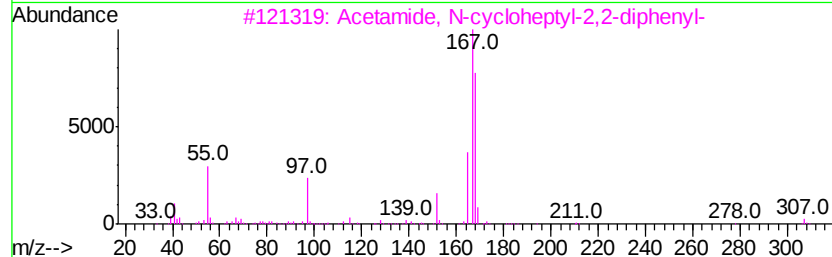
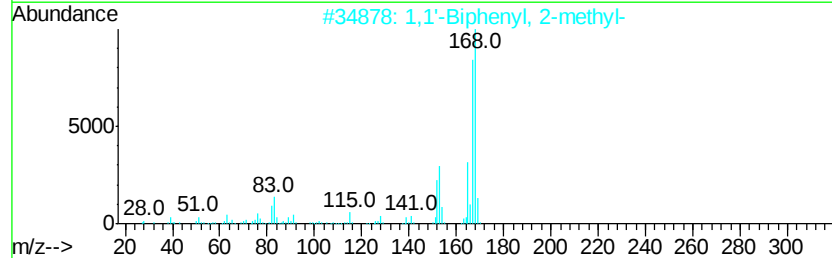
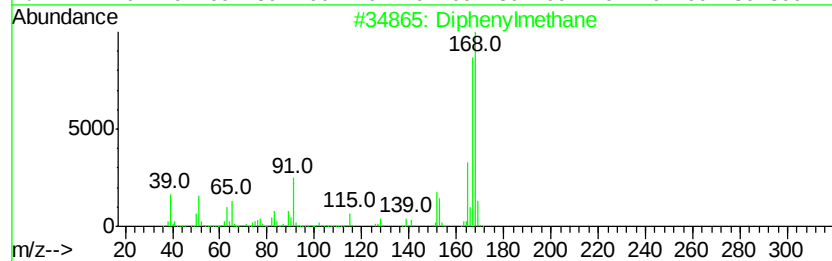
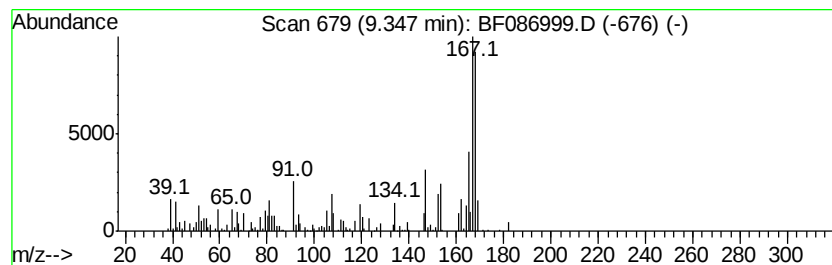
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Diphenylmethane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.35	2.68 ng	170430	Acenaphthene-d10	9.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diphenylmethane	168	C13H12	000101-81-5	64
2		1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	62
3		Acetamide, N-cycloheptyl-2,2-diphenyl-	307	C21H25NO	351062-01-6	53
4		Spiro(tricyclo[6.2.1.0(2,7)]undec-6-ene-2,2'-biphenyl)	168	C13H12	022003-58-3	52
5		Nordiphenamid	225	C15H15NO	000954-21-2	50



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051316\
Data File : BF086999.D
Acq On : 14 May 2016 5:07
Operator : UM/SJ
Sample : H2966-05
Misc :
ALS Vial : 25 Sample Multiplier: 1

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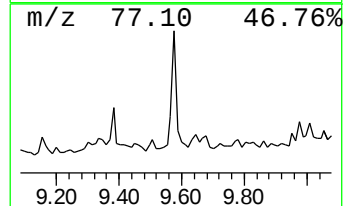
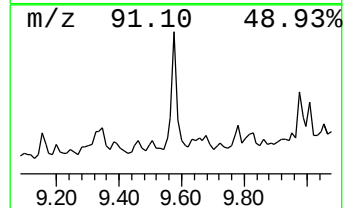
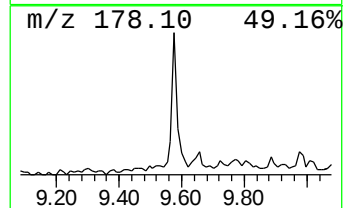
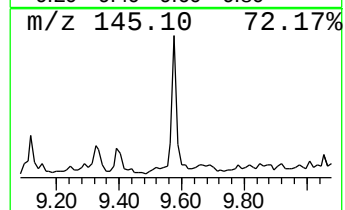
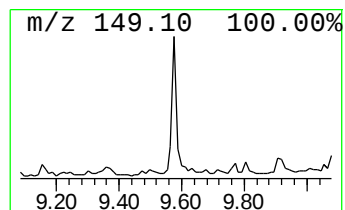
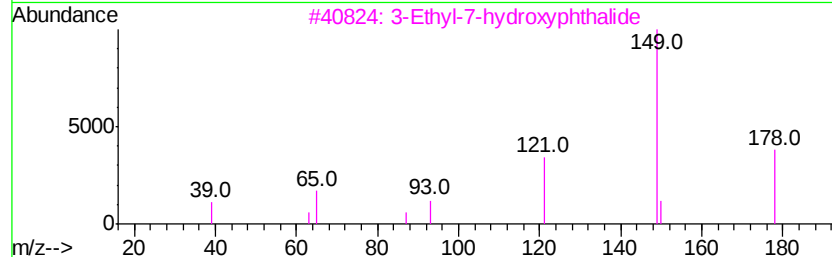
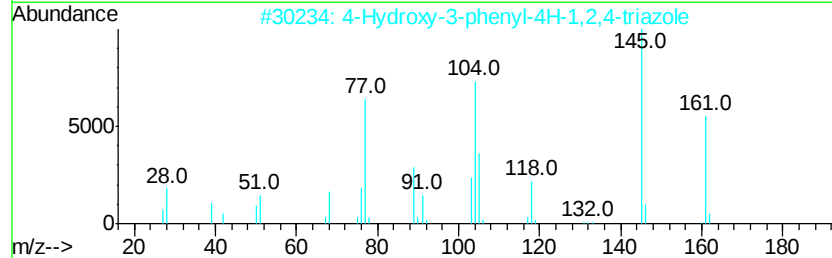
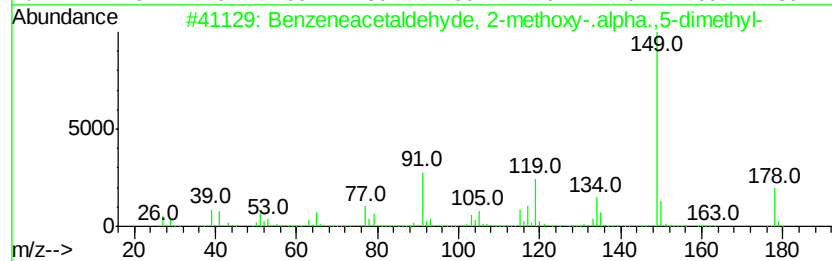
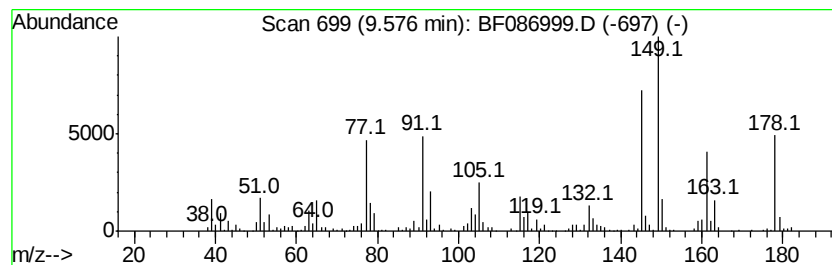
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Benzeneacetaldehyde, 2-meth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.58	4.73 ng	301262	Acenaphthene-d10	9.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzeneacetaldehyde, 2-methoxy-....	178	C11H14O2	053155-90-1	53
2		4-Hydroxy-3-phenyl-4H-1,2,4-tria...	161	C8H7N3O	031409-47-9	38
3		3-Ethyl-7-hydroxyphthalide	178	C10H10O3	000485-26-7	27
4		2-Acetylbenzoic acid	164	C9H8O3	000577-56-0	25
5		2-(1-Cyclohexenyl)cyclohexanone	178	C12H18O	001502-22-3	16



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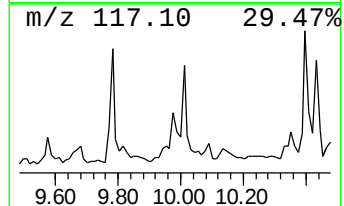
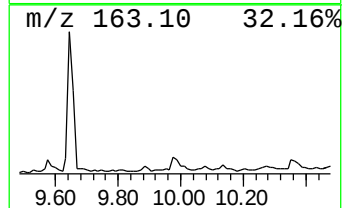
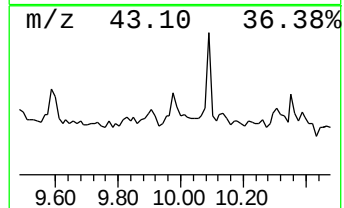
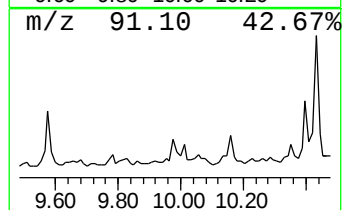
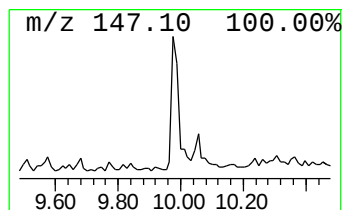
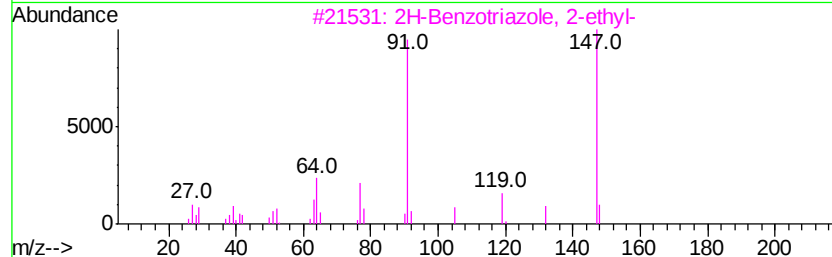
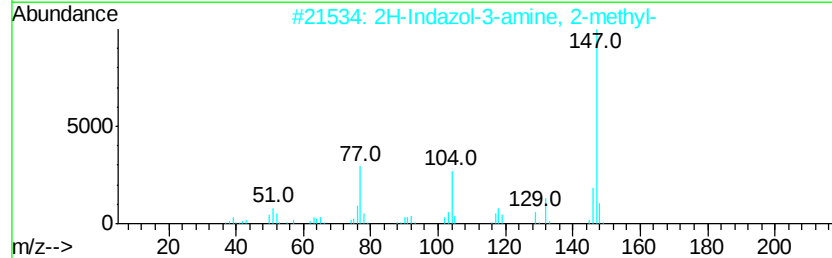
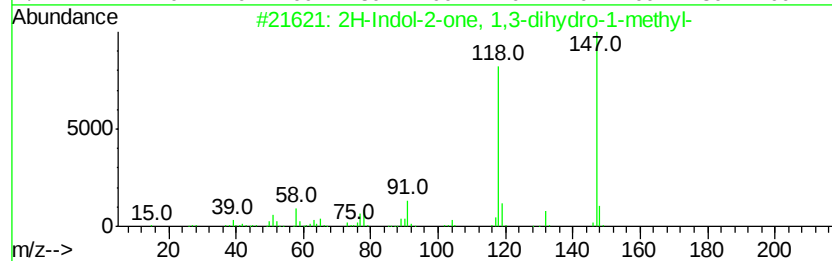
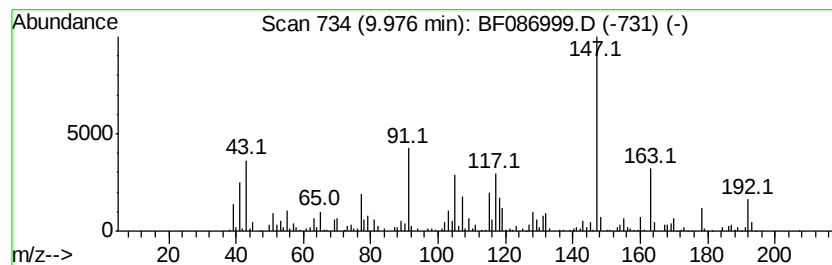
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown9.98 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.98	4.08 ng	259553	Acenaphthene-d10	9.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-Indol-2-one, 1,3-dihydro-1-methyl-	147	C9H9NO	000061-70-1	46
2		2H-Indazol-3-amine, 2-methyl-	147	C8H9N3	097990-19-7	43
3		2H-Benzotriazole, 2-ethyl-	147	C8H9N3	016584-04-6	43
4		2-Carboxycinnamic acid	192	C10H8O4	000612-40-8	43
5		Benzene, 1,4-bis(1-methylethyl)-	162	C12H18	000100-18-5	38



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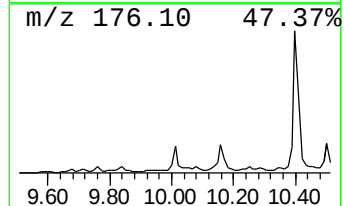
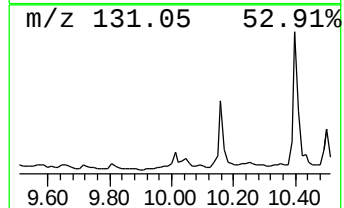
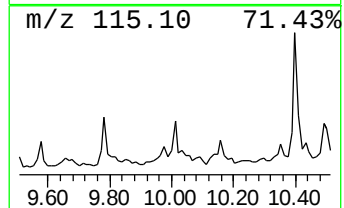
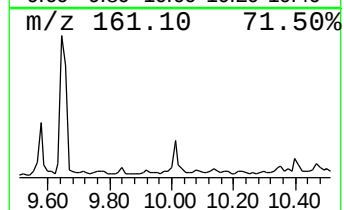
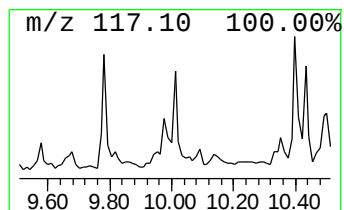
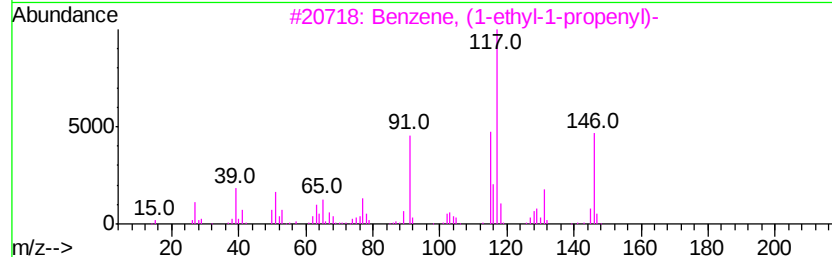
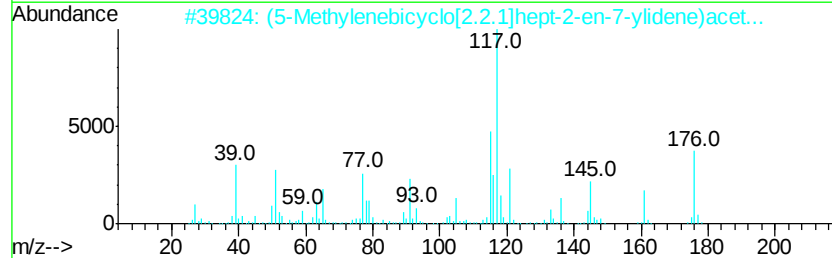
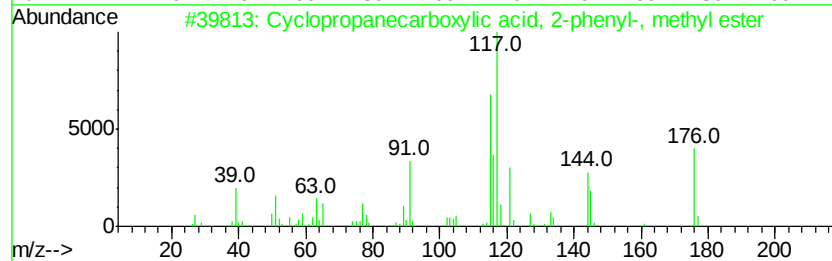
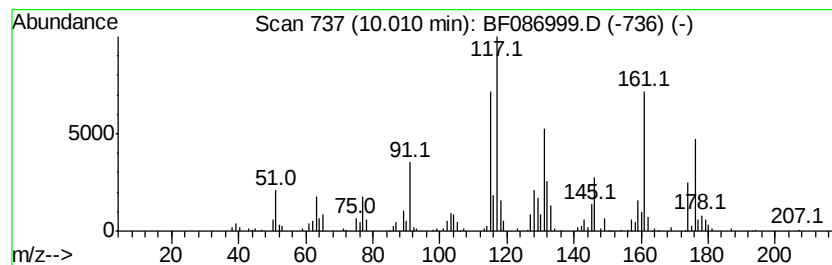
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown10.01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.01	2.67 ng	169856	Acenaphthene-d10	9.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropanecarboxylic acid, 2-p...	176	C11H12O2	020030-70-0	46
2		(5-Methylenebicyclo[2.2.1]hept-2...	176	C11H12O2	1000211-09-6	41
3		Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	38
4		2-Ethyl-2,3-dihydro-1H-indene	146	C11H14	056147-63-8	38
5		Cyclopropanecarboxylic acid, 2-p...	176	C11H12O2	010488-03-6	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051316\
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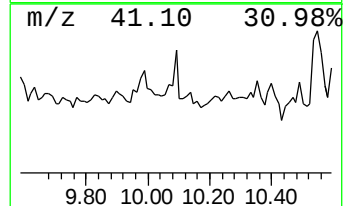
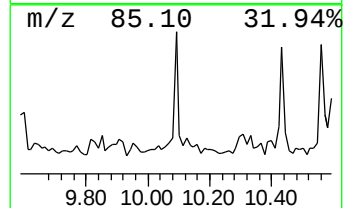
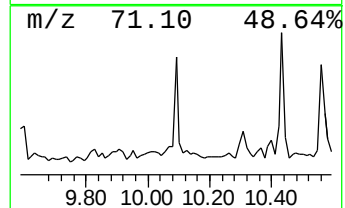
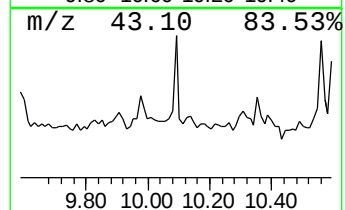
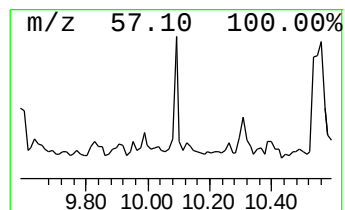
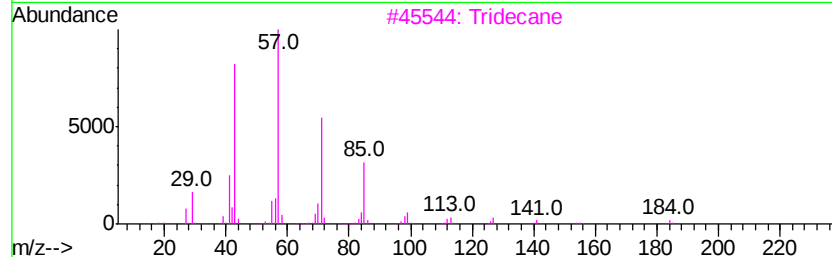
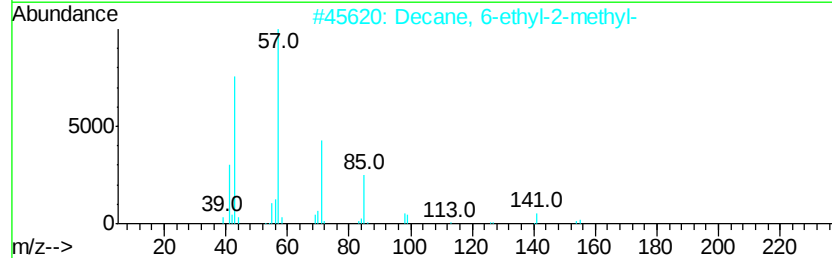
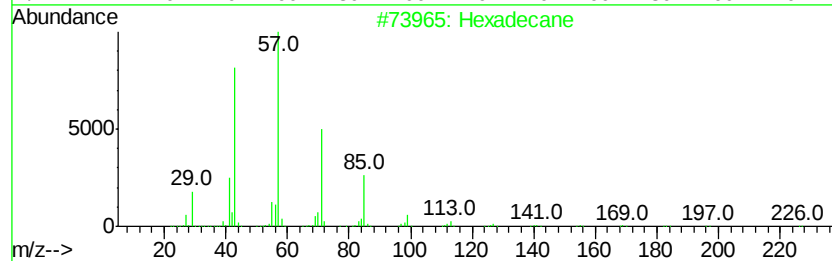
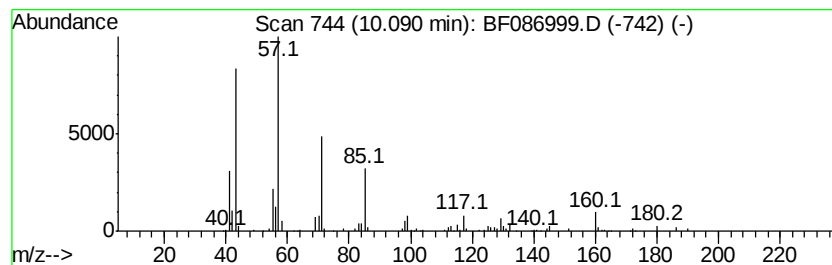
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Hexadecane Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.09	2.02 ng	128549	Acenaphthene-d10	9.64
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Hexadecane	226 C16H34	000544-76-3	64
2	Decane, 6-ethyl-2-methyl-	184 C13H28	062108-21-8	59
3	Tridecane	184 C13H28	000629-50-5	53
4	Tetradecane, 1-iodo-	324 C14H29I	019218-94-1	53
5	Dodecane, 2-methyl-6-propyl-	226 C16H34	055045-08-4	50



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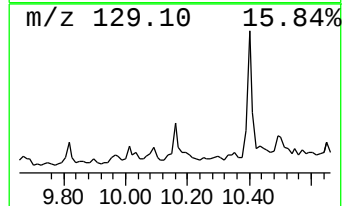
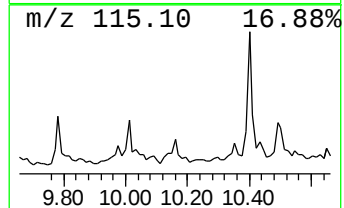
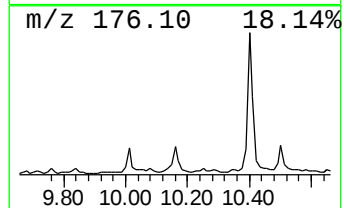
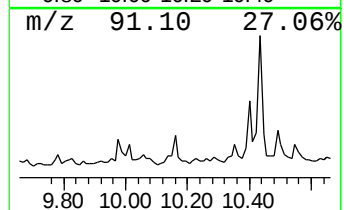
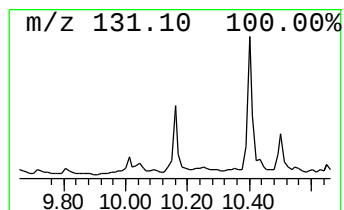
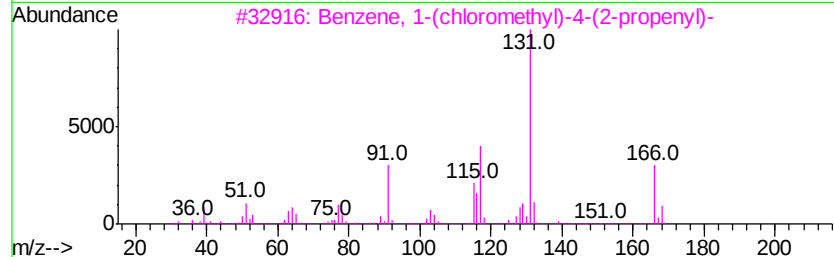
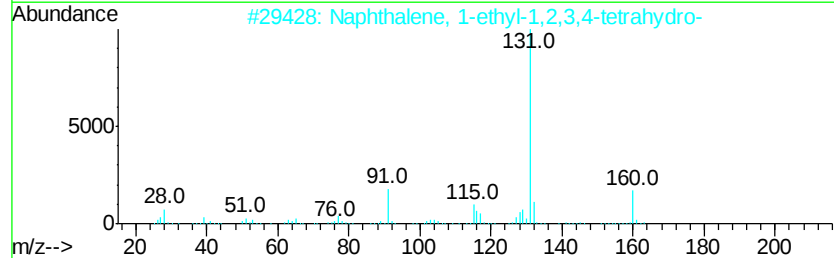
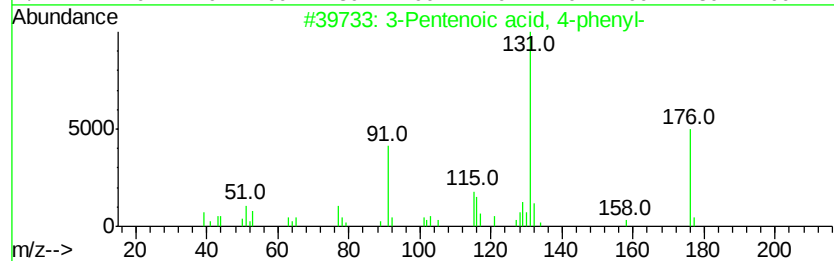
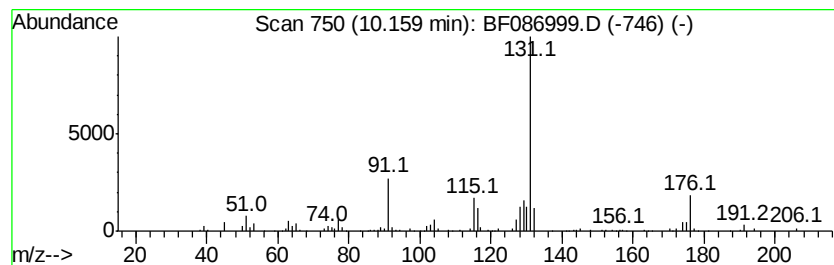
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 3-Pentenoic acid, 4-phenyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.16	2.58 ng	164213	Acenaphthene-d10	9.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Pentenoic acid, 4-phenyl-	176	C11H12O2	053774-19-9	72
2		Naphthalene, 1-ethyl-1,2,3,4-tet...	160	C12H16	013556-58-6	64
3		Benzene, 1-(chloromethyl)-4-(2-p...	166	C10H11Cl	036875-10-2	64
4		2-Butene, 3-chloro-1-phenyl-, (Z)-	166	C10H11Cl	016608-68-7	64
5		1H-Indene, 2,3-dihydro-4-propyl-	160	C12H16	092013-16-6	64



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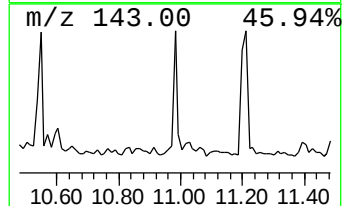
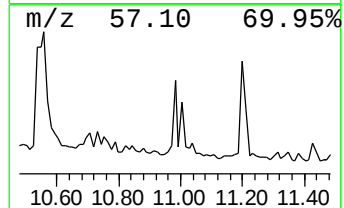
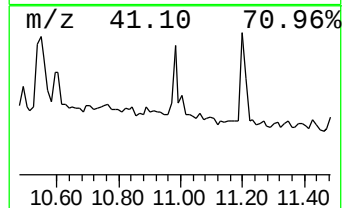
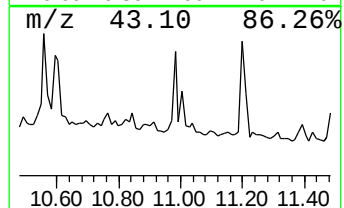
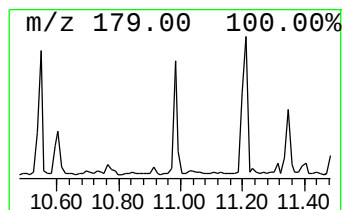
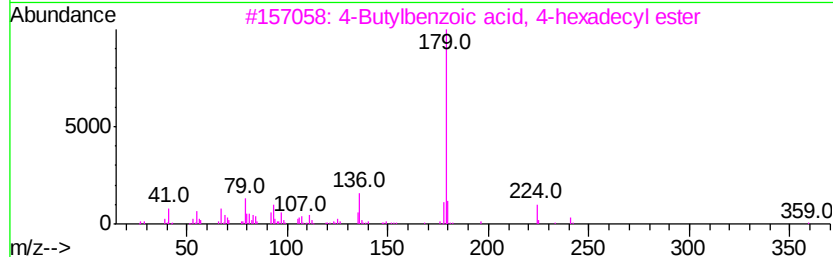
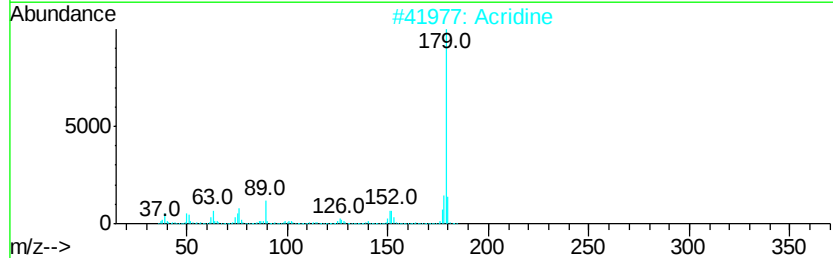
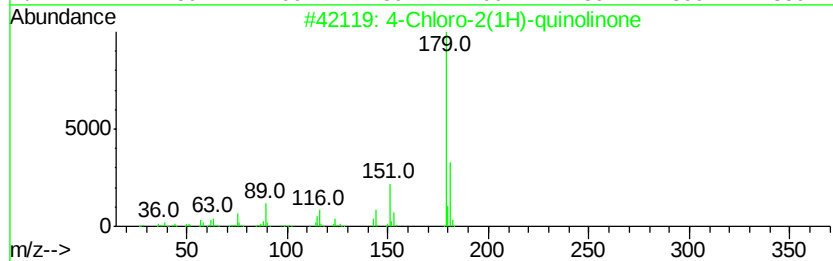
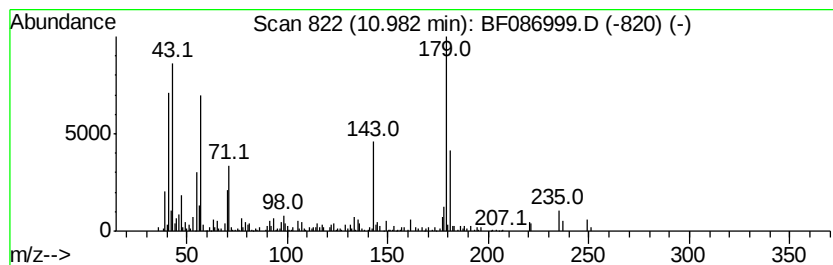
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown10.98 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.98	3.53 ng	206098	Phenanthrene-d10	11.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Chloro-2(1H)-quinolinone	179	C9H6ClNO	020146-59-2	43
2			Acridine	179	C13H9N	000260-94-6	27
3			4-Butylbenzoic acid, 4-hexadecyl...	402	C27H46O2	1000292-21-2	16
4			6H-Purin-6-one, 2-(dimethylamino...	179	C7H9N5O	001445-15-4	12
5			Benzoic acid, 4-nitroso-, ethyl ...	179	C9H9NO3	007476-79-1	12



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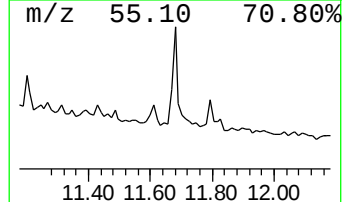
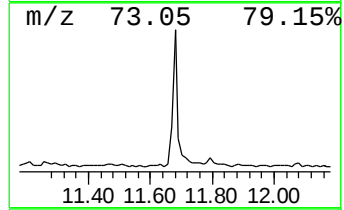
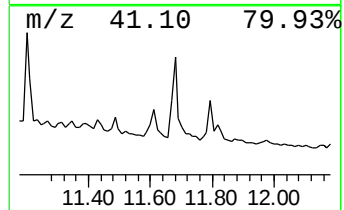
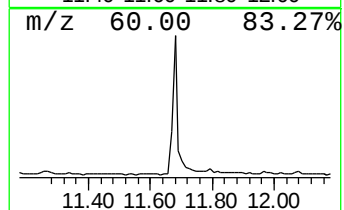
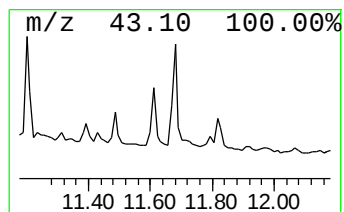
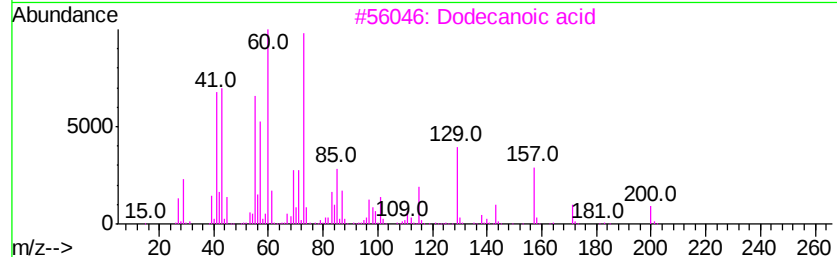
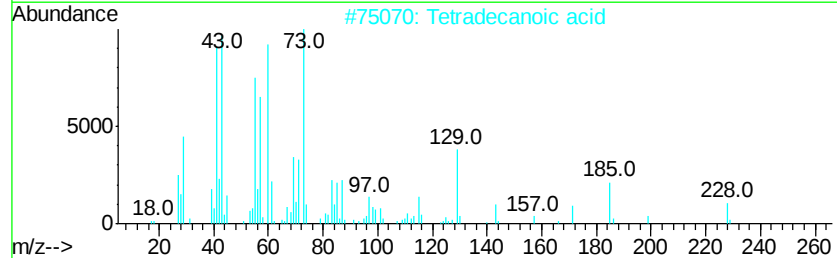
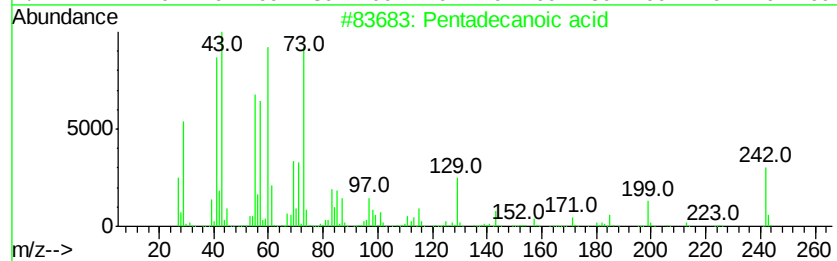
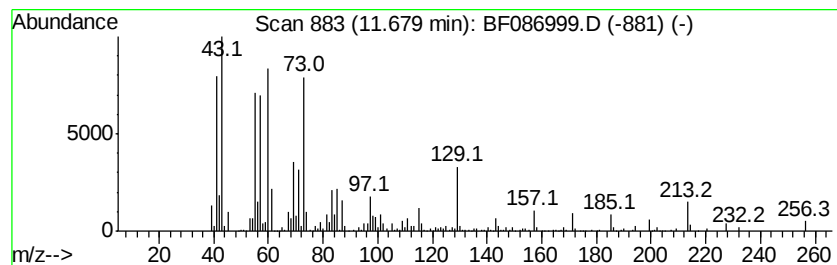
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Pentadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.68	4.42 ng	258084	Phenanthrene-d10	11.12

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid	242	C15H30O2	001002-84-2	80
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	74
3	Dodecanoic acid	200	C12H24O2	000143-07-7	72
4	Undecanoic acid	186	C11H22O2	000112-37-8	64
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051316\
Data File : BF086999.D
Acq On : 14 May 2016 5:07
Operator : UM/SJ
Sample : H2966-05
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-36

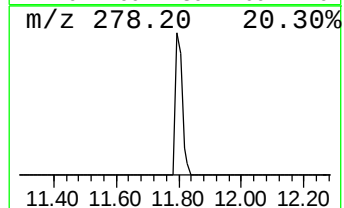
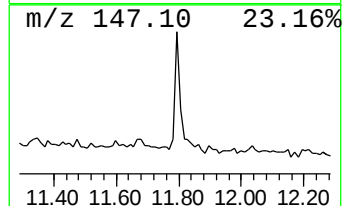
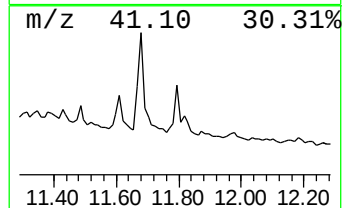
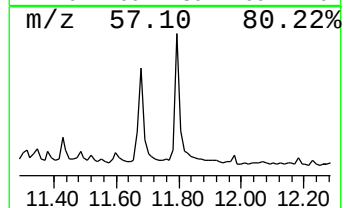
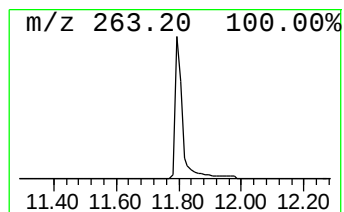
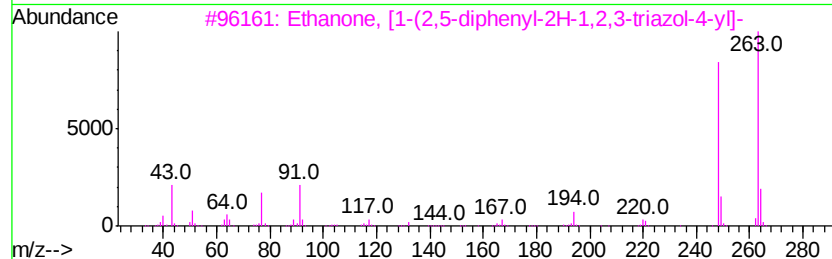
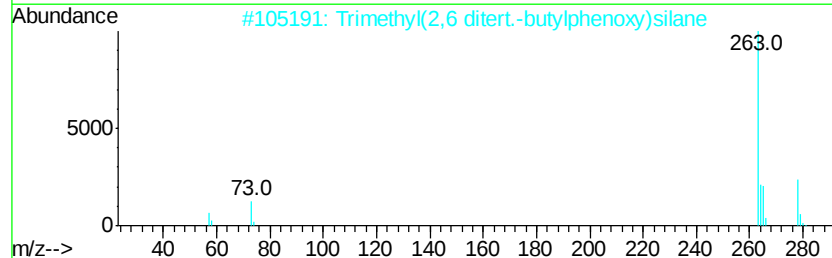
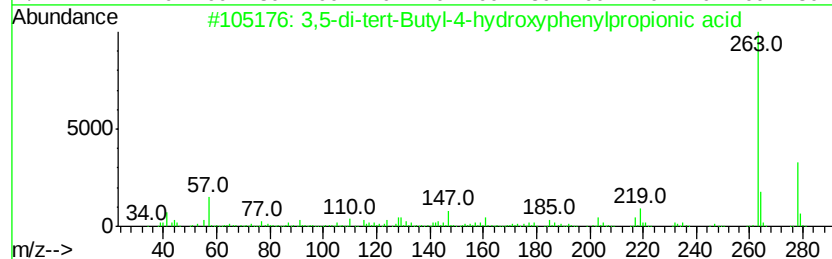
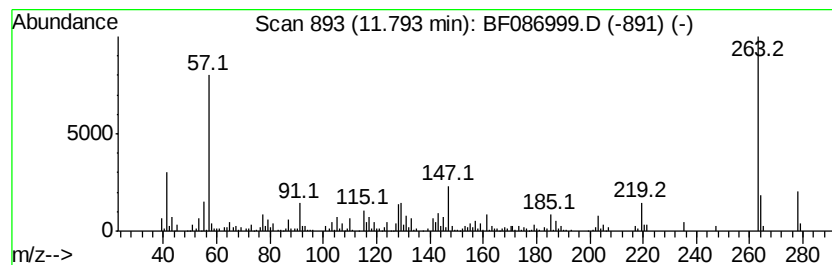
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	4.05 ng	236440	Phenanthrene-d10	11.12

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-di-tert-Butyl-4-hydroxyphenyl...	278	C17H26O3	020170-32-5	95
2			Trimethyl(2,6 ditert.-butylpheno...	278	C17H30OSi	010416-73-6	47
3			Ethanone, [1-(2,5-diphenyl-2H-1,...	263	C16H13N3O	322647-34-7	46
4			Isoquinoline, 5,6,7,8-tetrametho...	263	C14H17N04	093474-27-2	46
5			Pentan-2-one, 4-(2-fluorenylimino)-	263	C18H17NO	050651-58-6	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF051316\
Data File : BF086999.D
Acq On : 14 May 2016 5:07
Operator : UM/SJ
Sample : H2966-05
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-36

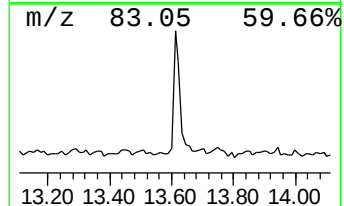
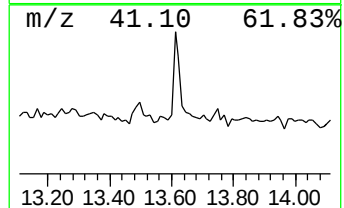
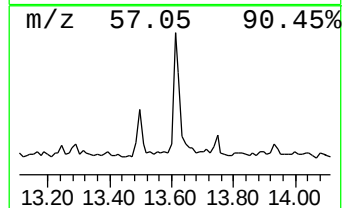
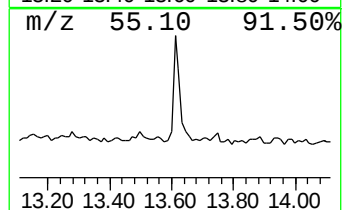
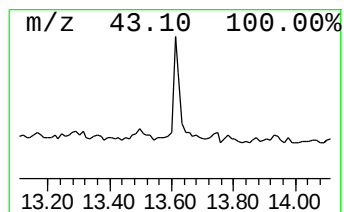
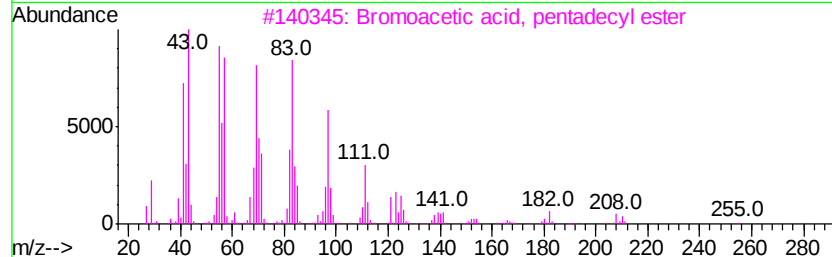
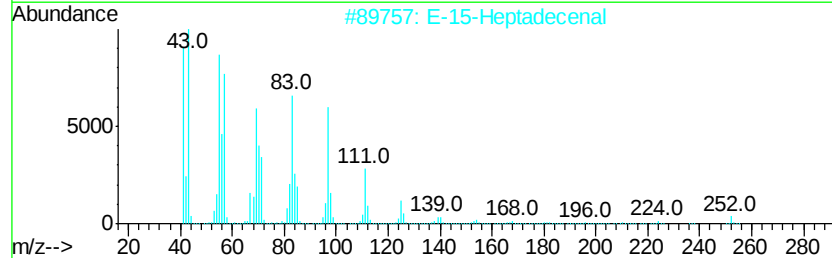
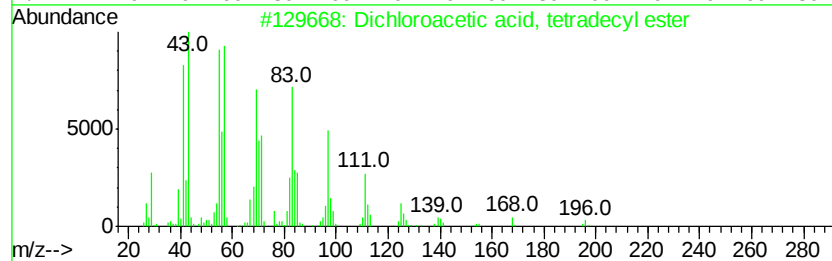
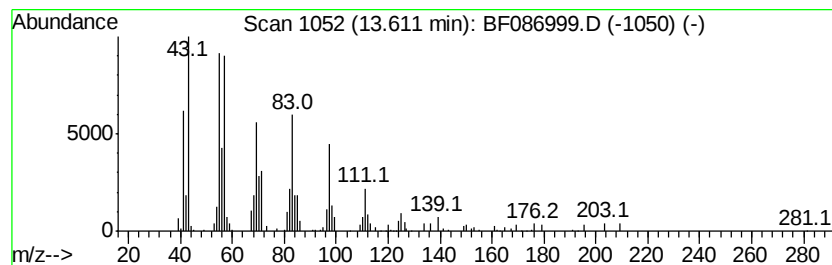
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Dichloroacetic acid, tetrad... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.61	2.50 ng	120783	Chrysene-d12	13.75

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dichloroacetic acid, tetradecyl ...	324	C16H30Cl2O2	083005-02-1	91
2			E-15-Heptadecenal	252	C17H32O	1000130-97-9	91
3			Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	91
4			Cyclohexadecane	224	C16H32	000295-65-8	91
5			Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	91



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051316\
Data File : BF086999.D
Acq On : 14 May 2016 5:07
Operator : UM/SJ
Sample : H2966-05
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-36

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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
unknown6.39	6.39	74.3	ng	3414520	1	6.62	919541	20.0
Benzene, 1-methyl...	7.37	2.0	ng	130023	2	7.90	1287510	20.0
Bicyclo[2.2.1]hep...	7.64	2.4	ng	156396	2	7.90	1287510	20.0
unknown8.66	8.66	4.8	ng	307617	2	7.90	1287510	20.0
Diphenylmethane	9.35	2.7	ng	170430	3	9.64	1272760	20.0
Benzeneacetaldehy...	9.58	4.7	ng	301262	3	9.64	1272760	20.0
unknown9.98	9.98	4.1	ng	259553	3	9.64	1272760	20.0
unknown10.01	10.01	2.7	ng	169856	3	9.64	1272760	20.0
Hexadecane	10.09	2.0	ng	128549	3	9.64	1272760	20.0
3-Pentenoic acid,...	10.16	2.6	ng	164213	3	9.64	1272760	20.0
unknown10.98	10.98	3.5	ng	206098	4	11.12	1168570	20.0
Pentadecanoic acid	11.68	4.4	ng	258084	4	11.12	1168570	20.0
3,5-di-tert-Butyl...	11.79	4.0	ng	236440	4	11.12	1168570	20.0
Dichloroacetic ac...	13.61	2.5	ng	120783	5	13.75	966212	20.0