

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
 Data File : BF080148.D
 Acq On : 25 Jun 2015 1:10
 Operator : TP/IZ
 Sample : G2717-03
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AC-UNDERCUT(0-2)-6

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-BF061715.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	5.533	211	214	216	rBV	192601	258043	18.76%	2.082%
2	5.921	246	248	251	rBV	1036881	948548	68.95%	7.655%
3	6.321	281	283	285	rBB	23127	19329	1.40%	0.156%
4	6.539	300	302	311	rBB	46819	52899	3.85%	0.427%
5	6.916	332	335	337	rBV	1255776	1051905	76.46%	8.489%
6	7.076	347	349	352	rBV	1106043	996215	72.41%	8.039%
7	7.316	367	370	372	rBB	266964	231493	16.83%	1.868%
8	7.476	382	384	386	rBV	786537	767339	55.77%	6.192%
9	7.864	416	418	421	rBV	593407	571571	41.55%	4.613%
10	8.607	481	483	485	rBV	316516	305951	22.24%	2.469%
11	9.156	528	531	532	rVB	13799	15942	1.16%	0.129%
12	9.556	564	566	569	rBV	20303	26249	1.91%	0.212%
13	9.682	573	577	579	rBV	1590607	1318152	95.81%	10.638%
14	10.070	609	611	613	rBV	93953	90167	6.55%	0.728%
15	10.379	635	638	640	rBV	360036	354299	25.75%	2.859%
16	11.088	697	700	702	rBV	43930	51949	3.78%	0.419%
17	11.168	704	707	709	rBV	851218	757926	55.09%	6.116%
18	11.259	713	715	719	rBV2	9493	15722	1.14%	0.127%
19	11.876	767	769	773	rBV	411466	449053	32.64%	3.624%
20	11.956	773	776	777	rBV	17040	15562	1.13%	0.126%
21	12.391	811	814	816	rBV3	33097	39963	2.90%	0.323%
22	12.436	816	818	820	rVV	152986	140755	10.23%	1.136%
23	12.516	823	825	829	rVB2	13727	25562	1.86%	0.206%
24	13.156	879	881	883	rVB	92234	95816	6.96%	0.773%
25	13.419	899	904	907	rBV	111180	151137	10.99%	1.220%
26	13.568	914	917	919	rBV	1093463	1375781	100.00%	11.103%
27	14.059	958	960	962	rBV2	7755	16366	1.19%	0.132%
28	14.105	962	964	967	rVB	34230	34856	2.53%	0.281%
29	14.619	1007	1009	1017	rBV2	173693	290254	21.10%	2.342%
30	14.791	1023	1024	1025	rBV	13944	15650	1.14%	0.126%
31	14.825	1025	1027	1029	rVV2	365683	500677	36.39%	4.040%
32	14.859	1029	1030	1032	rVB	51618	50482	3.67%	0.407%
33	15.442	1078	1081	1084	rBV2	50275	97055	7.05%	0.783%
34	16.162	1141	1144	1145	rBV	48439	78201	5.68%	0.631%

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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

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Peak separation: 5

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35	16.265	1151	1153	1155	rBV	36644	49451	3.59%	0.399%
36	16.540	1175	1177	1180	rVB	34818	52924	3.85%	0.427%
37	16.620	1181	1184	1186	rBV	38766	58560	4.26%	0.473%
38	16.700	1188	1191	1193	rBV	285379	395285	28.73%	3.190%
39	17.317	1242	1245	1248	rBV	23279	48904	3.55%	0.395%
40	17.374	1248	1250	1256	rVB5	28621	84918	6.17%	0.685%
41	17.603	1268	1270	1280	rVB7	28427	99787	7.25%	0.805%
42	18.003	1302	1305	1308	rBV	42735	82042	5.96%	0.662%
43	18.163	1316	1319	1324	rVB2	25888	63751	4.63%	0.514%
44	18.540	1348	1352	1360	rVB2	44079	160552	11.67%	1.296%
45	19.088	1396	1400	1403	rBV	33915	84474	6.14%	0.682%

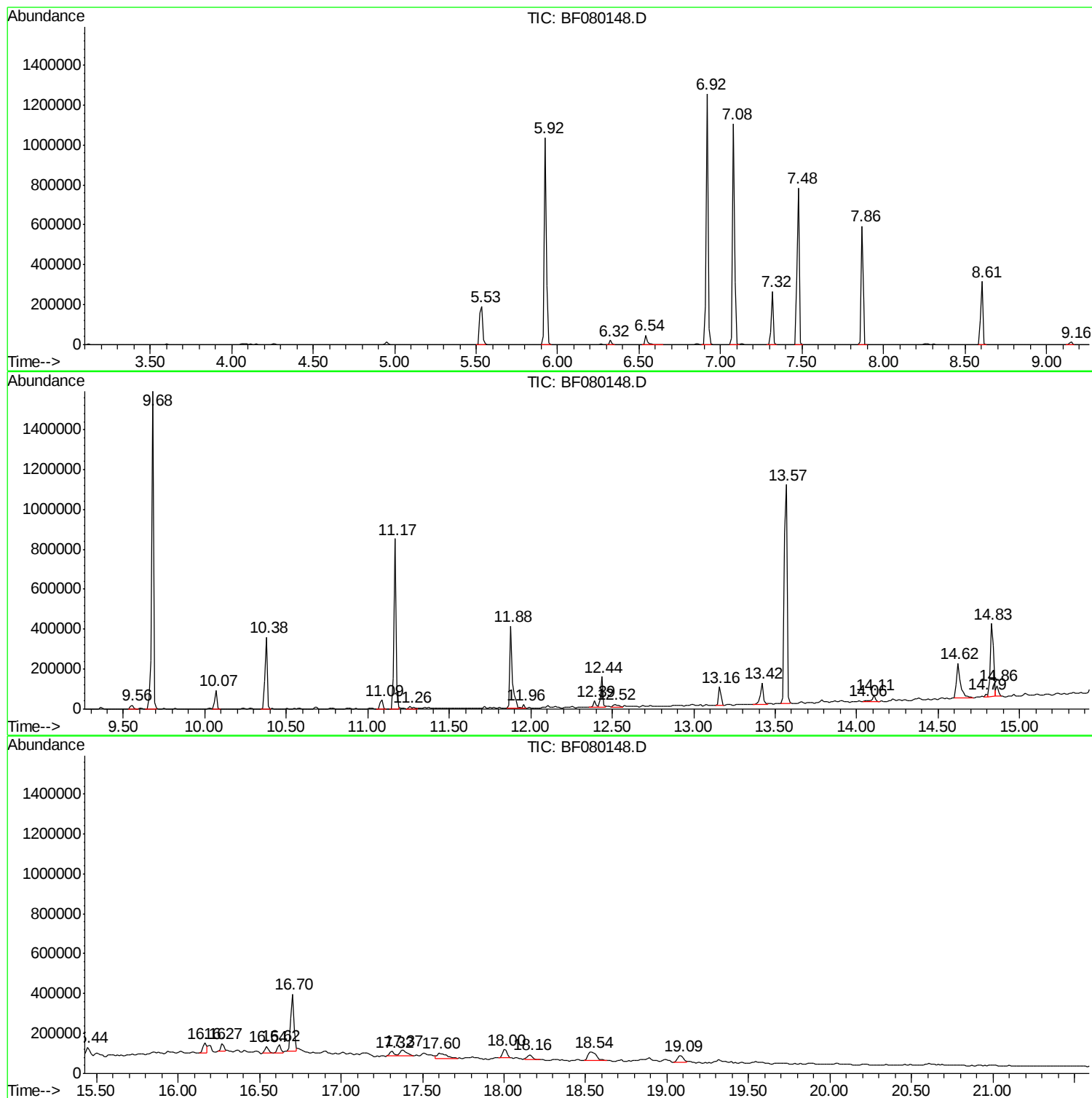
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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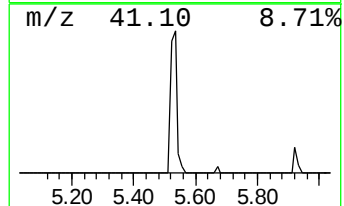
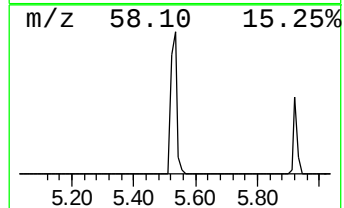
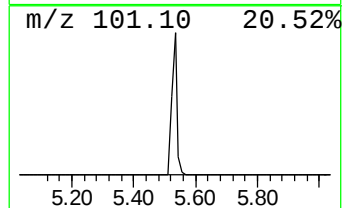
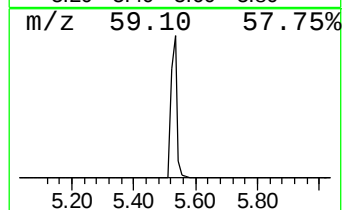
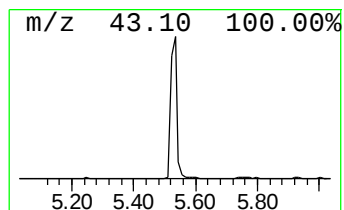
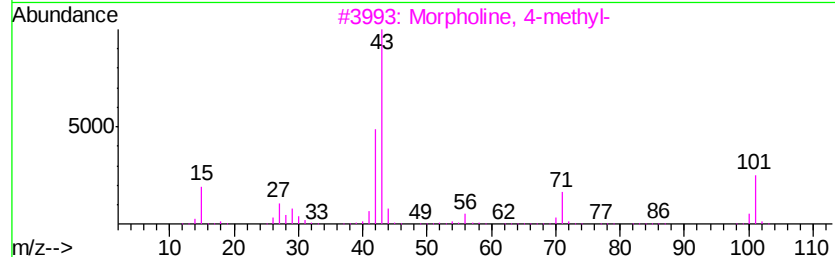
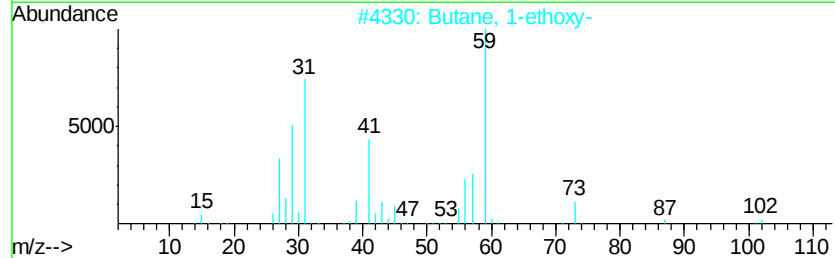
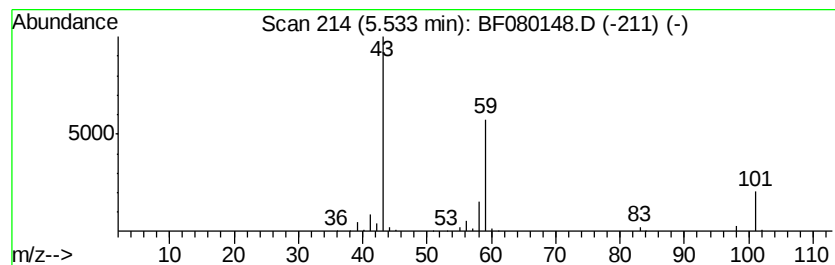
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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.53	22.29 ng	258043	1,4-Dichlorobenzene-d4	7.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
3		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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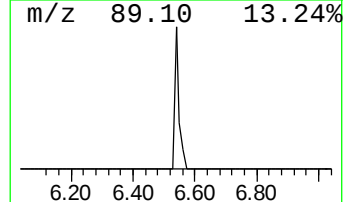
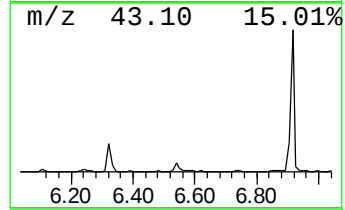
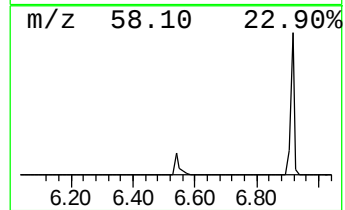
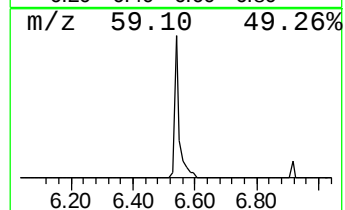
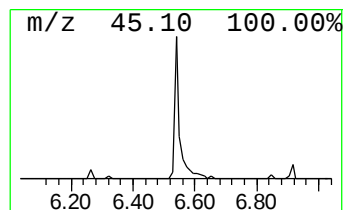
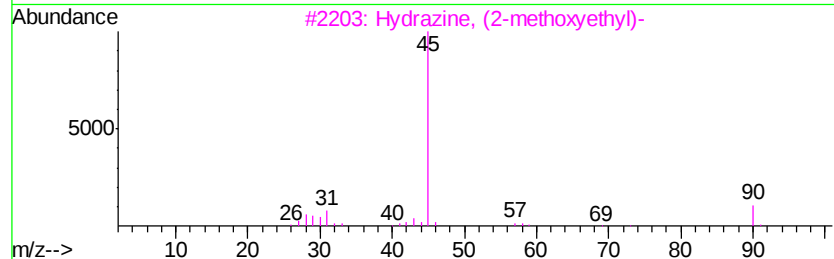
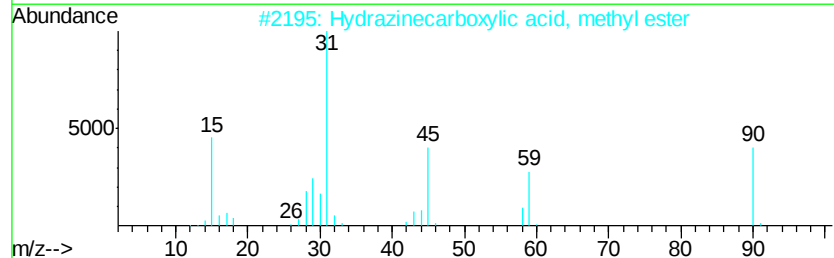
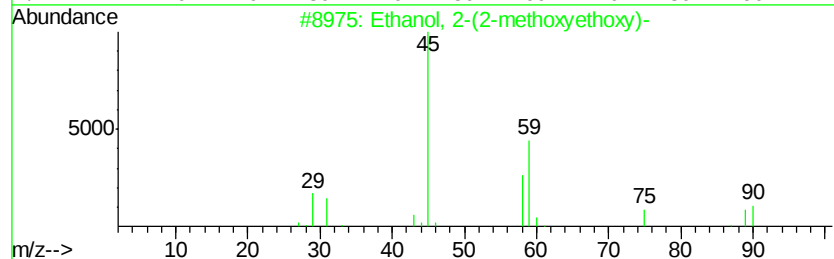
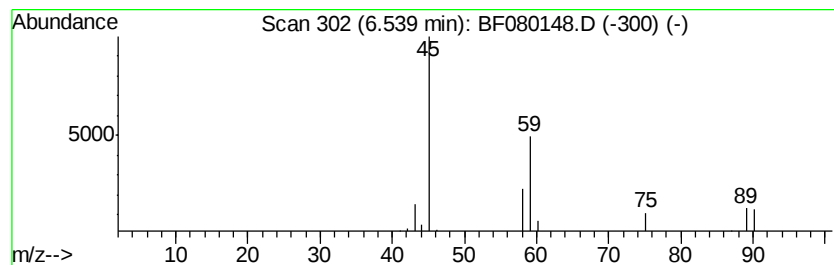
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 Ethanol, 2-(2-methoxyethoxy)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.54	4.57 ng	52899	1,4-Dichlorobenzene-d4	7.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ethanol, 2-(2-methoxyethoxy)-	120	C5H12O3	000111-77-3	83
2		Hydrazinecarboxylic acid, methyl...	90	C2H6N2O2	006294-89-9	39
3		Hydrazine, (2-methoxyethyl)-	90	C3H10N2O	003044-15-3	37
4		Triethylene glycol	150	C6H14O4	000112-27-6	36
5		Ethane, 1,2-dimethoxy-	90	C4H10O2	000110-71-4	28



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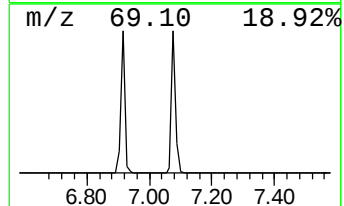
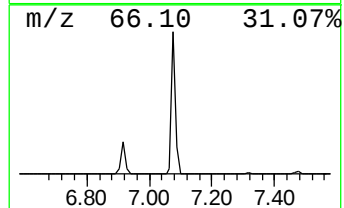
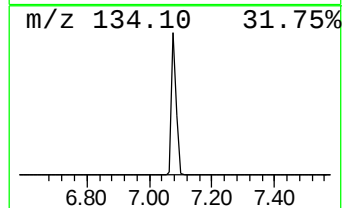
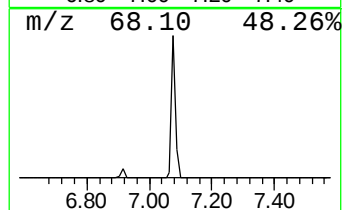
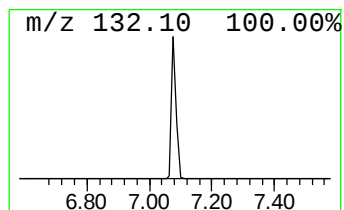
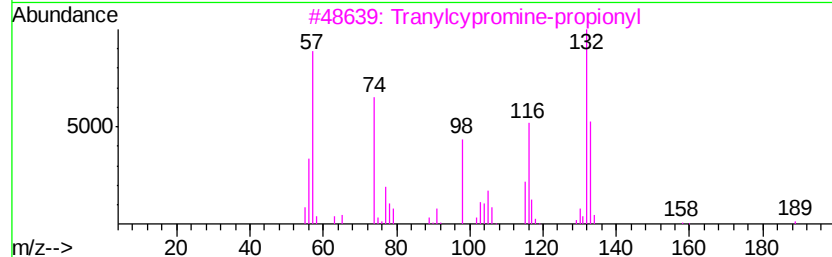
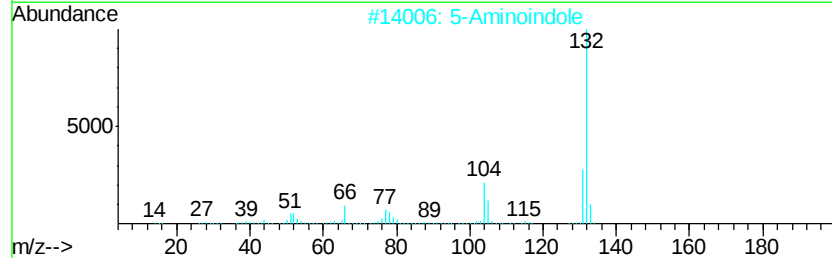
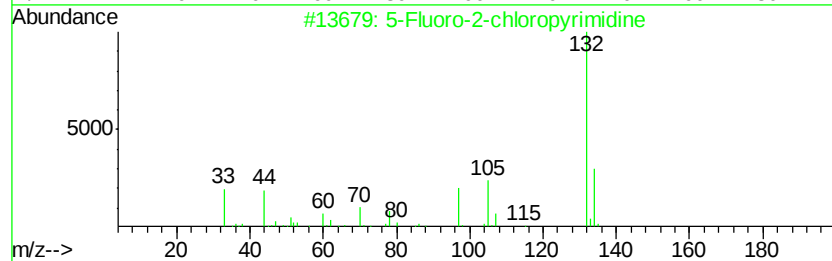
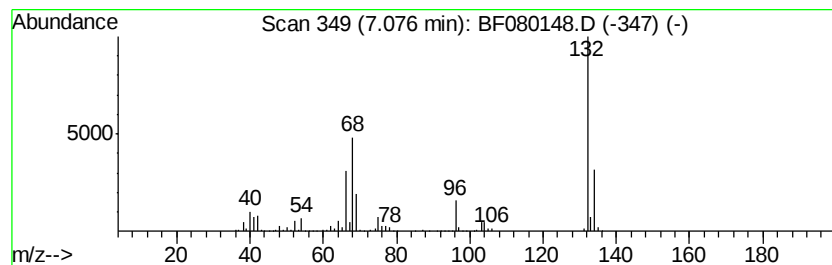
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown7.08 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.08	86.07 ng	996215	1,4-Dichlorobenzene-d4	7.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2			5-Aminoindole	132	C8H8N2	005192-03-0	10
3			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4			Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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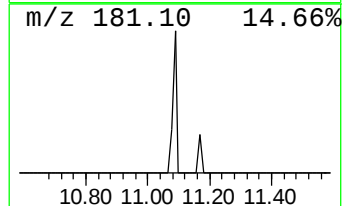
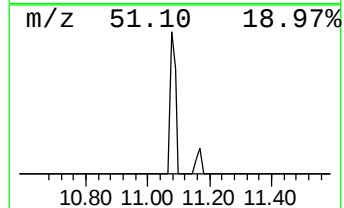
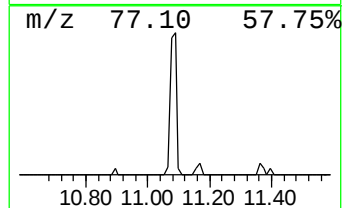
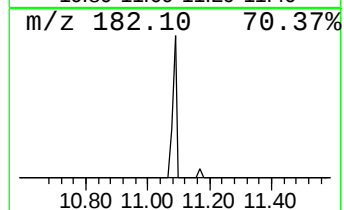
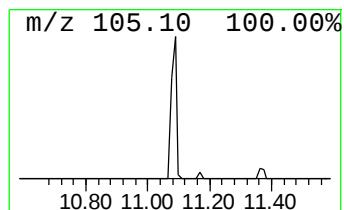
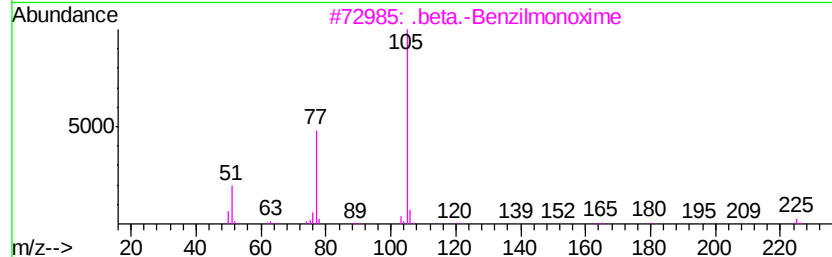
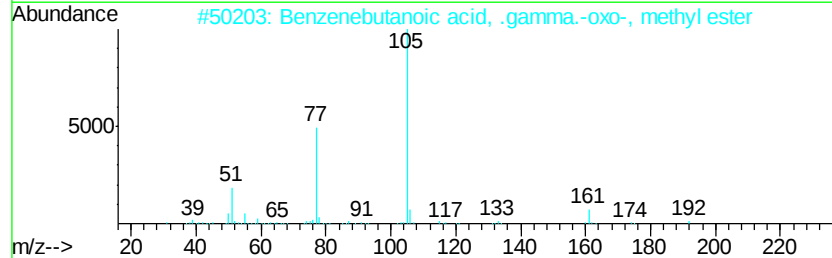
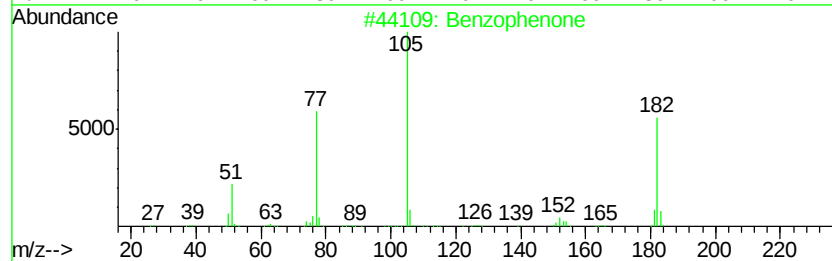
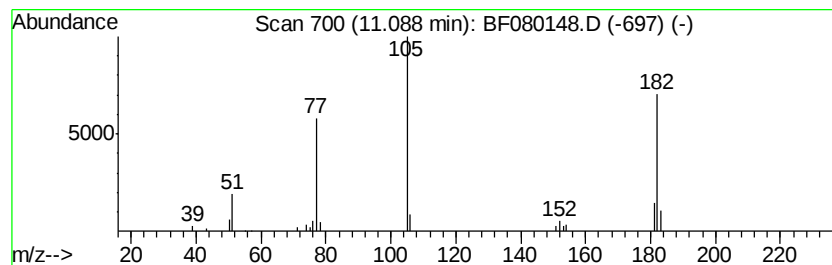
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 5 Benzophenone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.09	2.93 ng	51949	Acenaphthene-d10	10.38

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	96
2		Benzenebutanoic acid, .gamma.-ox...	192	C11H12O3	025333-24-8	38
3		.beta.-Benzilmonoxime	225	C14H11NO2	014090-77-8	38
4		Vinyl benzoate	148	C9H8O2	000769-78-8	38
5		1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	38



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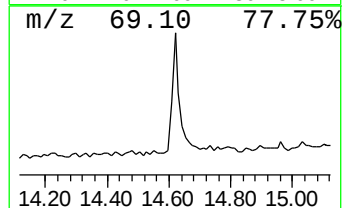
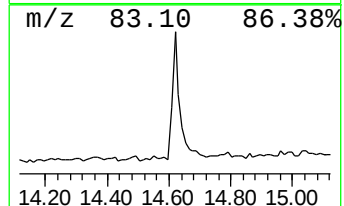
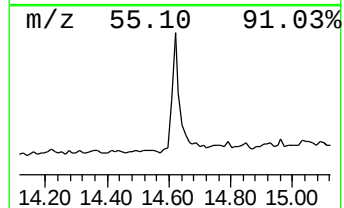
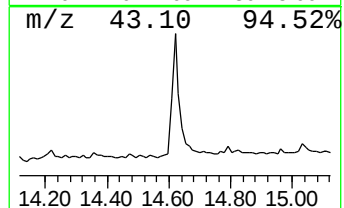
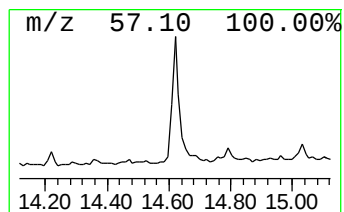
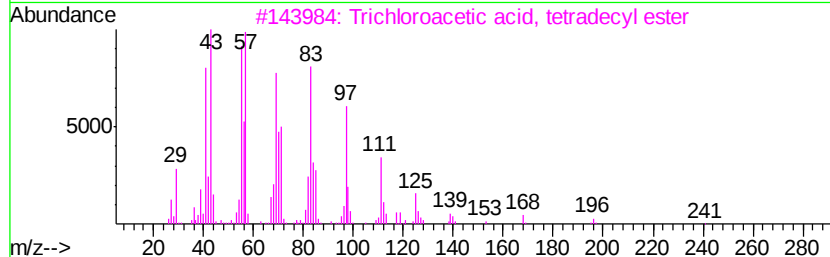
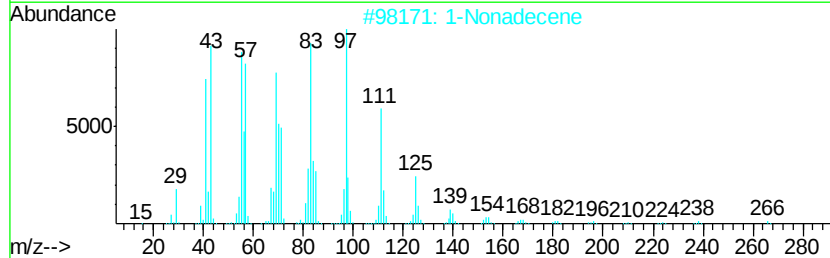
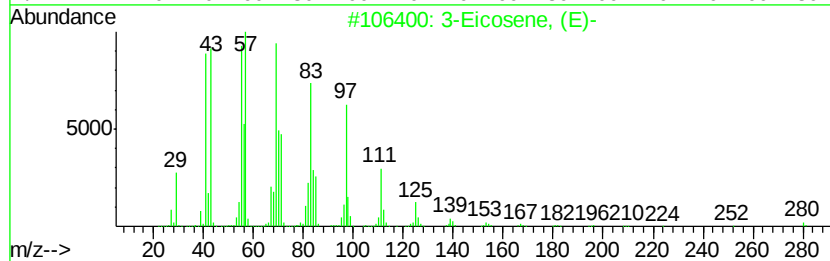
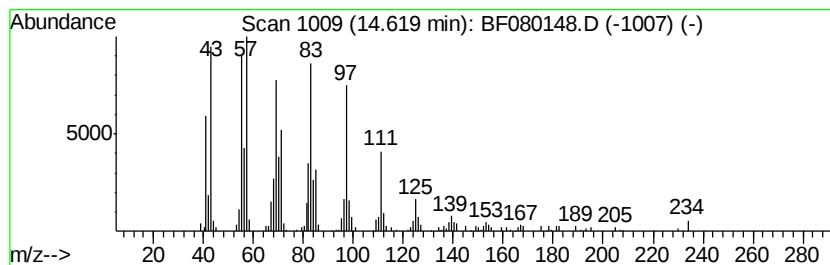
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Peak Number 6 3-Eicosene, (E)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.62	11.59 ng	290254	Chrysene-d12	14.83	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Eicosene, (E)-	280	C20H40	074685-33-9	91
2	1-Nonadecene	266	C19H38	018435-45-5	91
3	Trichloroacetic acid, tetradecyl...	358	C16H29Cl3O2	074339-52-9	91
4	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	87
5	Cyclopentadecane	210	C15H30	000295-48-7	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
Data File : BF080148.D
Acq On : 25 Jun 2015 1:10
Operator : TP/IZ
Sample : G2717-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
AC-UNDERCUT(0-2)-6

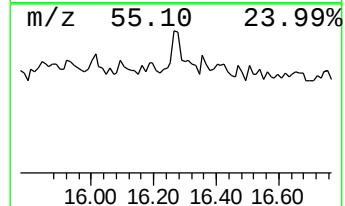
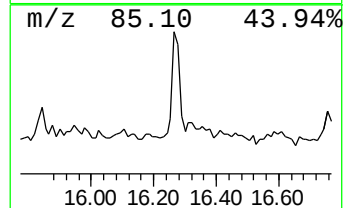
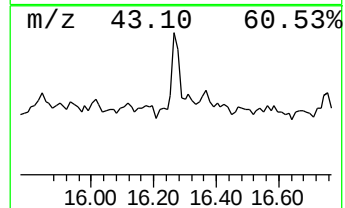
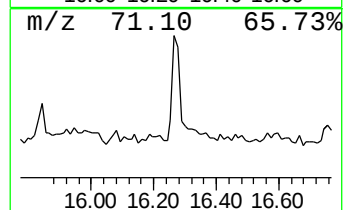
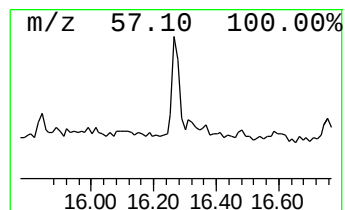
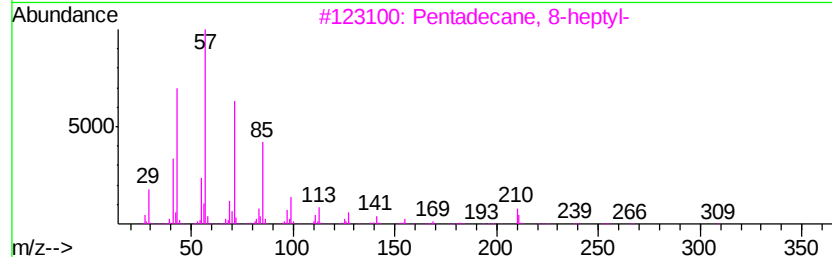
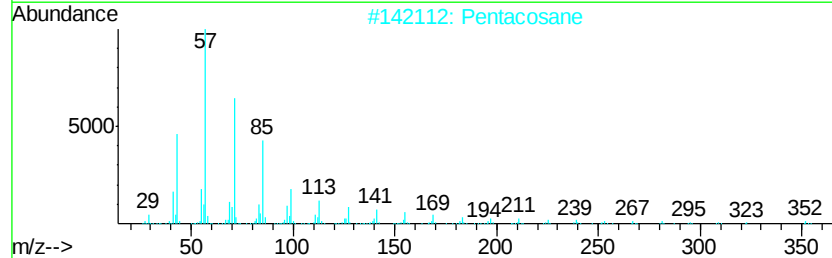
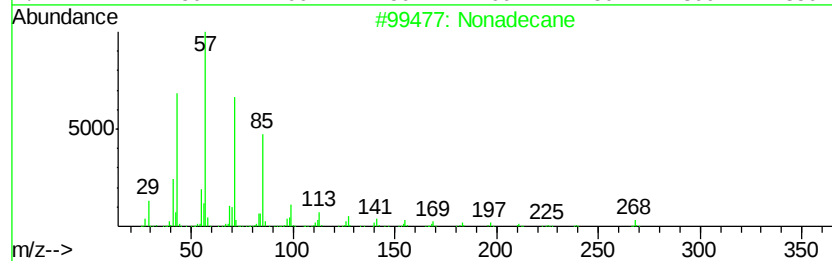
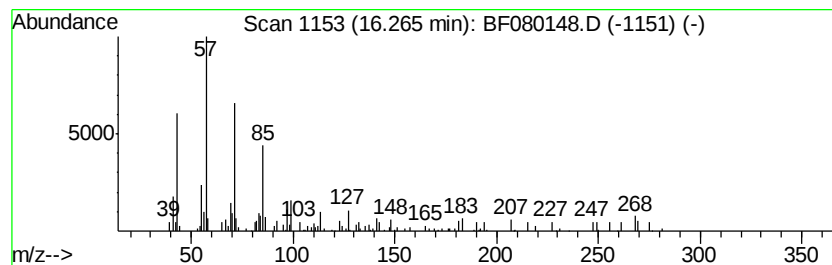
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Nonadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.27	2.50 ng	49451	Perylene-d12	16.70

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	76
2	Pentacosane		352	C25H52	000629-99-2	64
3	Pentadecane, 8-heptyl-		310	C22H46	071005-15-7	64
4	Eicosane, 7-hexyl-		366	C26H54	055333-99-8	64
5	Tetracosane		338	C24H50	000646-31-1	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
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Acq On : 25 Jun 2015 1:10
Operator : TP/IZ
Sample : G2717-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
AC-UNDERCUT(0-2)-6

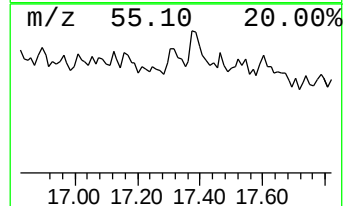
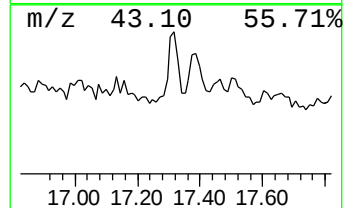
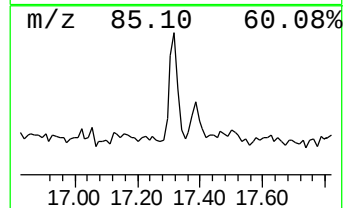
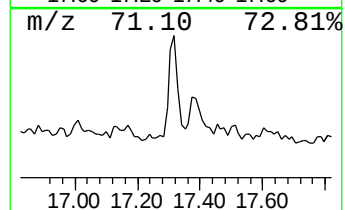
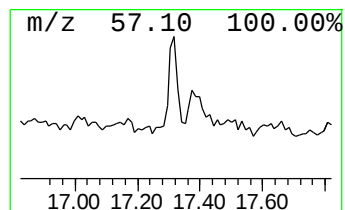
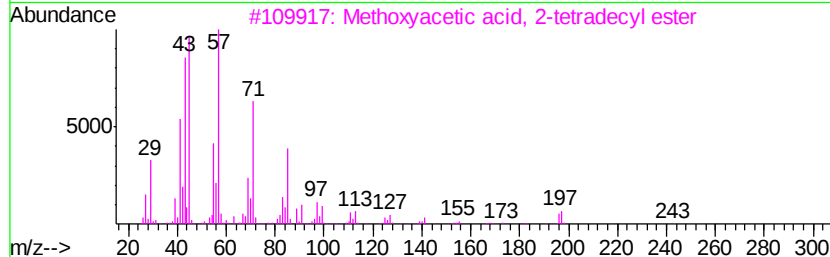
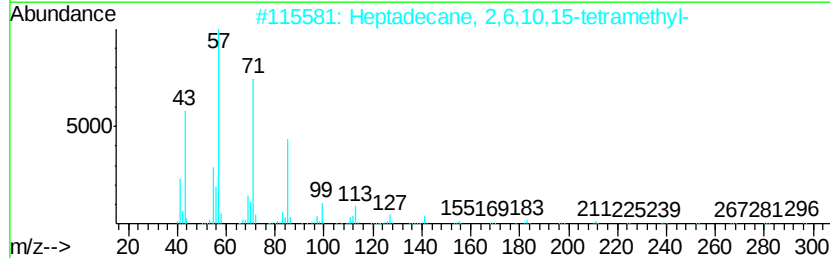
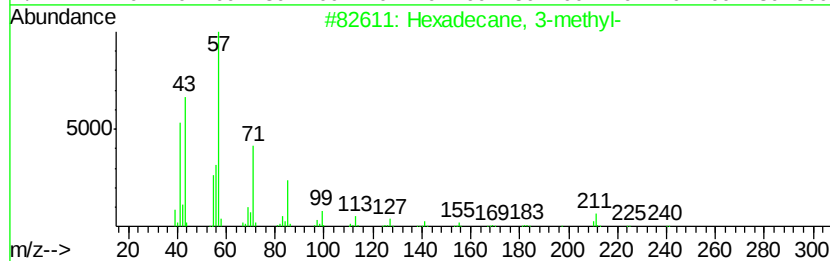
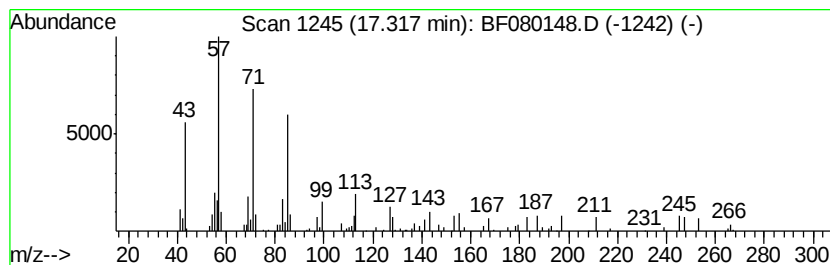
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexadecane, 3-methyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.32	2.47 ng	48904	Perylene-d12	16.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecane, 3-methyl-	240	C17H36	006418-43-5	64
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	59
3		Methoxyacetic acid, 2-tetradecyl...	286	C17H34O3	1000282-04-8	53
4		Pentadecane, 2-methyl-	226	C16H34	001560-93-6	53
5		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
Data File : BF080148.D
Acq On : 25 Jun 2015 1:10
Operator : TP/IZ
Sample : G2717-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
AC-UNDERCUT(0-2)-6

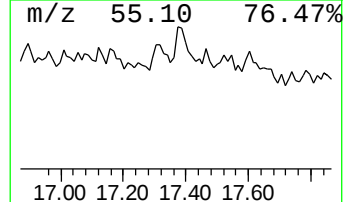
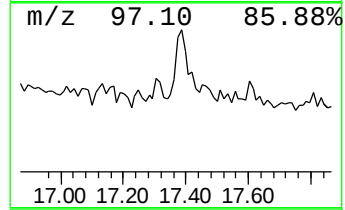
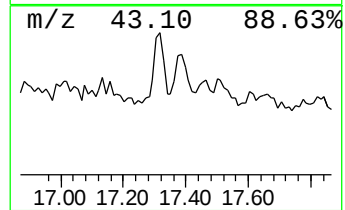
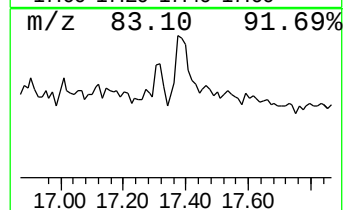
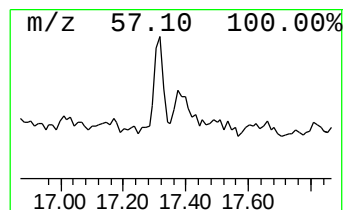
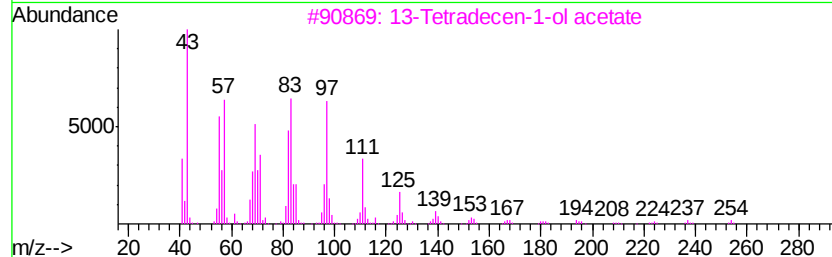
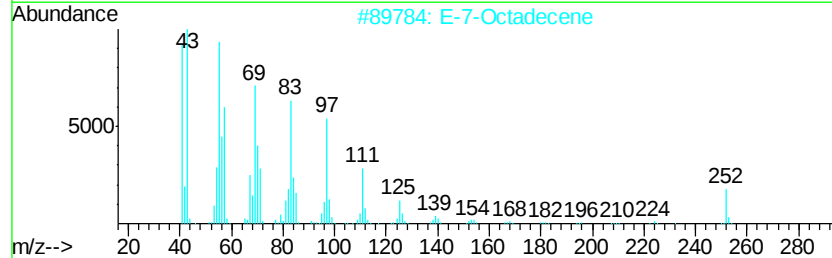
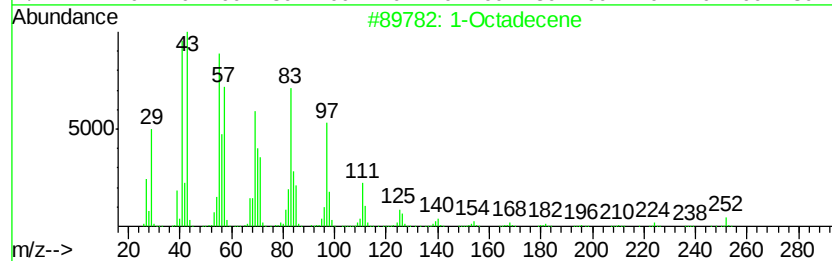
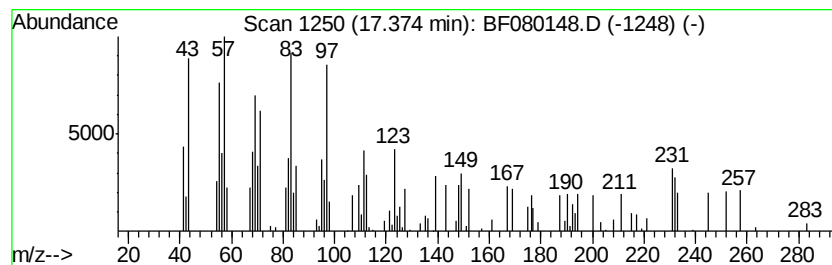
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 unknown17.37 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.37	4.30 ng	84918	Perylene-d12	16.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Octadecene	252	C18H36	000112-88-9	49
2			E-7-Octadecene	252	C18H36	1000130-92-0	47
3			13-Tetradecen-1-ol acetate	254	C16H30O2	056221-91-1	46
4			17-Pentatriacontene	491	C35H70	006971-40-0	46
5			11-Dodecen-2-one, 7,7-dimethyl-	210	C14H26O	035194-22-0	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
Data File : BF080148.D
Acq On : 25 Jun 2015 1:10
Operator : TP/IZ
Sample : G2717-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
AC-UNDERCUT(0-2)-6

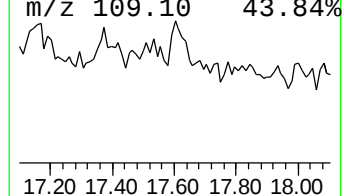
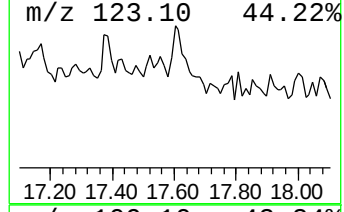
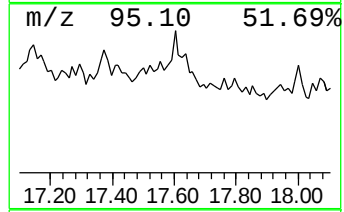
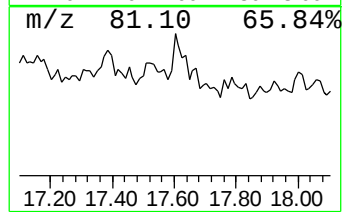
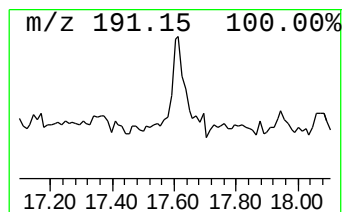
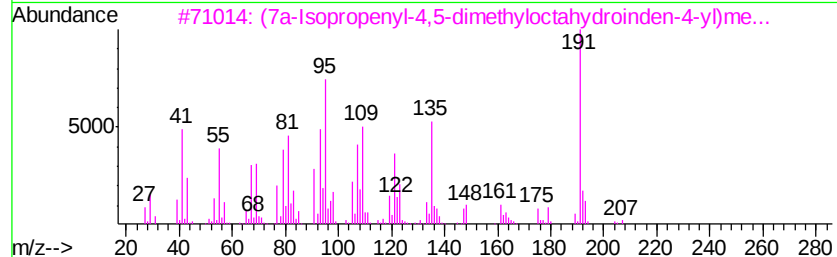
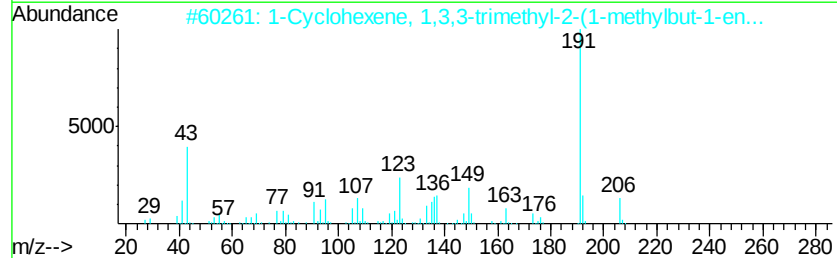
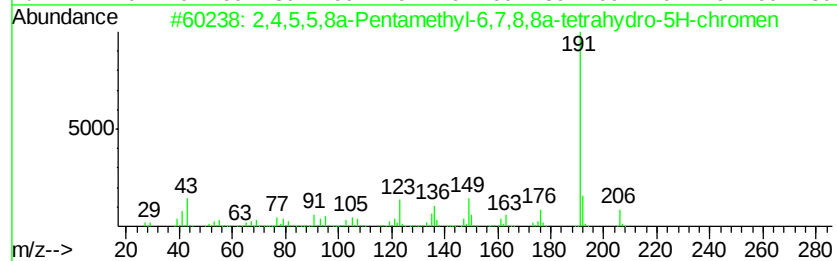
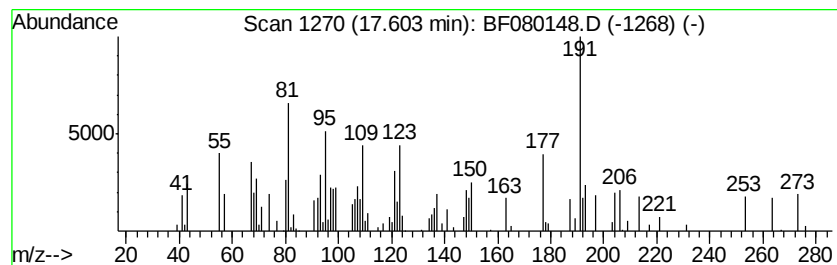
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2,4,5,5,8a-Pentamethyl-6,7,... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.60	5.05 ng	99787	Perylene-d12	16.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	55		
2	1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	43		
3	(7a-Isopropenyl-4,5-dimethylocta...	222	C15H26O	1000187-02-9	32		
4	1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	22		
5	1,2-Dithiolane-3-pentanoic acid	206	C8H14O2S2	000062-46-4	20		



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
Data File : BF080148.D
Acq On : 25 Jun 2015 1:10
Operator : TP/IZ
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Misc :
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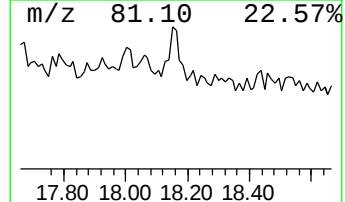
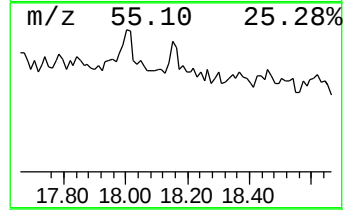
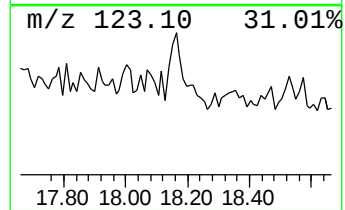
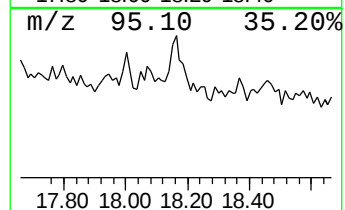
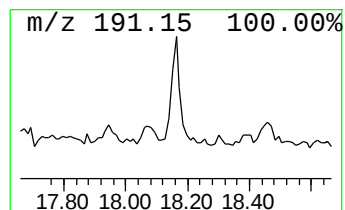
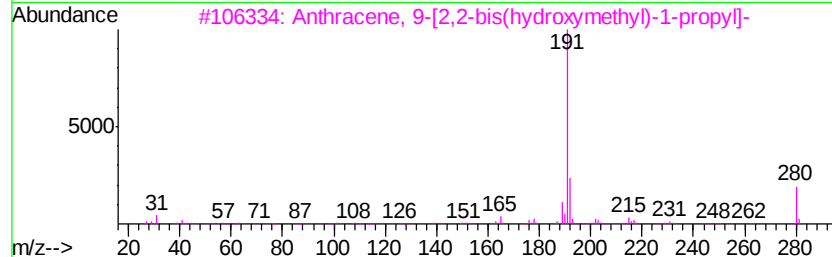
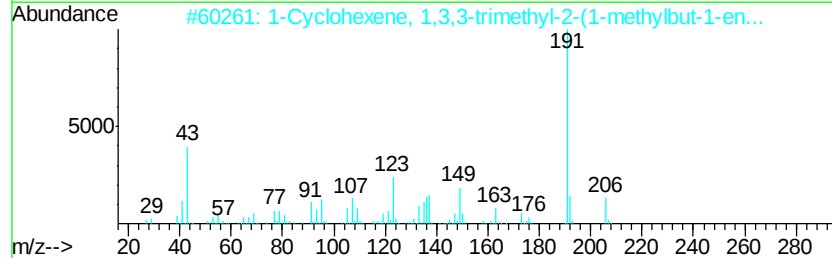
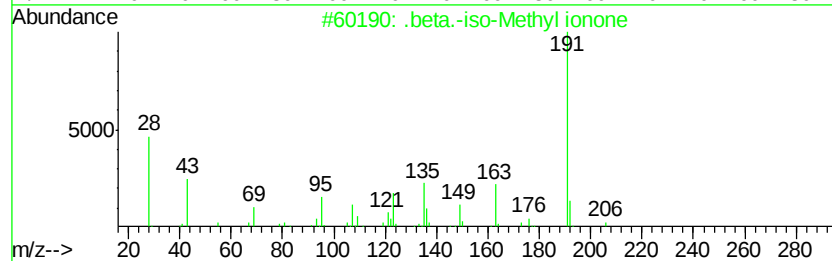
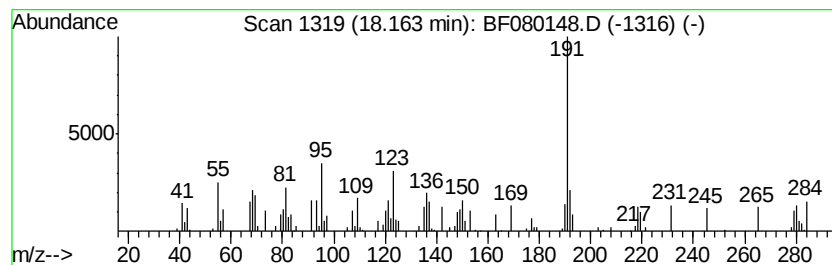
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 .beta.-iso-Methyl ionone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.16	3.23 ng	63751	Perylene-d12	16.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	50
2		1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	42
3		Anthracene, 9-[2,2-bis(hydroxyme...	280	C19H20O2	144000-85-1	38
4		1,1-Diphenyl-2-propynyl dimethyl...	279	C18H17NO2	010473-88-8	38
5		Cyclopropene, 3-bromo-1,2-diphenyl-	270	C15H11Br	029559-17-9	37



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.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...	5.53	22.3	ng	258043	1	7.32	231493	20.0
Ethanol, 2-(2-met...	6.54	4.6	ng	52899	1	7.32	231493	20.0
unknown7.08	7.08	86.1	ng	996215	1	7.32	231493	20.0
Benzophenone	11.09	2.9	ng	51949	3	10.38	354299	20.0
3-Eicosene, (E)-	14.62	11.6	ng	290254	5	14.83	500677	20.0
Nonadecane	16.27	2.5	ng	49451	6	16.70	395285	20.0
Hexadecane, 3-met...	17.32	2.5	ng	48904	6	16.70	395285	20.0
unknown17.37	17.37	4.3	ng	84918	6	16.70	395285	20.0
2,4,5,5,8a-Pentam...	17.60	5.0	ng	99787	6	16.70	395285	20.0
.beta.-iso-Methyl...	18.16	3.2	ng	63751	6	16.70	395285	20.0