

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050515\
 Data File : BF078937.D
 Acq On : 5 May 2015 12:59
 Operator : TP/IZ
 Sample : G2115-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUMBO-EP-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-BF043015.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	1.267	5	7	26	rVB3	256381	1120626	20.86%	3.516%
2	1.507	26	28	31	rVB	4061048	5371612	100.00%	16.855%
3	1.656	38	41	50	rVB	313469	548427	10.21%	1.721%
4	1.770	50	51	55	rVV	287801	420390	7.83%	1.319%
5	1.838	55	57	61	rVV	205419	311401	5.80%	0.977%
6	1.907	61	63	87	rVB	1899969	4521564	84.18%	14.188%
7	5.164	346	348	353	rVB	177998	255158	4.75%	0.801%
8	5.656	388	391	393	rBV	1241099	1508566	28.08%	4.734%
9	6.902	497	500	503	rBV	1750733	1638060	30.49%	5.140%
10	7.062	512	514	517	rBV	1706938	1657289	30.85%	5.200%
11	7.359	538	540	542	rBV	403477	397774	7.41%	1.248%
12	7.542	554	556	558	rBV	1444199	1290495	24.02%	4.049%
13	8.045	597	600	602	rBV	1211707	1037172	19.31%	3.254%
14	8.936	675	678	680	rBV	435136	527702	9.82%	1.656%
15	9.633	737	739	742	rBV	90901	87536	1.63%	0.275%
16	10.274	793	795	798	rBV	1889170	2059595	38.34%	6.463%
17	10.445	808	810	812	rVB	61722	56485	1.05%	0.177%
18	10.776	837	839	842	rBV	135346	141275	2.63%	0.443%
19	11.108	866	868	870	rBV	639184	588661	10.96%	1.847%
20	12.011	944	947	949	rBV	417726	401963	7.48%	1.261%
21	12.091	951	954	956	rBV	1091636	1250895	23.29%	3.925%
22	12.537	990	993	995	rBV	144191	168246	3.13%	0.528%
23	12.948	1027	1029	1031	rBV	475641	642088	11.95%	2.015%
24	12.982	1031	1032	1034	rVB	147847	112863	2.10%	0.354%
25	13.200	1048	1051	1054	rBV	149014	198630	3.70%	0.623%
26	14.320	1146	1149	1150	rBV3	53778	90533	1.69%	0.284%
27	14.457	1159	1161	1163	rVB	106307	127359	2.37%	0.400%
28	14.560	1167	1170	1175	rVB	89593	121006	2.25%	0.380%
29	14.743	1181	1186	1189	rBV	144155	167305	3.11%	0.525%
30	14.948	1201	1204	1207	rBV	2265936	2415411	44.97%	7.579%
31	15.165	1215	1223	1224	rBV6	24842	77727	1.45%	0.244%
32	15.440	1245	1247	1253	rVB2	33910	65029	1.21%	0.204%
33	15.543	1253	1256	1259	rBV	81600	107487	2.00%	0.337%
34	15.874	1282	1285	1289	rBV5	20356	57573	1.07%	0.181%

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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

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	16.137	1303	1308	1314	rBV2	163482	300897	5.60%	0.944%
36	16.263	1316	1319	1321	rVV	568269	843851	15.71%	2.648%
37	16.297	1321	1322	1329	rVB2	78512	137901	2.57%	0.433%
38	16.571	1344	1346	1353	rVB4	26579	57785	1.08%	0.181%
39	17.623	1436	1438	1447	rVB	51007	136084	2.53%	0.427%
40	18.034	1472	1474	1477	rBV	50194	69840	1.30%	0.219%
41	18.103	1477	1480	1483	rBV	584775	779657	14.51%	2.446%

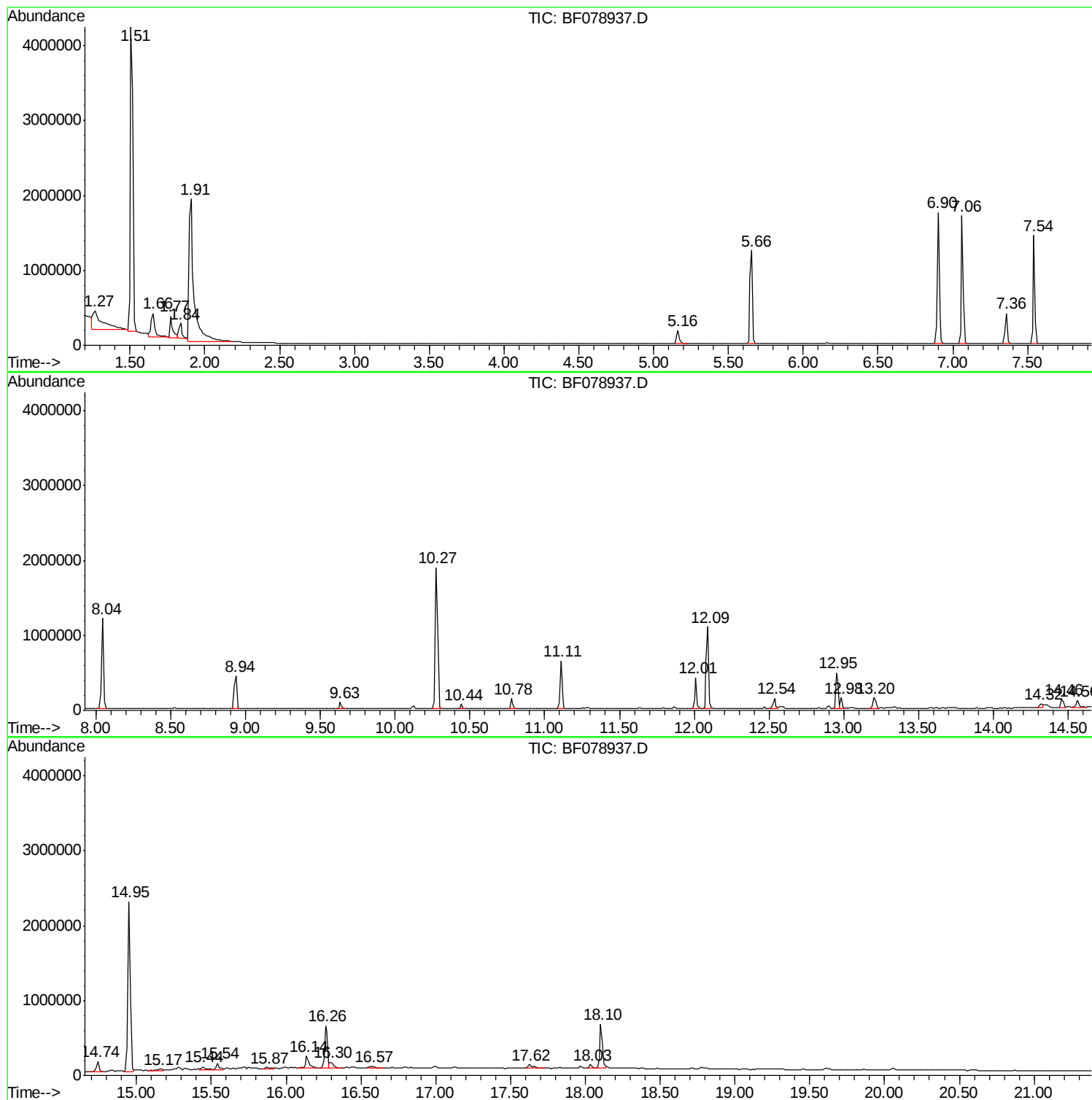
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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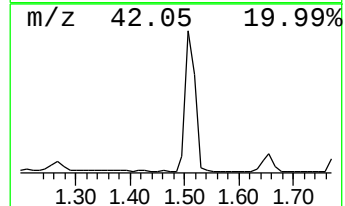
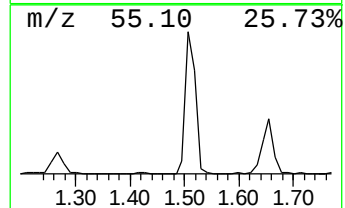
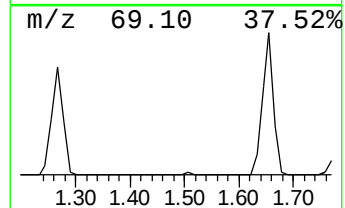
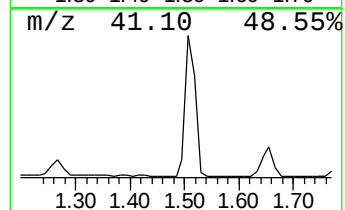
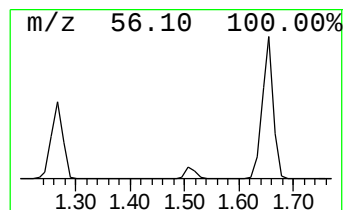
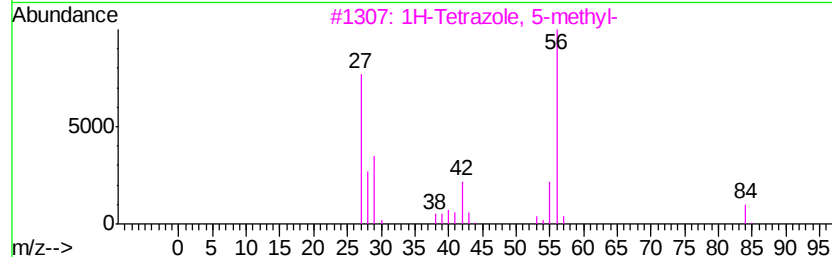
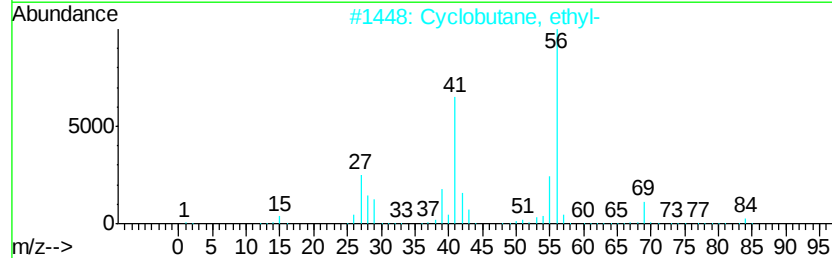
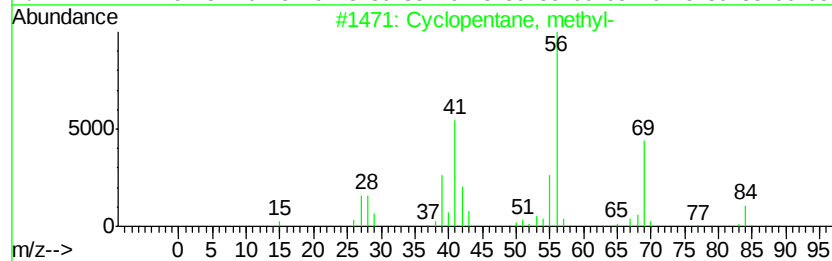
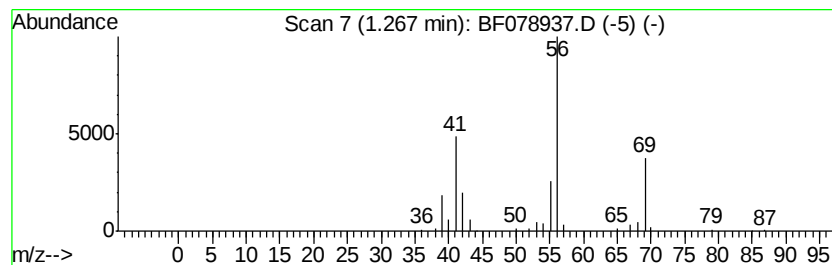
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Cyclopentane, methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.27	56.34 ng	1120630	1,4-Dichlorobenzene-d4	7.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2		Cyclobutane, ethyl-	84	C6H12	004806-61-5	56
3		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	40
4		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	33
5		2-Azetidinone, 3,3-dimethyl-	99	C5H9NO	007486-91-1	9



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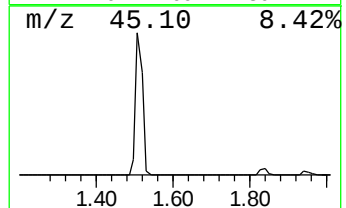
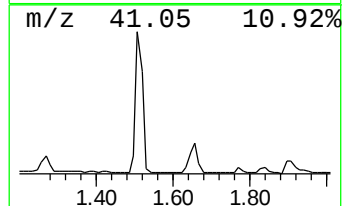
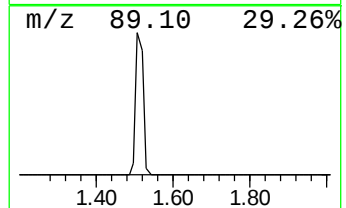
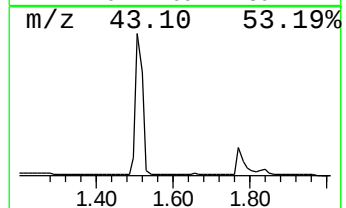
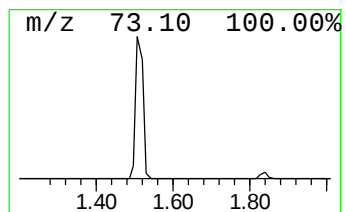
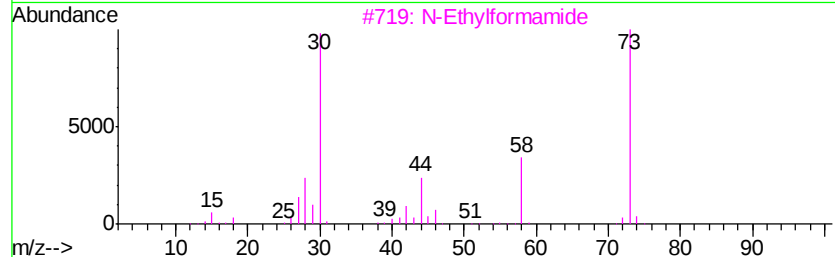
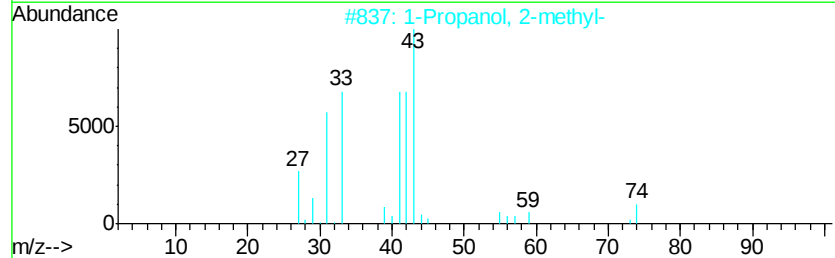
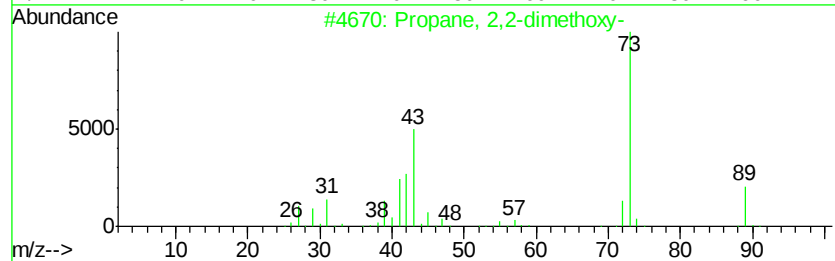
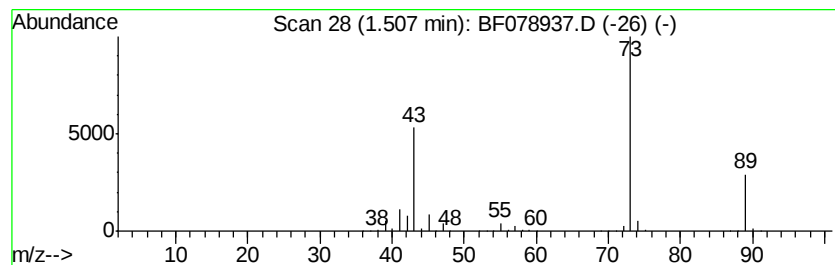
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 Propane, 2,2-dimethoxy- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.51	270.08 ng	5371610	1,4-Dichlorobenzene-d4	7.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	72
2		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
3		N-Ethylformamide	73	C3H7NO	000627-45-2	9
4		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	9
5		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9



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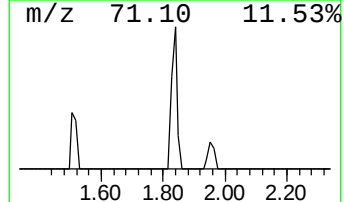
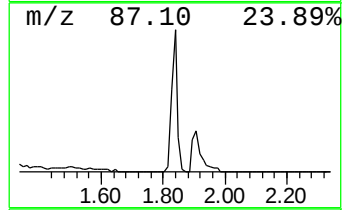
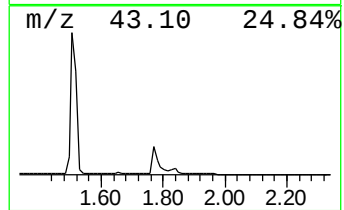
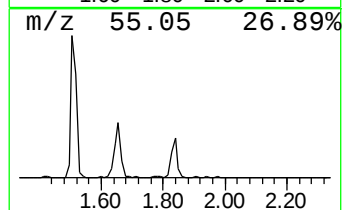
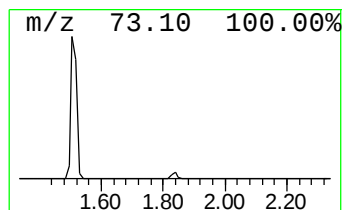
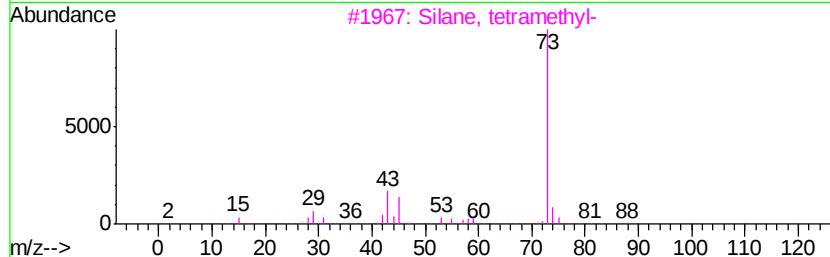
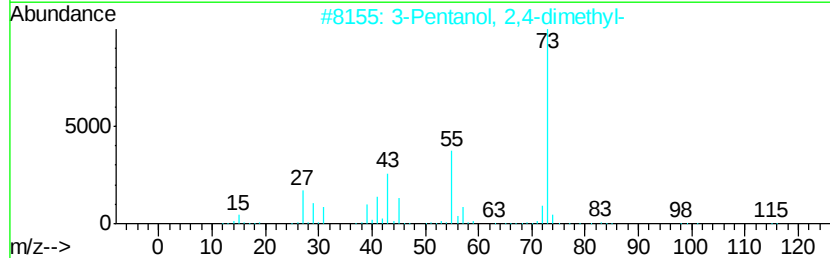
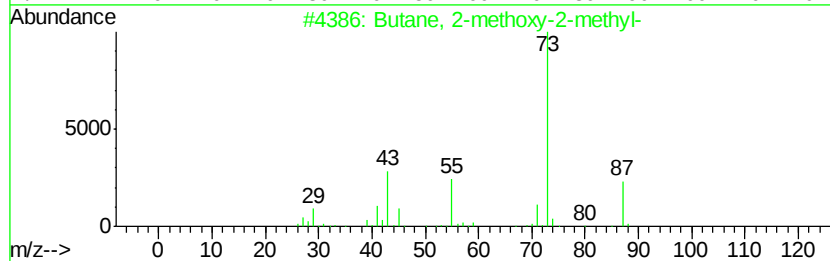
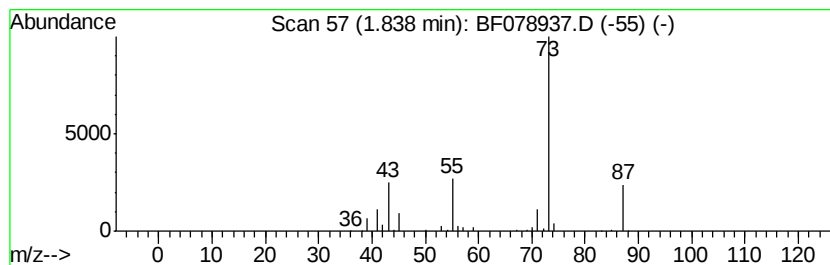
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Butane, 2-methoxy-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.84	15.66 ng	311401	1,4-Dichlorobenzene-d4	7.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	25
3		Silane, tetramethyl-	88	C4H12Si	000075-76-3	12
4		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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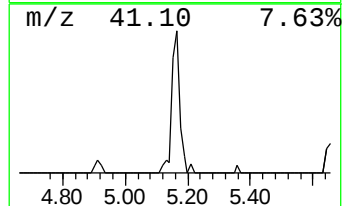
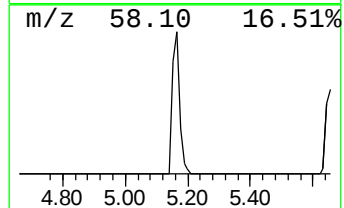
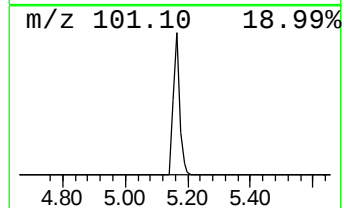
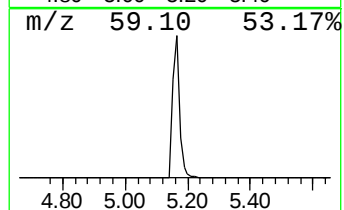
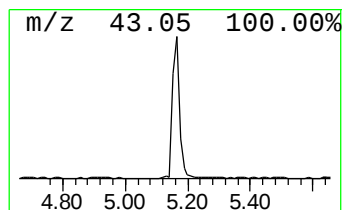
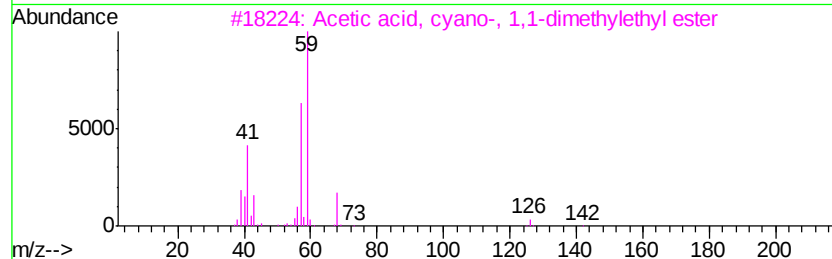
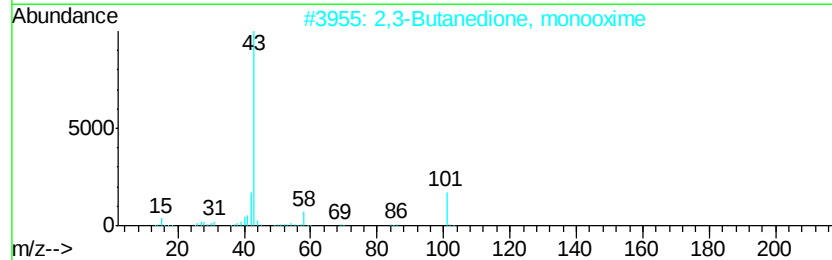
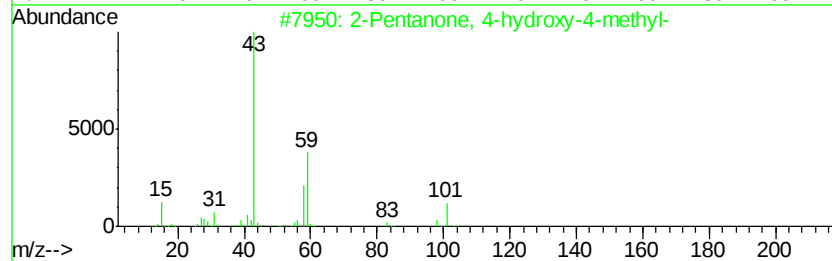
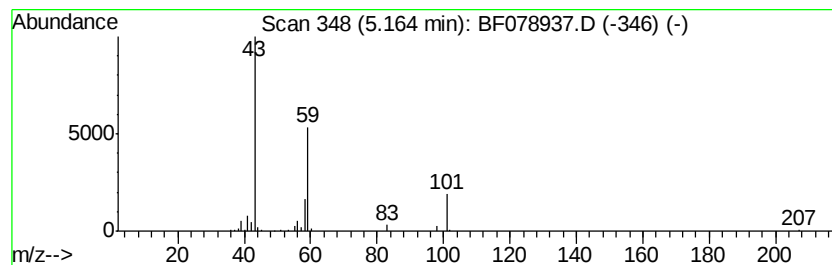
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	12.83 ng	255158	1,4-Dichlorobenzene-d4	7.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	27
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4		5-Hexen-2-one	98	C6H10O	000109-49-9	10
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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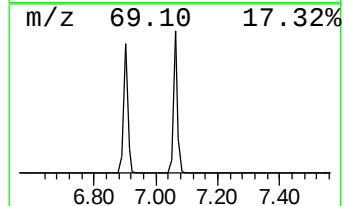
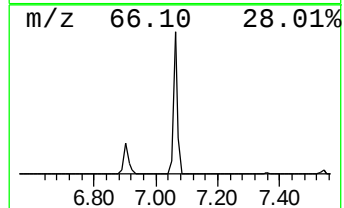
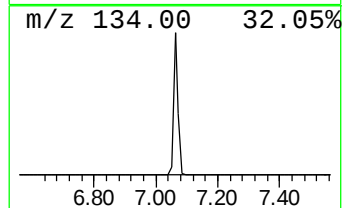
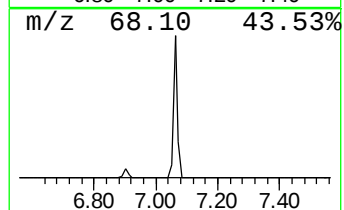
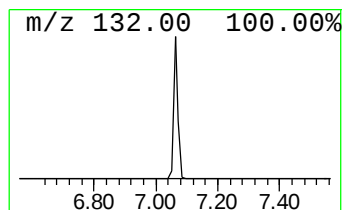
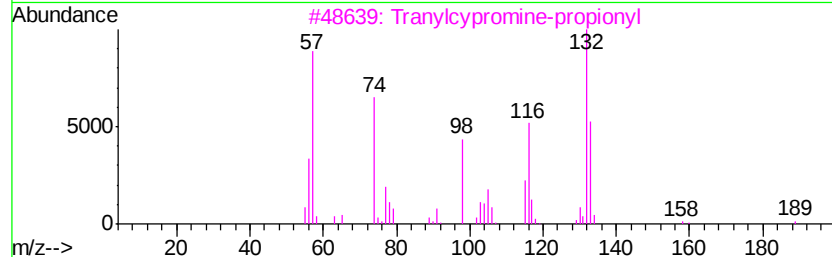
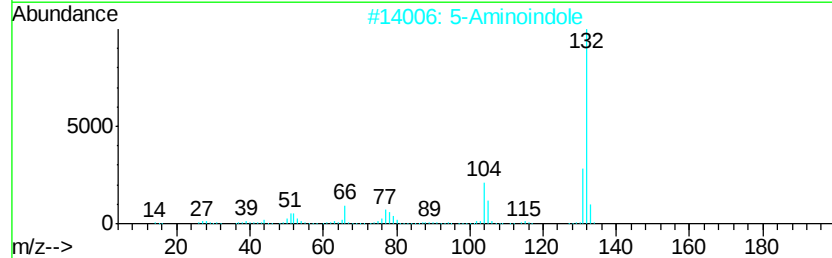
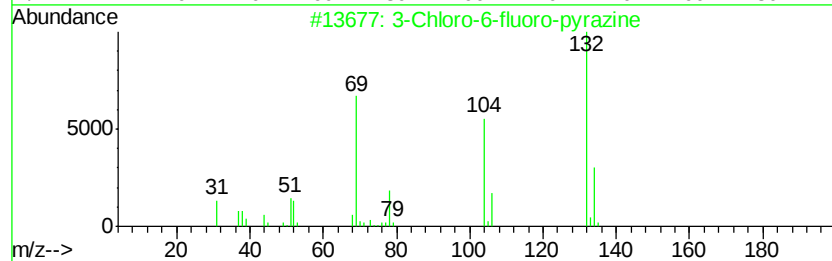
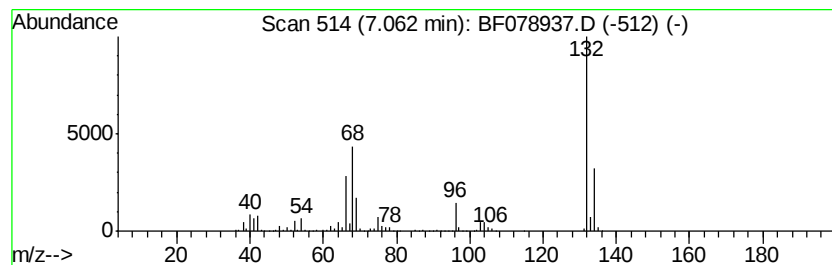
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown7.06 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	83.33 ng	1657290	1,4-Dichlorobenzene-d4	7.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole			132	C8H8N2	005192-03-0	12
3	Tranvlcvpromine-propionyl			189	C12H15NO	1000123-86-3	10
4	Bicyclo[4.2.0]octa-1,3,5-triene,...			132	C10H12	028749-81-7	9
5	5-Fluoro-2-chloropyrimidine			132	C4H2ClFN2	062802-42-0	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050515\
Data File : BF078937.D
Acq On : 5 May 2015 12:59
Operator : TP/IZ
Sample : G2115-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUMBO-EP-6

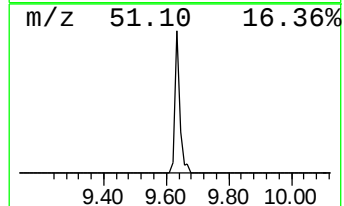
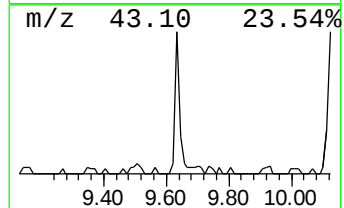
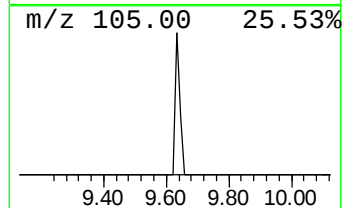
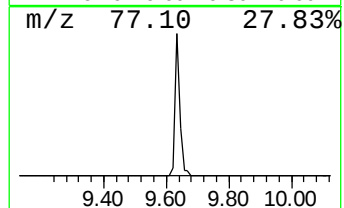
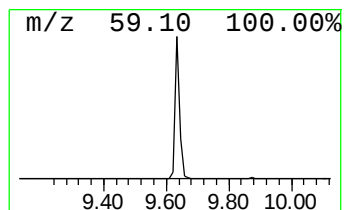
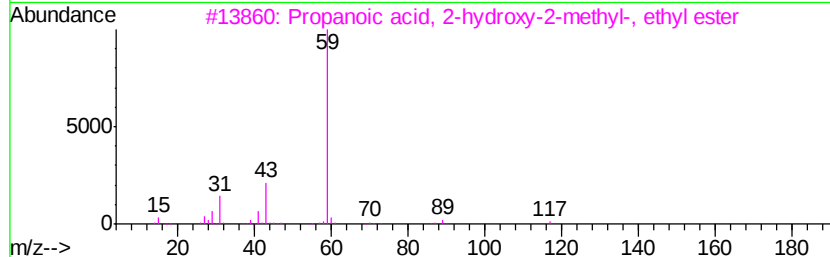
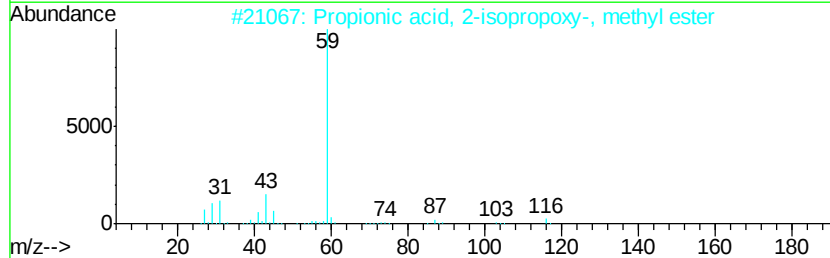
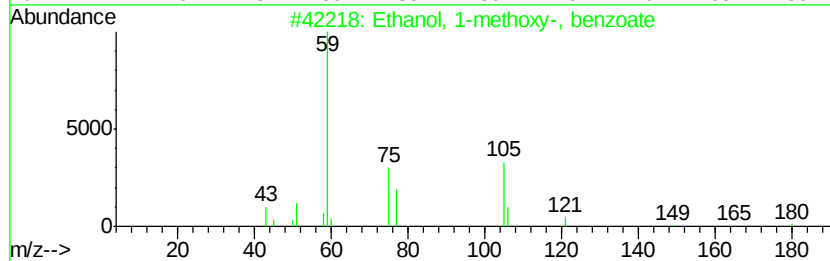
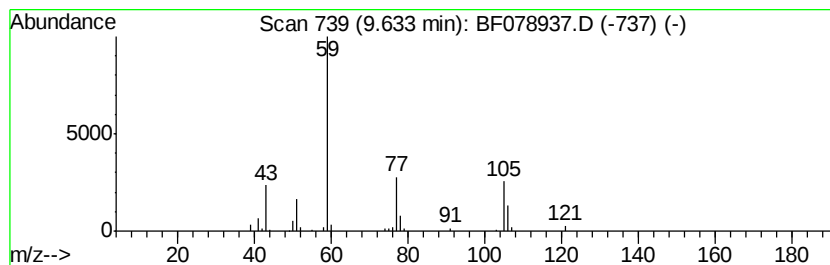
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Ethanol, 1-methoxy-, benzoate Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.63	3.32 ng	87536	Naphthalene-d8	8.94

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 1-methoxy-, benzoate	180	C10H12O3	051835-44-0	72
2		Propionic acid, 2-isopropoxy-, m...	146	C7H14O3	1000151-15-1	23
3		Propanoic acid, 2-hydroxy-2-meth...	132	C6H12O3	000080-55-7	9
4		Acetic acid, chloro-, methyl ester	108	C3H5ClO2	000096-34-4	9
5		Guanidine	59	CH5N3	000113-00-8	7



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050515\
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ClientSampleId :
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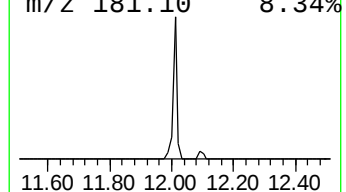
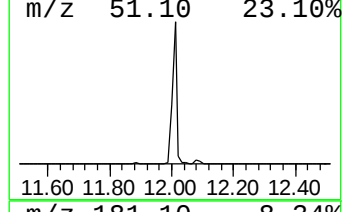
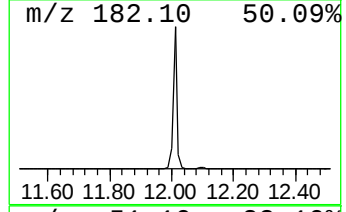
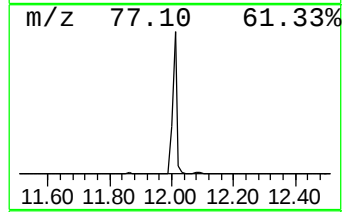
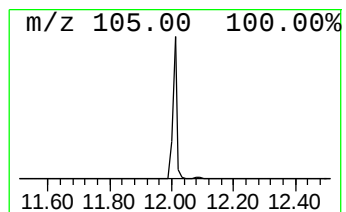
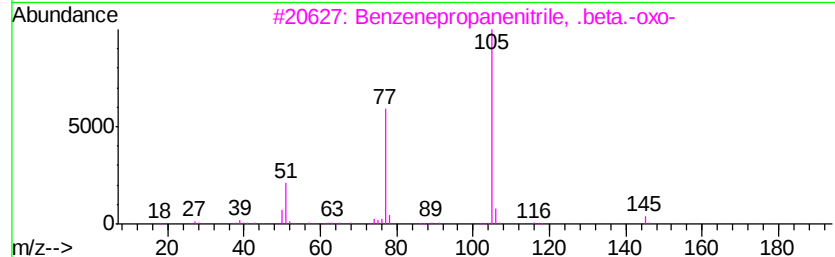
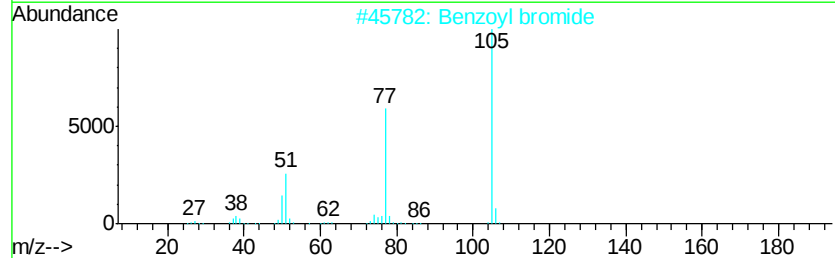
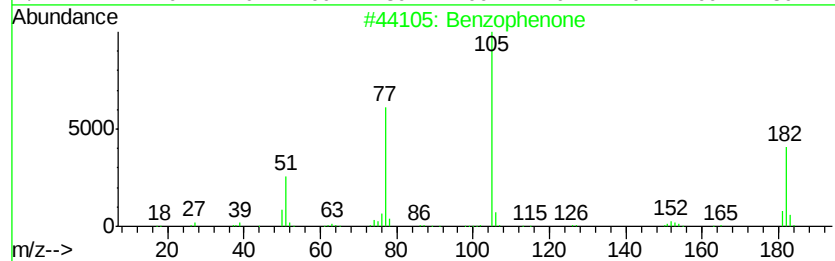
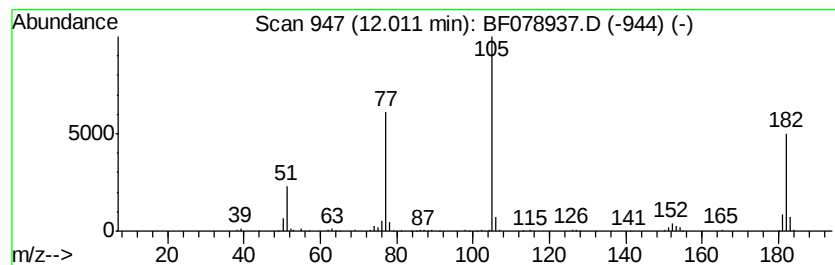
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzophenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.01	13.66 ng	401963	Acenaphthene-d10	11.11

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzophenone	182	C13H10O	000119-61-9	97
2		Benzoyl bromide	184	C7H5BrO	000618-32-6	47
3		Benzenepropanenitrile, .beta.-oxo-	145	C9H7NO	000614-16-4	47
4		1-Propanone, 2-bromo-1-phenyl-	212	C9H9BrO	002114-00-3	47
5		1-Propanone, 3-chloro-1-phenyl-	168	C9H9ClO	000936-59-4	47



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050515\
Data File : BF078937.D
Acq On : 5 May 2015 12:59
Operator : TP/IZ
Sample : G2115-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

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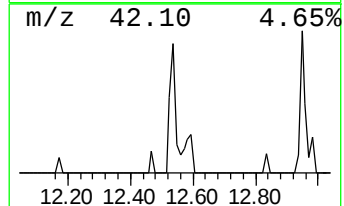
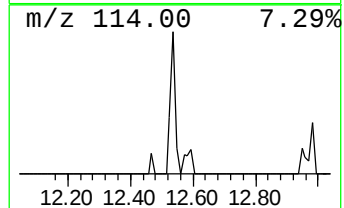
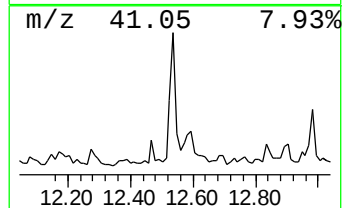
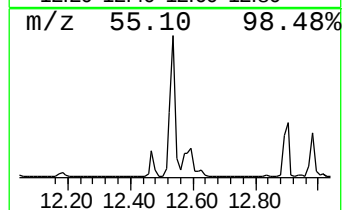
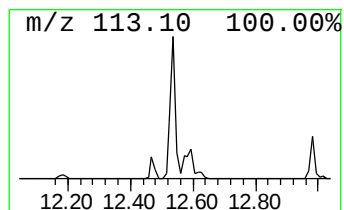
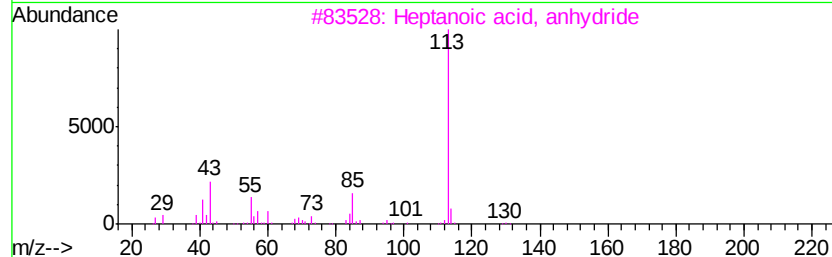
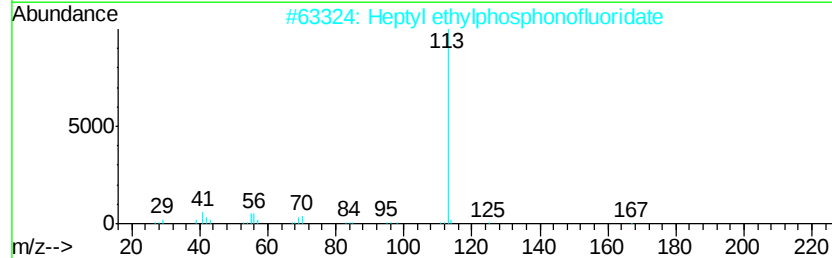
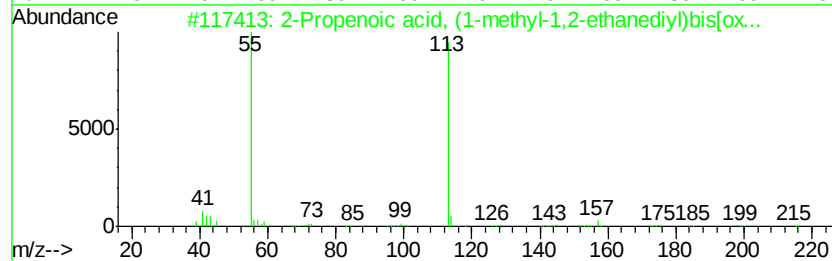
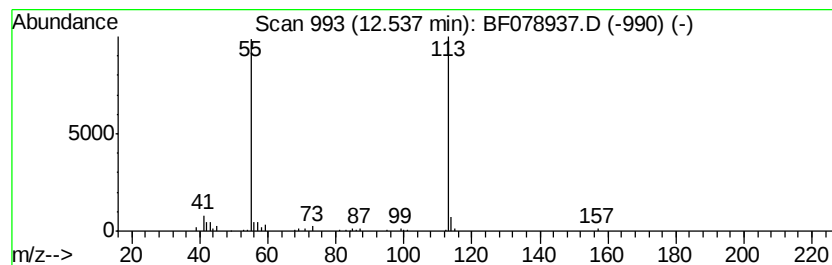
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2-Propenoic acid, (1-methyl... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	5.24 ng	168246	Phenanthrene-d10	12.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propenoic acid, (1-methyl-1,2-...	300	C15H24O6	042978-66-5	78
2		Heptyl ethylphosphonofluoridate	210	C9H20F02P	162085-85-0	36
3		Heptanoic acid, anhydride	242	C14H26O3	000626-27-7	17
4		Propanedioic acid, 2-propenyl-, ...	172	C8H12O4	040637-56-7	9
5		1H-Pyrazole, 4-nitro-	113	C3H3N3O2	002075-46-9	9



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050515\
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Operator : TP/IZ
Sample : G2115-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

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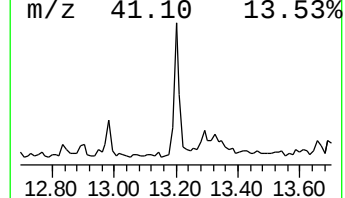
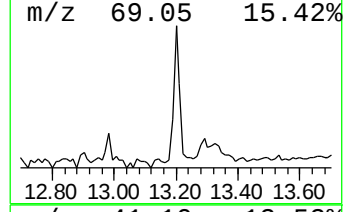
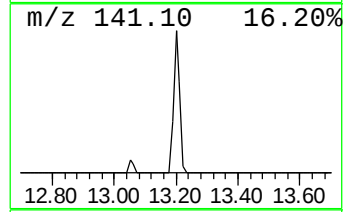
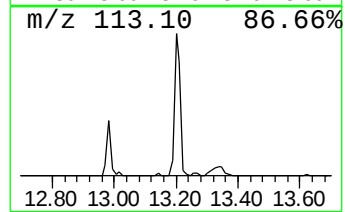
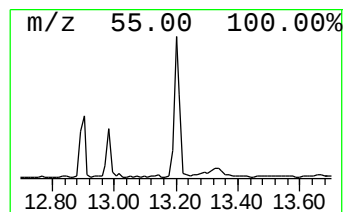
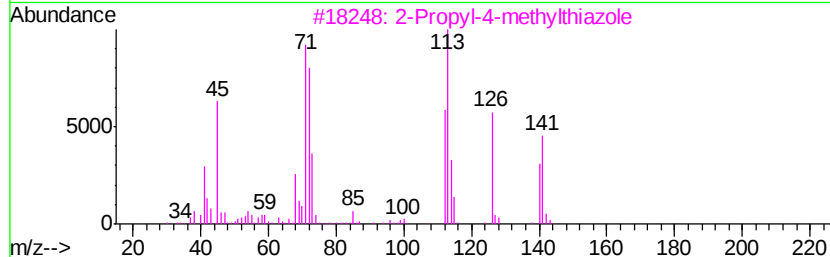
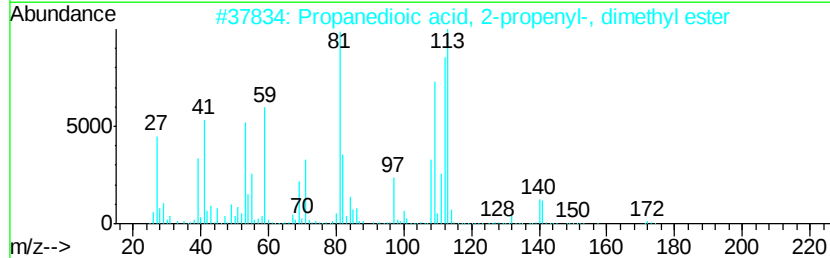
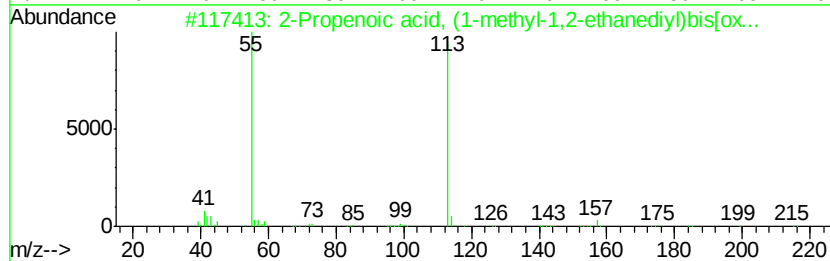
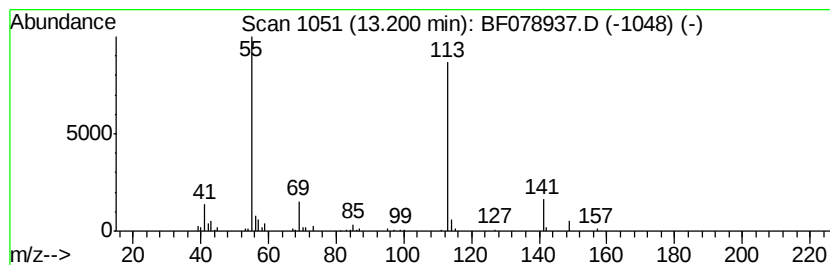
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown13.20 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.20	6.19 ng	198630	Phenanthrene-d10	12.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenoic acid, (1-methyl-1,2-...	300	C15H24O6	042978-66-5	59		
2	Propanedioic acid, 2-propenyl-, ...	172	C8H12O4	040637-56-7	38		
3	2-Propyl-4-methylthiazole	141	C7H11NS	052414-87-6	38		
4	Heptanoic acid, anhydride	242	C14H26O3	000626-27-7	37		
5	3-Methylpentyl ethylphosphonoflu...	196	C8H18F02P	1000298-39-1	36		



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050515\
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Misc :
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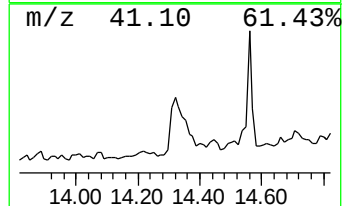
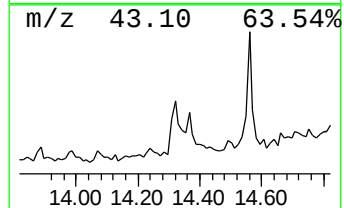
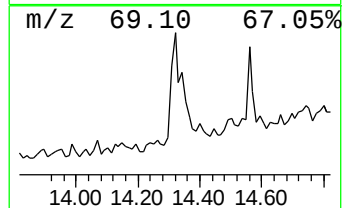
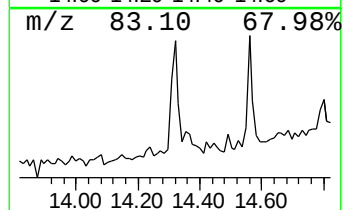
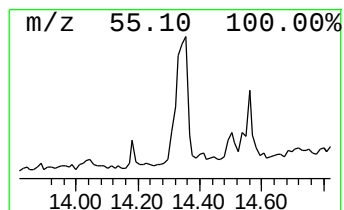
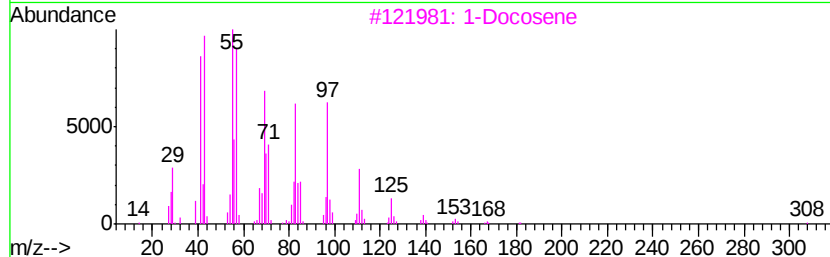
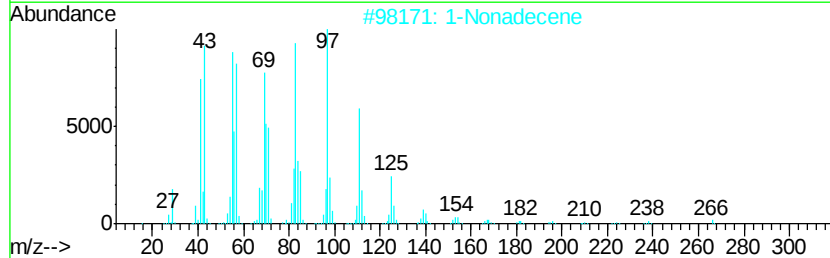
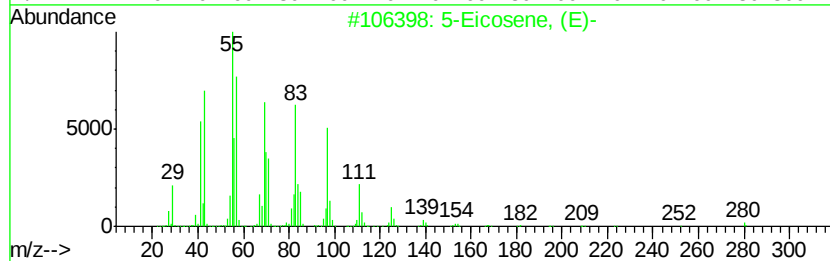
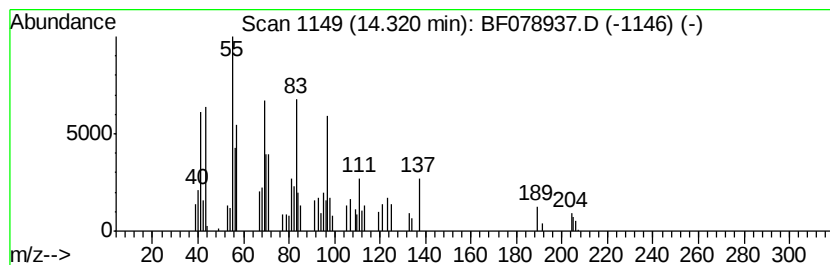
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 5-Eicosene, (E)- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.32	2.82 ng	90533	Phenanthrene-d10	12.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Eicosene, (E)-	280	C20H40	074685-30-6	80
2			1-Nonadecene	266	C19H38	018435-45-5	80
3			1-Docosene	308	C22H44	001599-67-3	80
4			3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	64
5			1-Hexadecanol	242	C16H34O	036653-82-4	58



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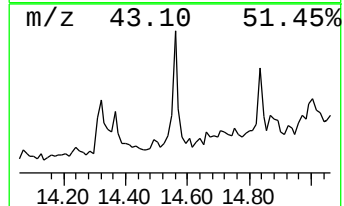
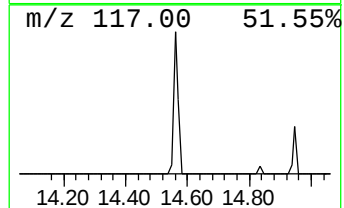
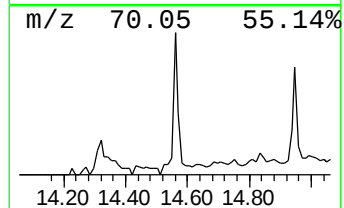
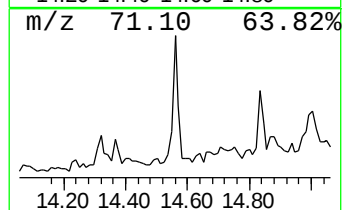
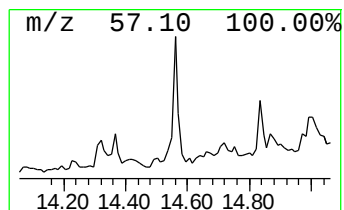
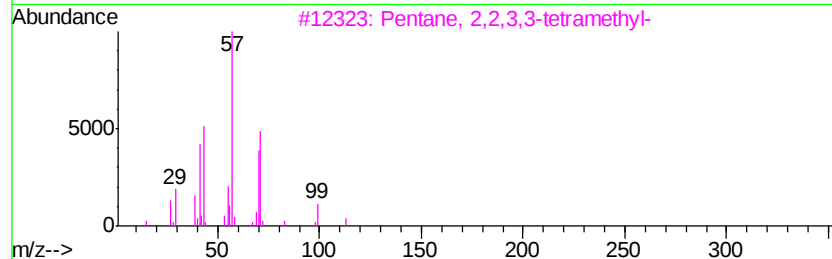
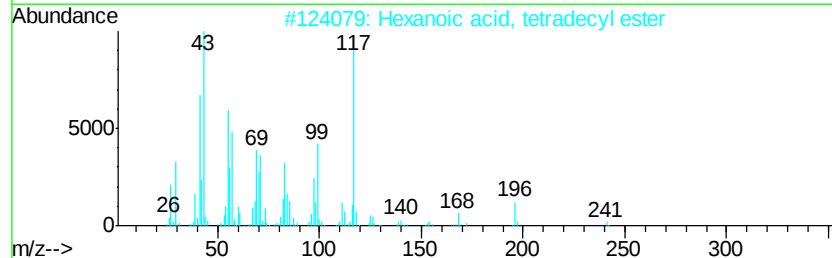
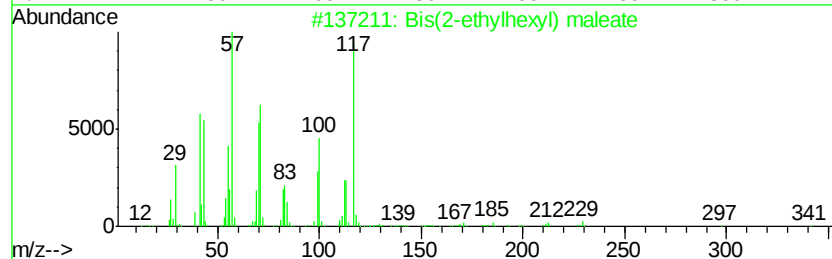
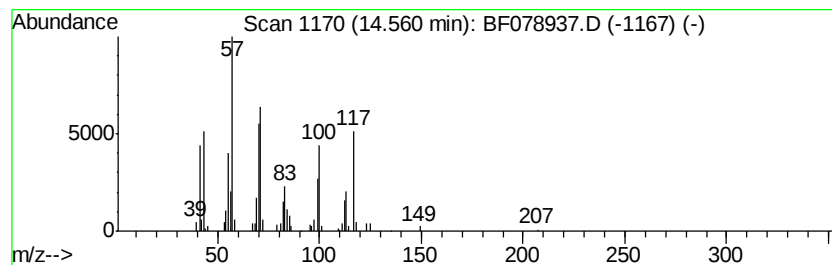
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Bis(2-ethylhexyl) maleate Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.56	3.77 ng	121006	Phenanthrene-d10	12.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Bis(2-ethylhexyl) maleate	340	C20H36O4	000142-16-5	90
2		Hexanoic acid, tetradecyl ester	312	C20H40O2	071801-23-5	35
3		Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	35
4		Acetic acid, cyano-, 2-ethylhexyl-	197	C11H19NO2	013361-34-7	27
5		dl-2-Ethylhexyl chloroformate	192	C9H17ClO2	024468-13-1	27



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Data File : BF078937.D
Acq On : 5 May 2015 12:59
Operator : TP/IZ
Sample : G2115-02
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ClientSampleId :
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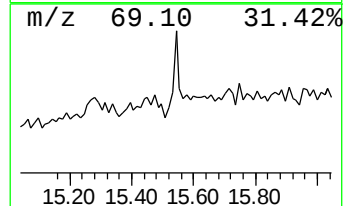
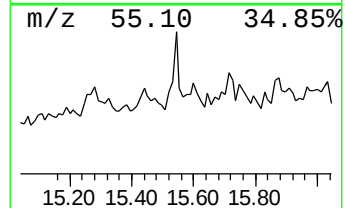
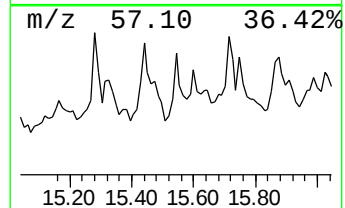
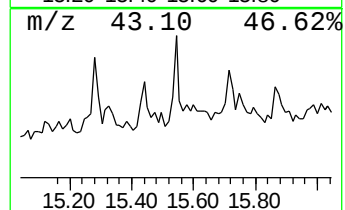
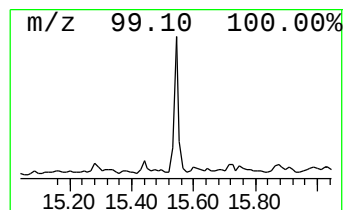
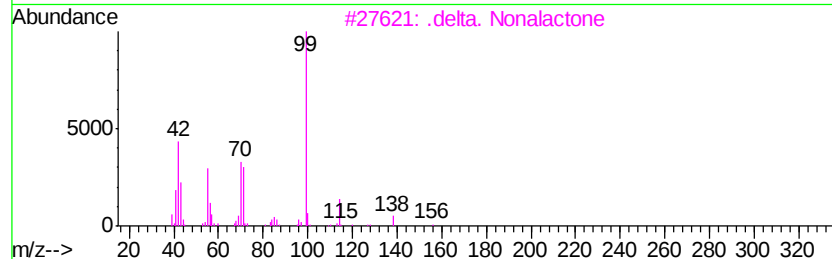
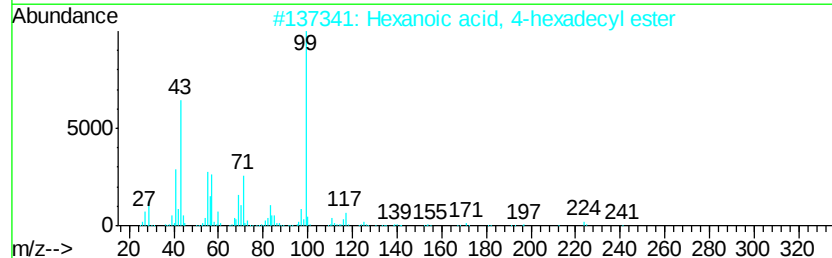
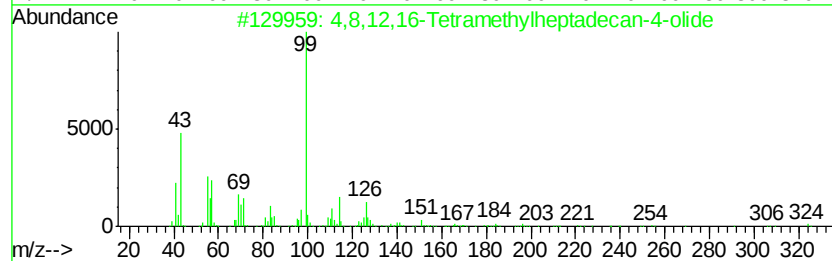
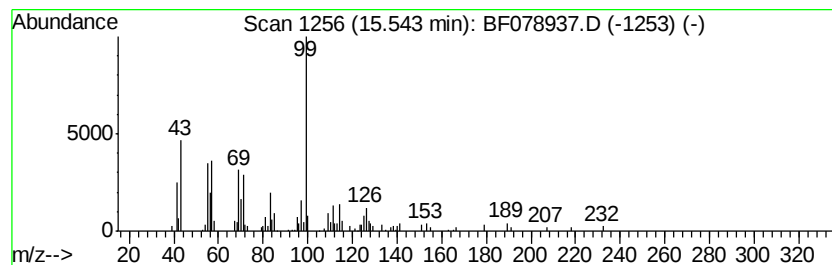
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 4,8,12,16-Tetramethylheptad... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.54	2.55 ng	107487	Chrysene-d12	16.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4,8,12,16-Tetramethylheptadecan-...	324	C21H40O2	096168-15-9	80
2			Hexanoic acid, 4-hexadecyl ester	340	C22H44O2	1000282-83-8	59
3			.delta. Nonalactone	156	C9H16O2	003301-94-8	47
4			2(3H)-Furanone, 5-hexyldihydro-4...	184	C11H20O2	125133-98-4	47
5			2H-Pyran-2-one, 6-heptyltetrahydro-	198	C12H22O2	000713-95-1	47



Data Path : Z:\HPCHEM1\BNA F\DATA\BF050515\
Data File : BF078937.D
Acq On : 5 May 2015 12:59
Operator : TP/IZ
Sample : G2115-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUMBO-EP-6

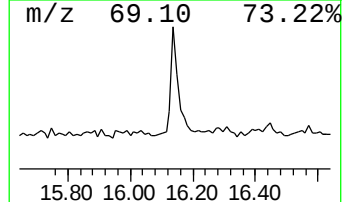
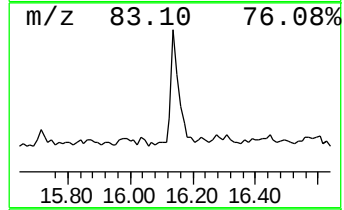
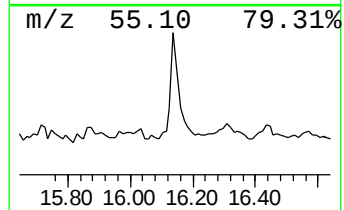
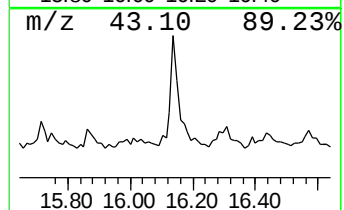
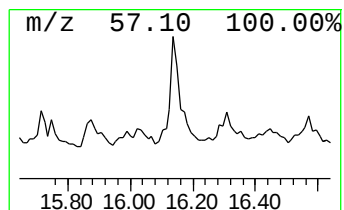
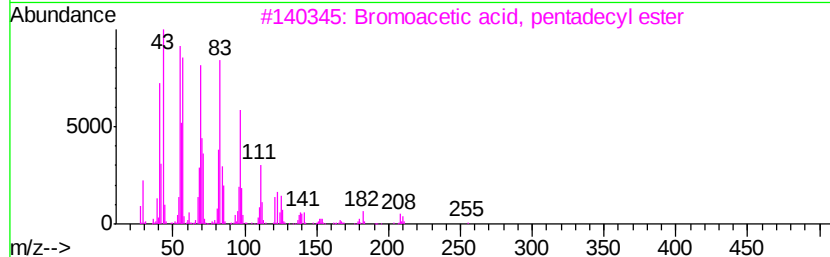
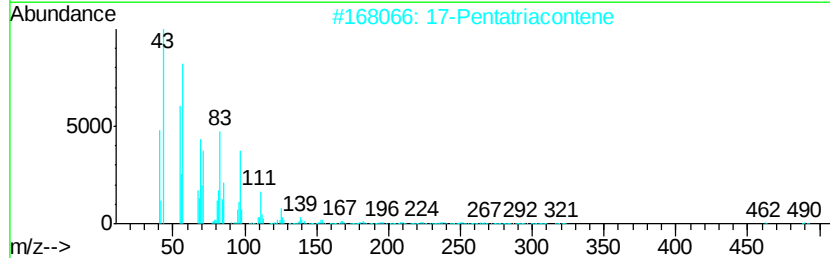
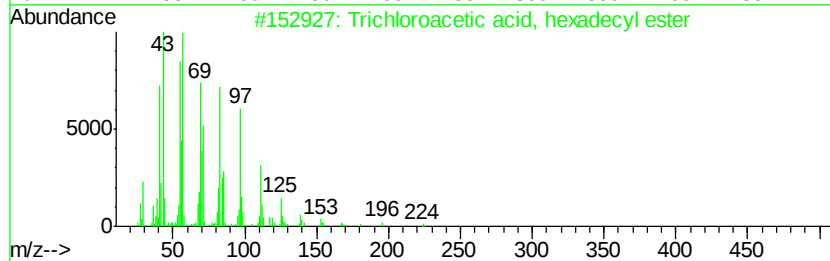
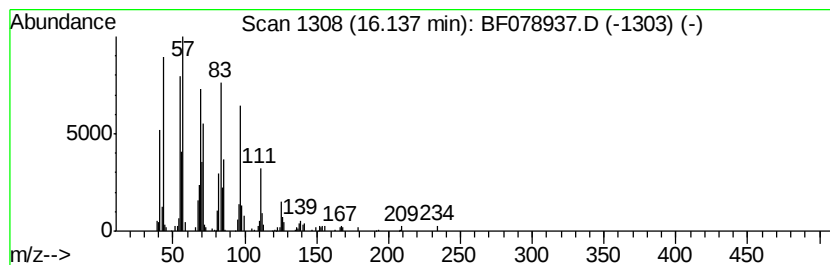
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Trichloroacetic acid, hexad... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.14	7.13 ng	300897	Chrysene-d12	16.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
2		17-Pentatriacontene	491	C35H70	006971-40-0	91
3		Bromoacetic acid, pentadecyl ester	348	C17H33BrO2	131143-01-6	90
4		3-Eicosene, (E)-	280	C20H40	074685-33-9	90
5		9-Eicosene, (E)-	280	C20H40	074685-29-3	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF050515\
 Data File : BF078937.D
 Acq On : 5 May 2015 12:59
 Operator : TP/IZ
 Sample : G2115-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUMBO-EP-6

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.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
Cyclopentane, met...	1.27	56.3	ng	1120630	1	7.36	397774	20.0
Propane, 2,2-dime...	1.51	270.1	ng	5371610	1	7.36	397774	20.0
Butane, 2-methoxy...	1.84	15.7	ng	311401	1	7.36	397774	20.0
2-Pentanone, 4-hy...	5.16	12.8	ng	255158	1	7.36	397774	20.0
unknown7.06	7.06	83.3	ng	1657290	1	7.36	397774	20.0
Ethanol, 1-methox...	9.63	3.3	ng	87536	2	8.94	527702	20.0
Benzophenone	12.01	13.7	ng	401963	3	11.11	588661	20.0
2-Propenoic acid, ...	12.54	5.2	ng	168246	4	12.95	642088	20.0
unknown13.20	13.20	6.2	ng	198630	4	12.95	642088	20.0
5-Eicosene, (E)-	14.32	2.8	ng	90533	4	12.95	642088	20.0
Bis(2-ethylhexyl)...	14.56	3.8	ng	121006	4	12.95	642088	20.0
4,8,12,16-Tetrame...	15.54	2.5	ng	107487	5	16.27	843851	20.0
Trichloroacetic a...	16.14	7.1	ng	300897	5	16.27	843851	20.0