

Data Path : S:\HPCHEM1\BNA_F\DATA\BF021815\
 Data File : BF077355.D
 Acq On : 18 Feb 2015 16:40
 Operator : TP/IZ
 Sample : G1300-03
 Misc : S
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-04-005-B

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-BF021715.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	1.516	52	55	58	rVB	4430937	4993994	100.00%	22.859%
2	1.653	64	67	75	rVB	105963	212222	4.25%	0.971%
3	1.767	75	77	81	rVV	89383	181796	3.64%	0.832%
4	1.836	81	83	87	rVB	145587	187048	3.75%	0.856%
5	1.904	87	89	102	rVB	682519	979883	19.62%	4.485%
6	5.150	371	373	376	rBV	360254	412891	8.27%	1.890%
7	5.642	413	416	419	rBV	1209566	1191948	23.87%	5.456%
8	6.899	523	526	528	rBV	1430260	1365778	27.35%	6.252%
9	7.059	537	540	543	rBV	1436777	1331849	26.67%	6.096%
10	7.356	563	566	568	rBV	228986	277872	5.56%	1.272%
11	7.539	579	582	584	rBV	1191556	1041389	20.85%	4.767%
12	8.042	623	626	628	rBV	925211	802240	16.06%	3.672%
13	8.933	701	704	706	rBV	361553	410054	8.21%	1.877%
14	10.282	819	822	824	rVB	1362193	1678554	33.61%	7.683%
15	11.105	892	894	897	rVB	498267	484230	9.70%	2.217%
16	12.088	977	980	982	rBV	1029121	1053539	21.10%	4.822%
17	12.957	1053	1056	1057	rBV	473720	528804	10.59%	2.421%
18	13.665	1116	1118	1120	rBV	98629	117501	2.35%	0.538%
19	14.465	1185	1188	1190	rBV	80510	97797	1.96%	0.448%
20	14.648	1202	1204	1209	rBV	77679	128364	2.57%	0.588%
21	14.740	1209	1212	1214	rVV	80421	108072	2.16%	0.495%
22	14.957	1228	1231	1233	rBV	1683384	1980808	39.66%	9.067%
23	15.665	1291	1293	1295	rBV	47284	54213	1.09%	0.248%
24	16.134	1332	1334	1340	rBV	288308	385209	7.71%	1.763%
25	16.271	1343	1346	1348	rBV	562919	653038	13.08%	2.989%
26	17.003	1407	1410	1413	rBV	45622	95602	1.91%	0.438%
27	17.620	1462	1464	1466	rBV	48477	91664	1.84%	0.420%
28	17.826	1480	1482	1488	rBV2	37395	117981	2.36%	0.540%
29	18.020	1497	1499	1501	rVB	43605	54333	1.09%	0.249%
30	18.089	1502	1505	1509	rVV	403642	621589	12.45%	2.845%
31	18.763	1562	1564	1566	rVB	42459	65217	1.31%	0.299%
32	19.243	1604	1606	1608	rBV2	29215	59588	1.19%	0.273%
33	20.283	1694	1697	1699	rBV4	32748	81465	1.63%	0.373%

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ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
FK-GF-04-005-B

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-BF021715.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

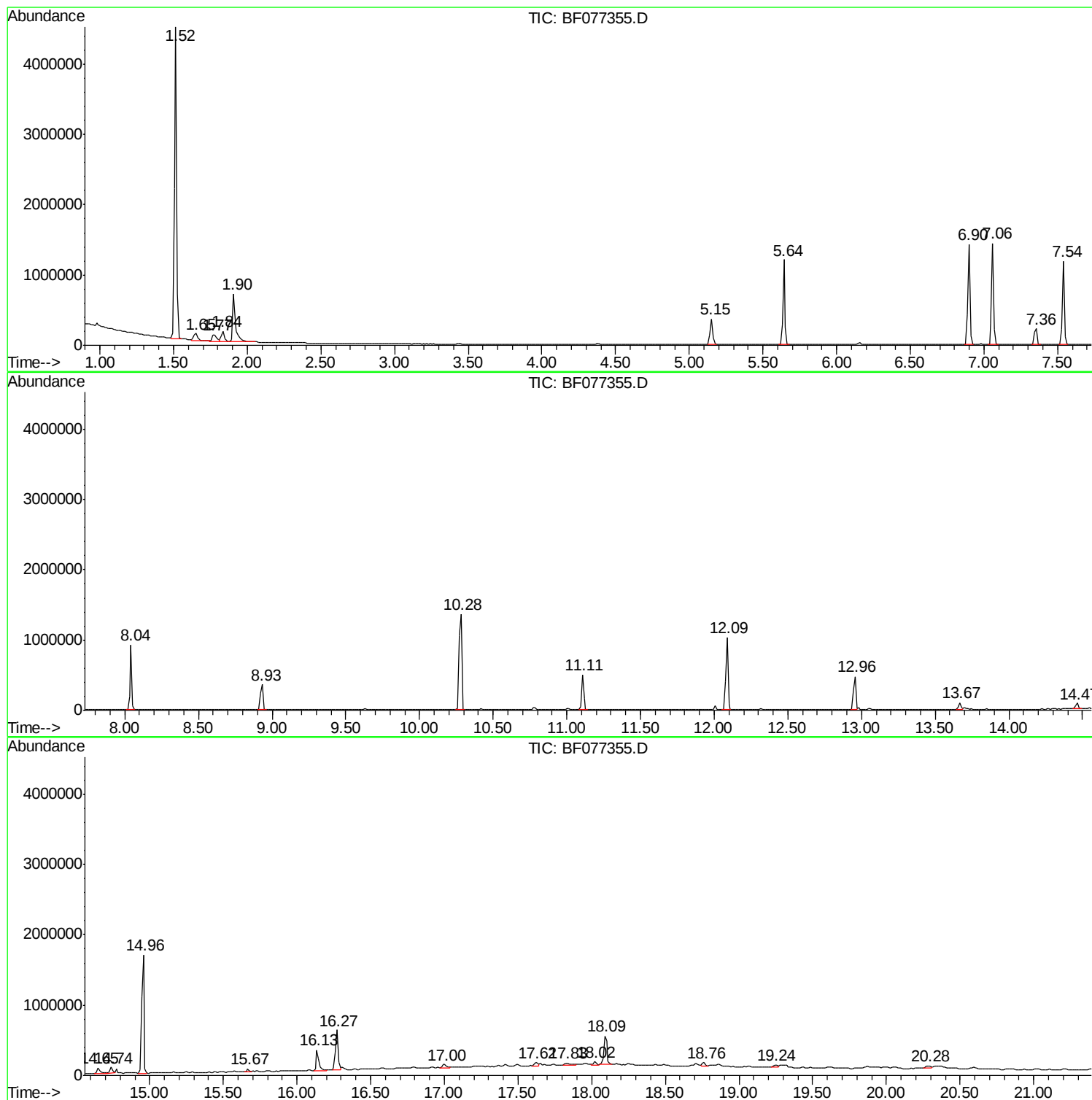
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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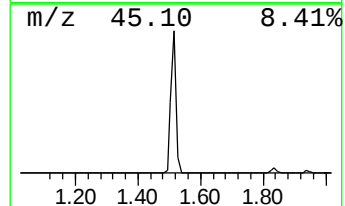
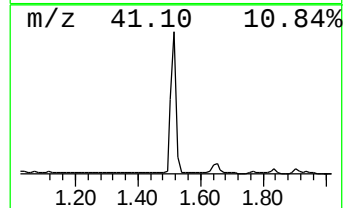
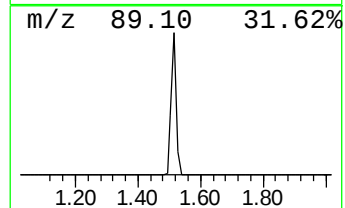
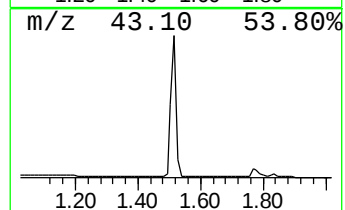
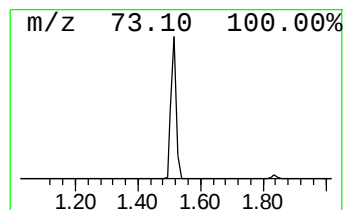
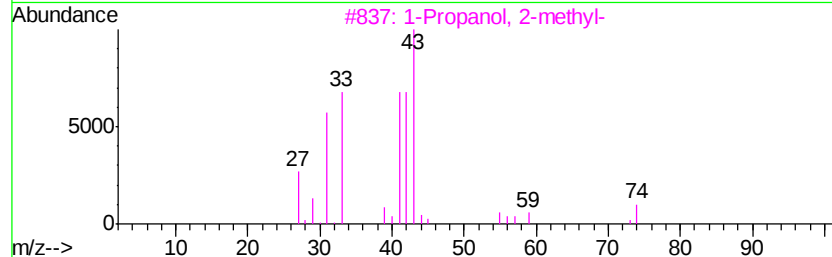
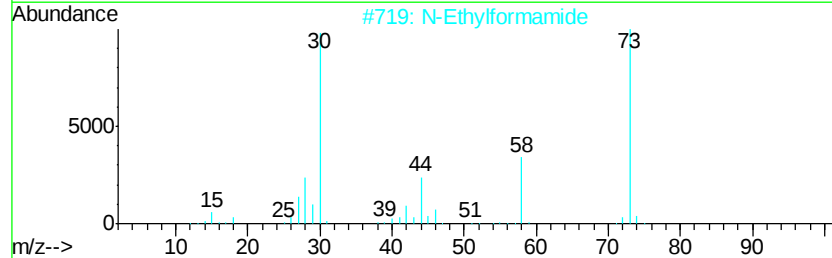
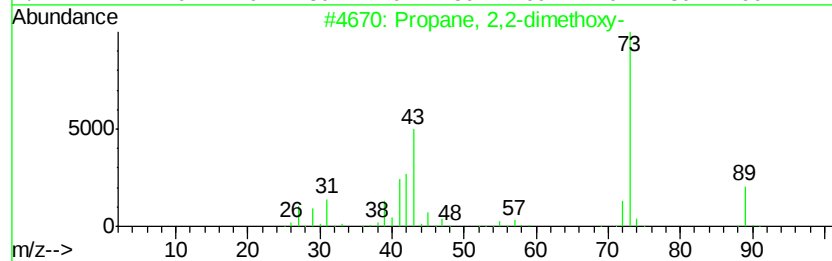
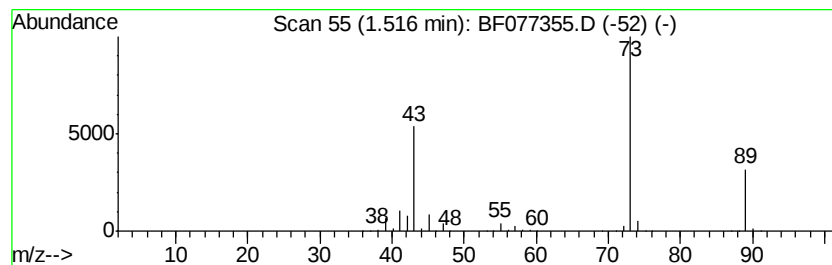
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown1.52 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.52	359.45 ng	4993990	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	38
2		N-Ethylformamide	73	C3H7NO	000627-45-2	9
3		1-Propanol, 2-methyl-	74	C4H10O	000078-83-1	9
4		1,3-Diaminoguanidine	89	CH7N5	004364-78-7	9
5		Propane, 2-methoxy-2-methyl-	88	C5H12O	001634-04-4	9



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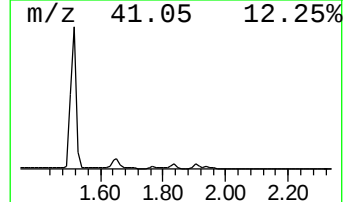
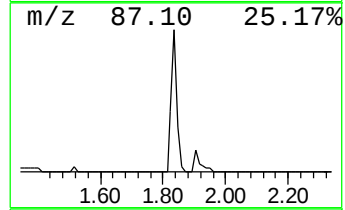
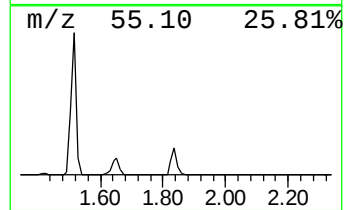
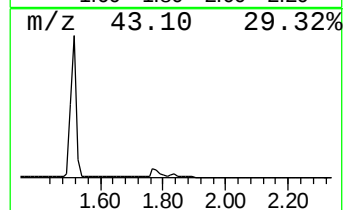
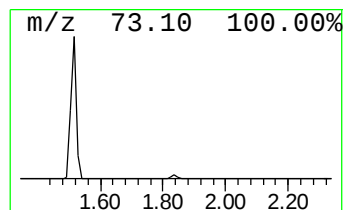
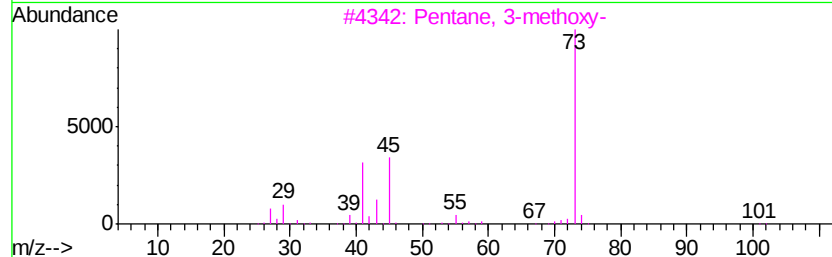
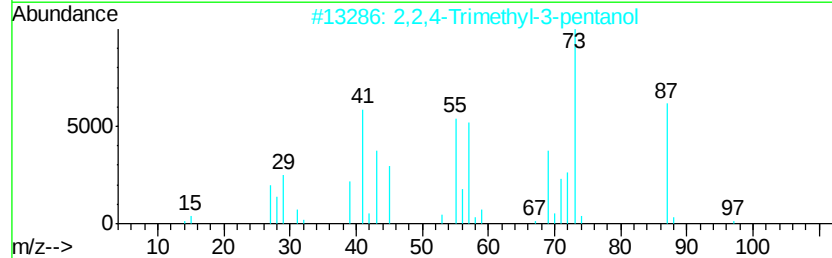
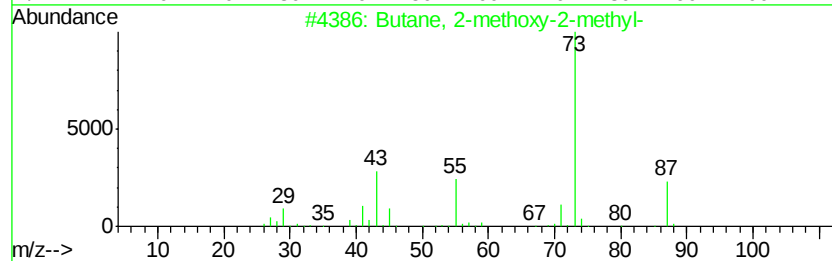
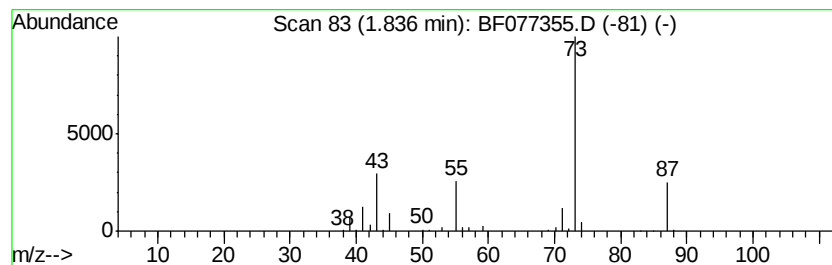
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.84	13.46 ng	187048	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	90
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	43
3		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4		Silane, tetramethyl-	88	C4H12Si	000075-76-3	16
5		N-Ethylformamide	73	C3H7NO	000627-45-2	9



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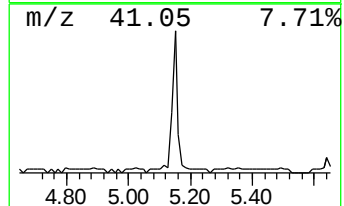
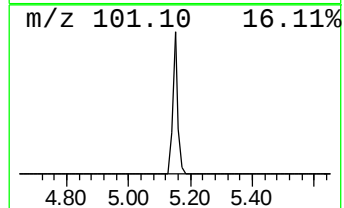
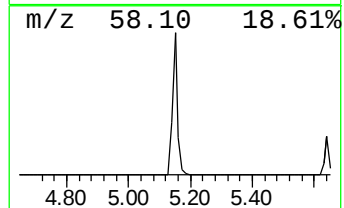
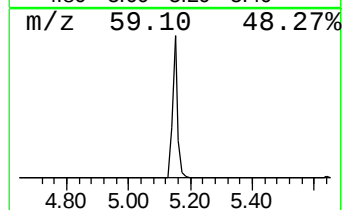
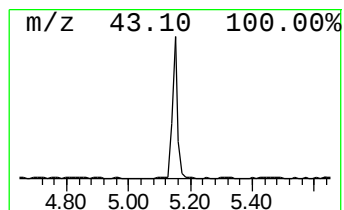
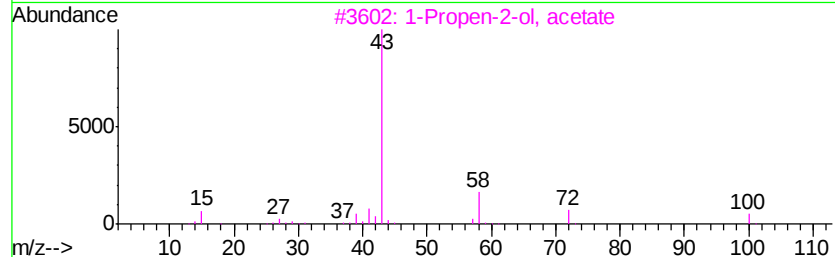
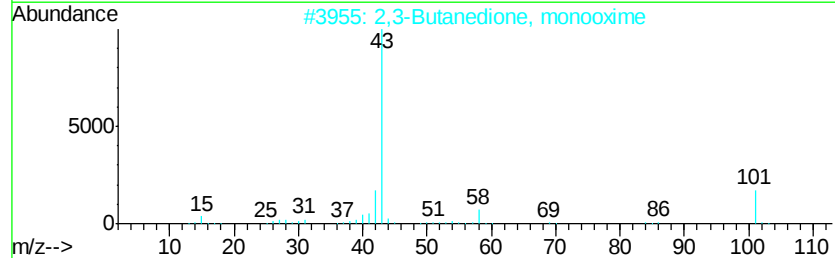
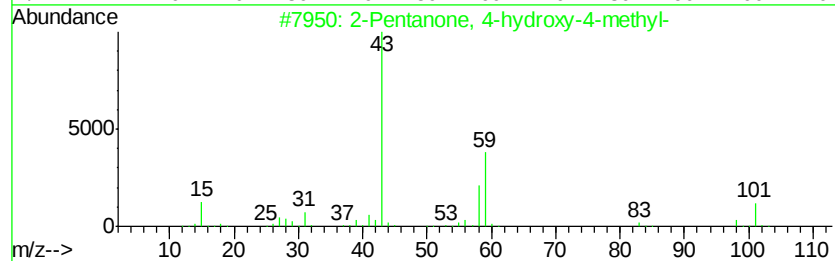
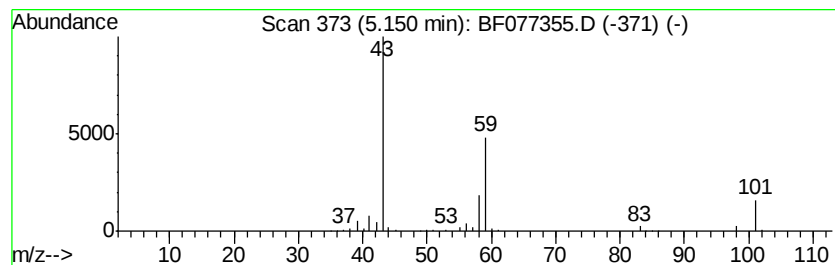
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.15	29.72 ng	412891	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	16
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10
4		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	9
5		Thiocyanic acid, propyl ester	101	C4H7NS	004251-16-5	7



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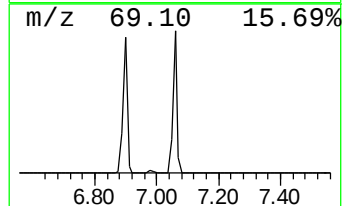
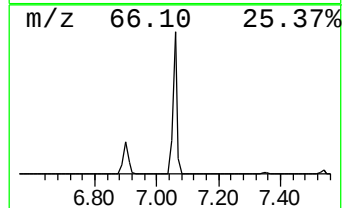
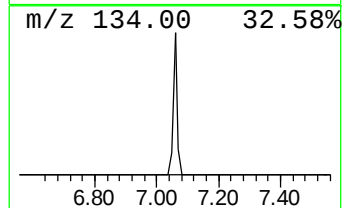
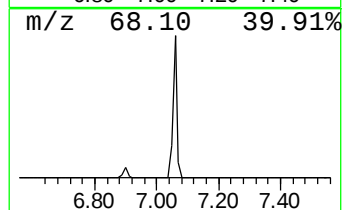
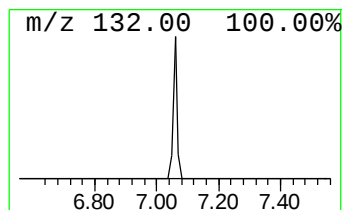
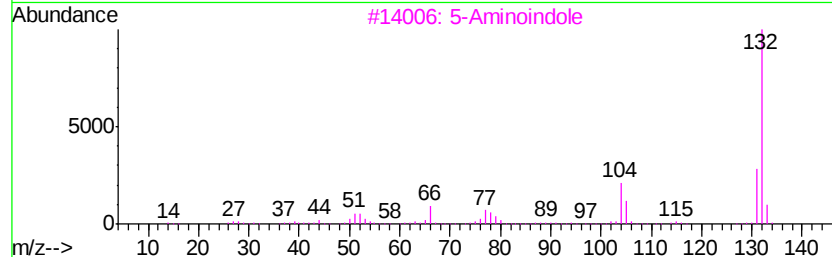
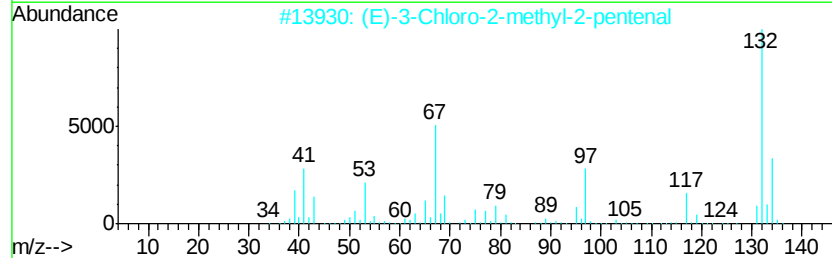
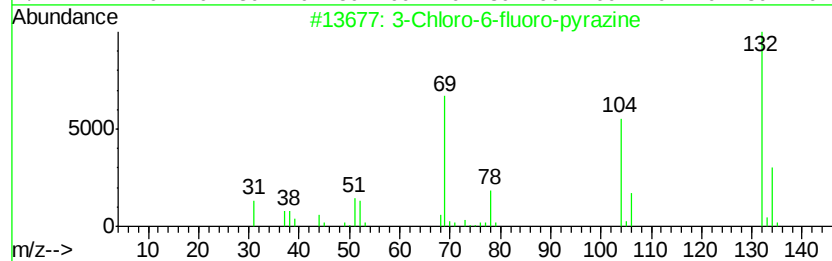
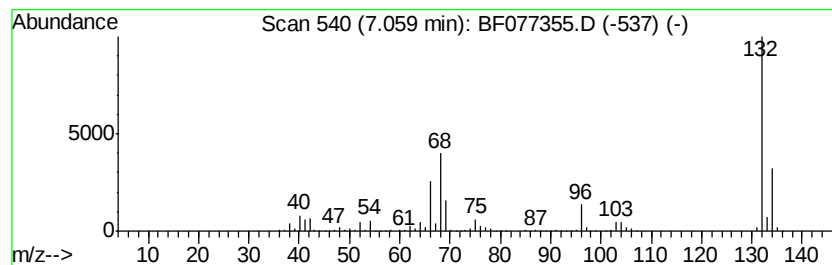
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 7 unknown7.06 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.06	95.86 ng	1331850	1,4-Dichlorobenzene-d4	7.36

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	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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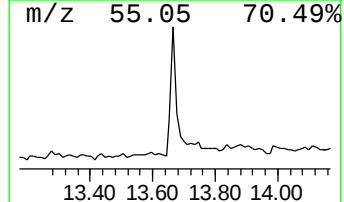
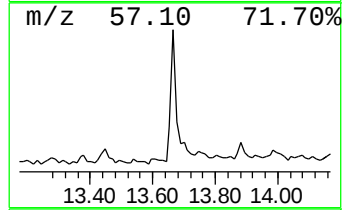
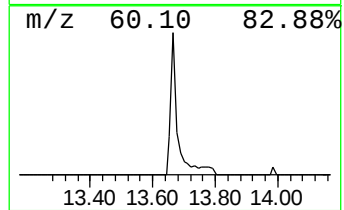
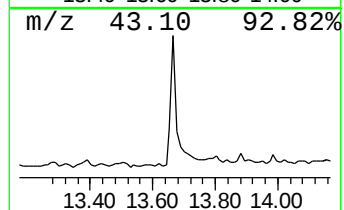
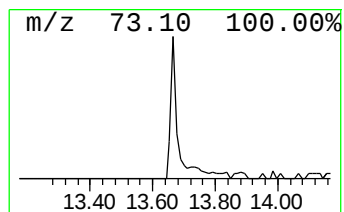
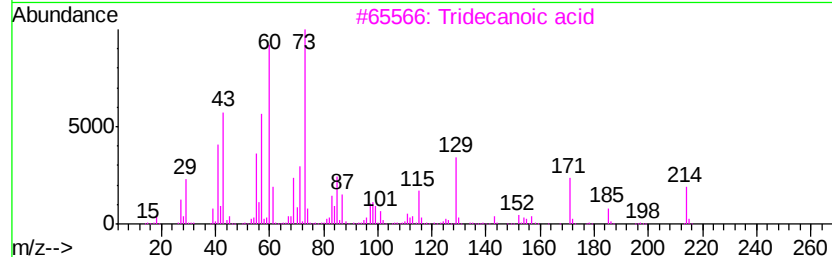
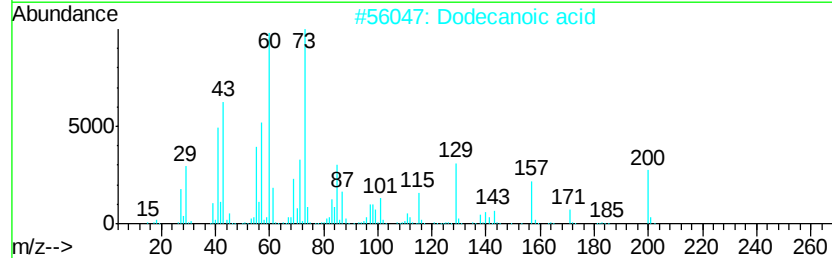
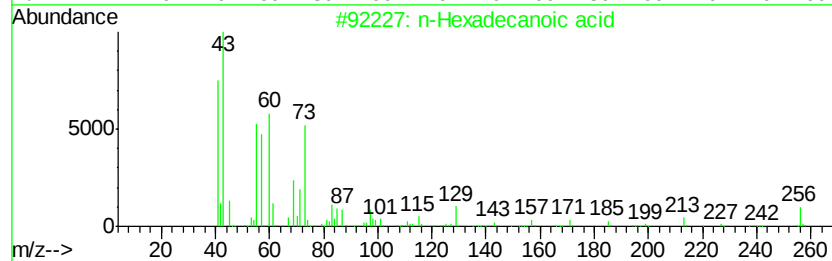
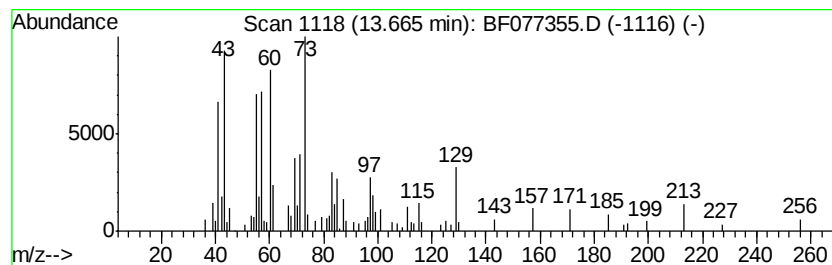
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 9 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.67	4.44 ng	117501	Phenanthrene-d10	12.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
2		Dodecanoic acid	200	C12H24O2	000143-07-7	80
3		Tridecanoic acid	214	C13H26O2	000638-53-9	80
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5		Undecane	156	C11H24	001120-21-4	11



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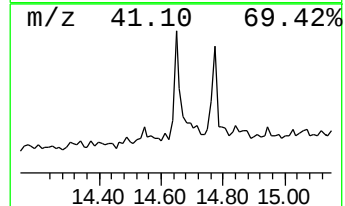
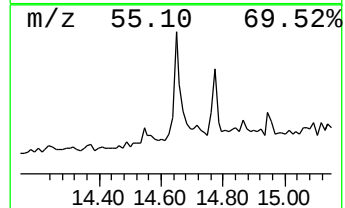
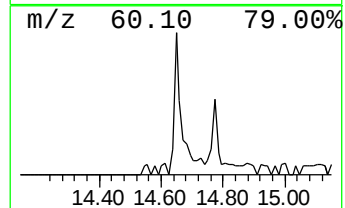
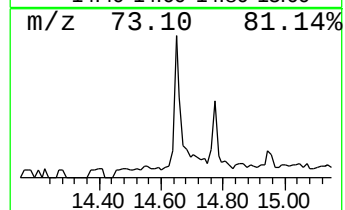
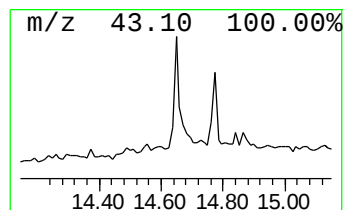
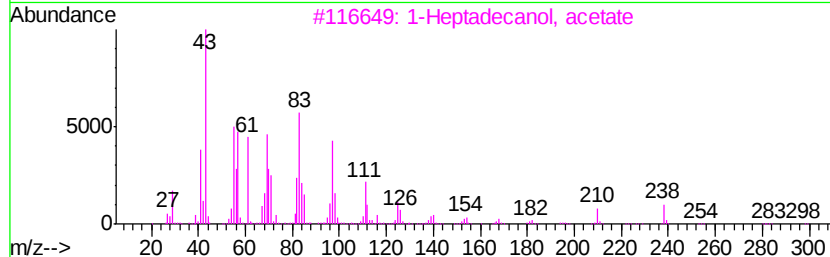
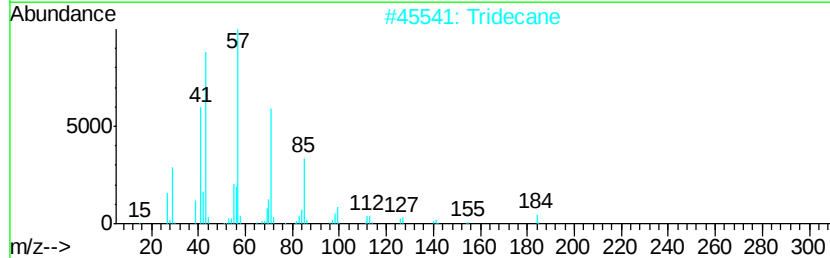
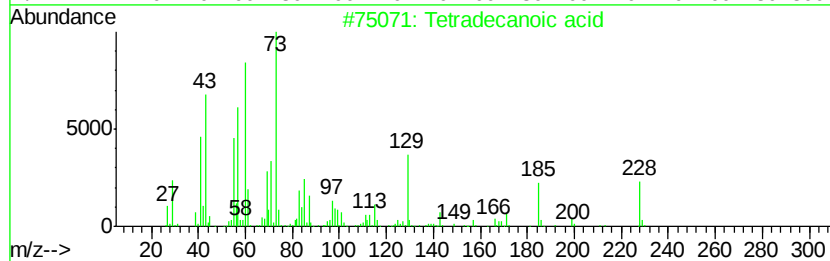
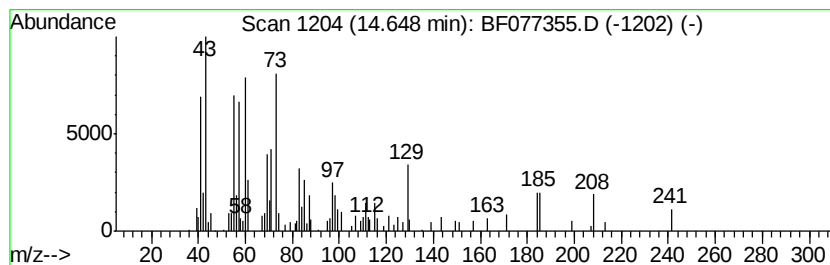
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetradecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.65	3.93 ng	128364	Chrysene-d12	16.27

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanoic acid	228	C14H28O2	000544-63-8	64
2			Tridecane	184	C13H28	000629-50-5	35
3			1-Heptadecanol, acetate	298	C19H38O2	000822-20-8	15
4			Tetracontane, 3,5,24-trimethyl-	605	C43H88	055162-61-3	14
5			Nonahexacontanoic acid	999	C69H138O2	040710-32-5	14



Data Path : S:\HPCHEM1\BNA_F\DATA\BF021815\
Data File : BF077355.D
Acq On : 18 Feb 2015 16:40
Operator : TP/IZ
Sample : G1300-03
Misc : S
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-GF-04-005-B

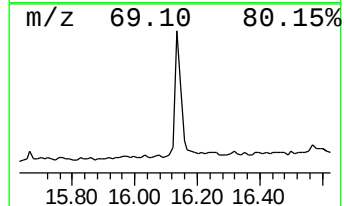
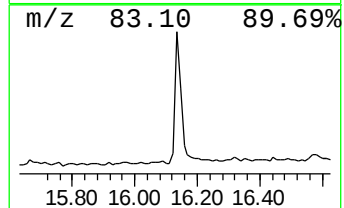
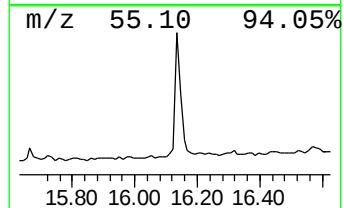
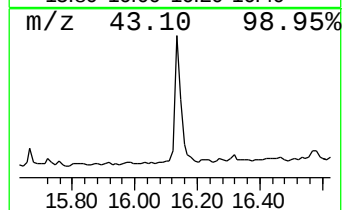
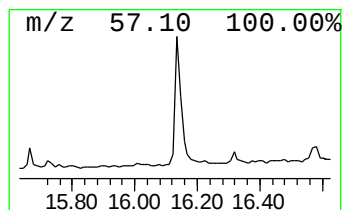
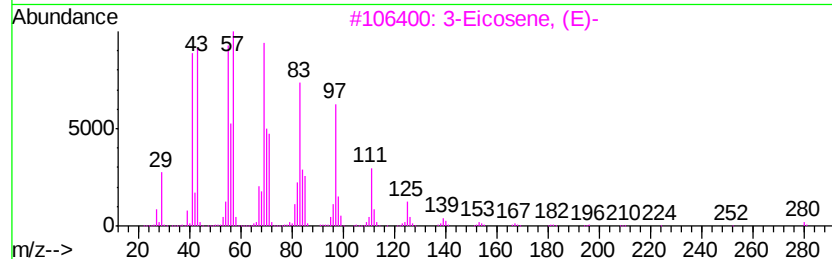
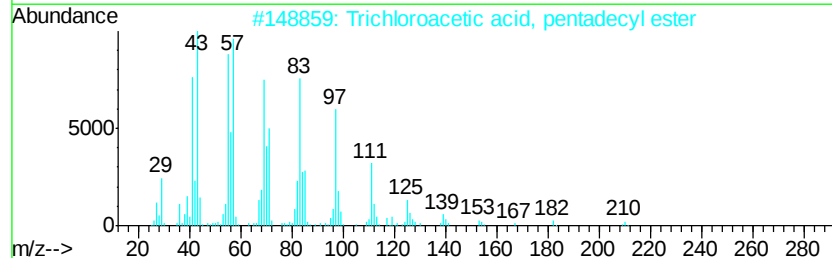
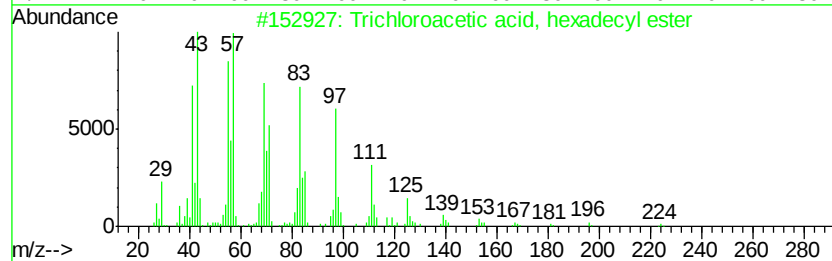
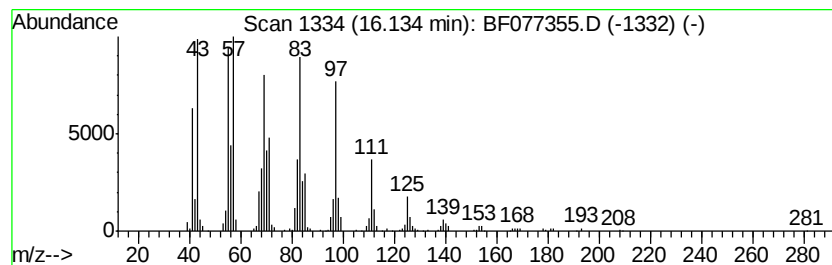
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Trichloroacetic acid, hexad... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.13	11.80 ng	385209	Chrysene-d12	16.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
2		Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
3		3-Eicosene, (E)-	280	C20H40	074685-33-9	91
4		1-Nonadecene	266	C19H38	018435-45-5	91
5		11-Tricosene	322	C23H46	052078-56-5	91



Data Path : S:\HPCHEM1\BNA_F\DATA\BF021815\
Data File : BF077355.D
Acq On : 18 Feb 2015 16:40
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Misc : S
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-GF-04-005-B

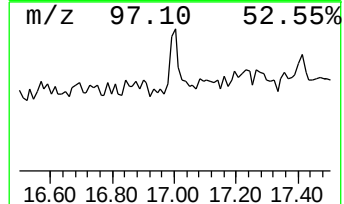
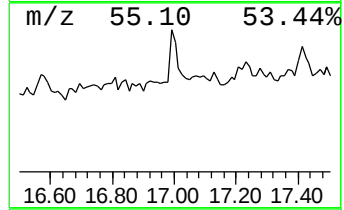
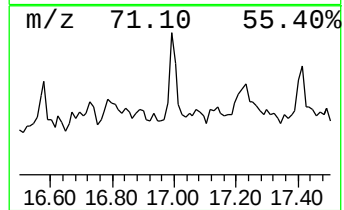
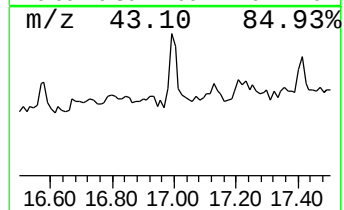
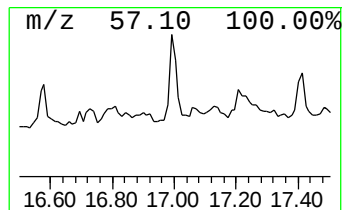
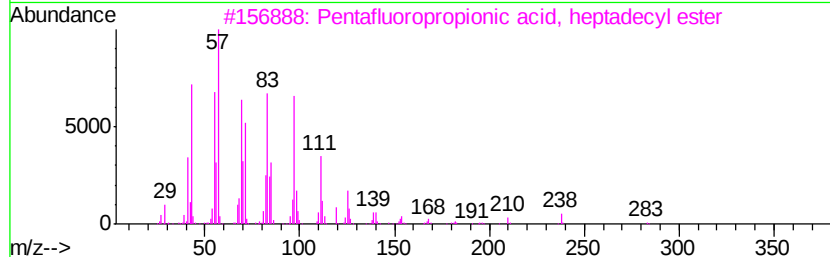
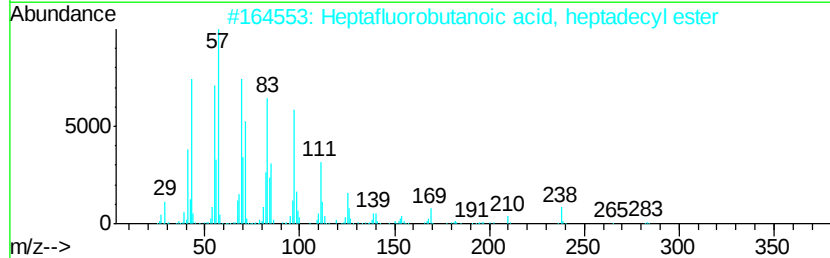
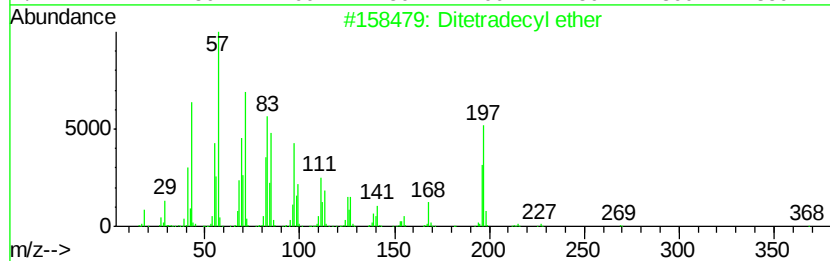
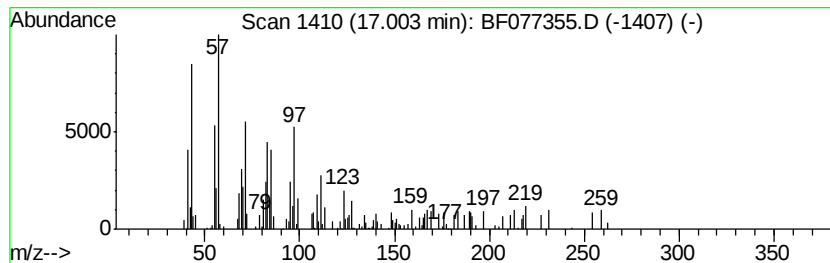
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown17.00 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.00	2.93 ng	95602	Chrysene-d12	16.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ditetradecyl ether	410	C28H58O	005412-98-6	47
2		Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	43
3		Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	43
4		1-Nonadecene	266	C19H38	018435-45-5	38
5		Hexatriacontane	507	C36H74	000630-06-8	35



Data Path : S:\HPCHEM1\BNA_F\DATA\BF021815\
Data File : BF077355.D
Acq On : 18 Feb 2015 16:40
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Sample : G1300-03
Misc : S
ALS Vial : 3 Sample Multiplier: 1

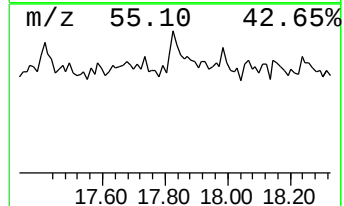
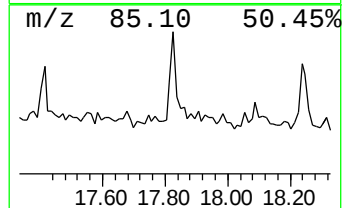
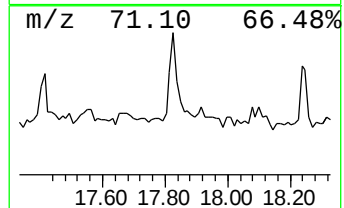
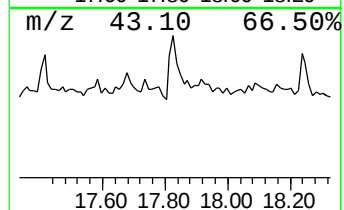
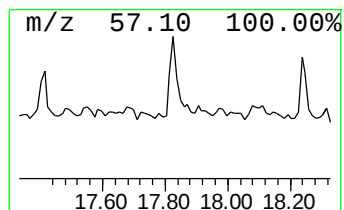
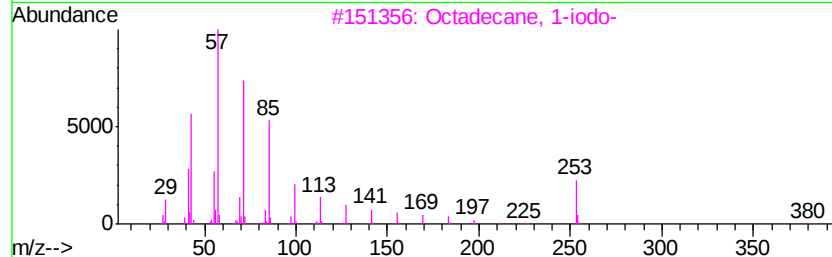
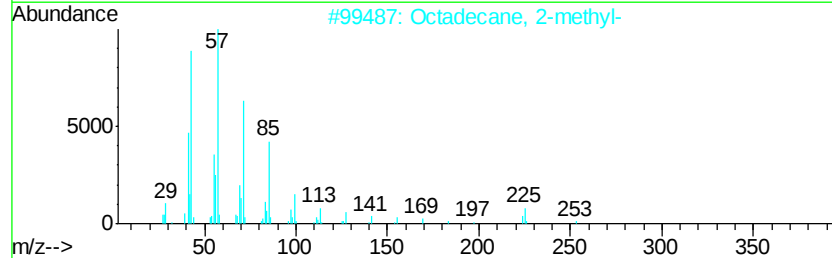
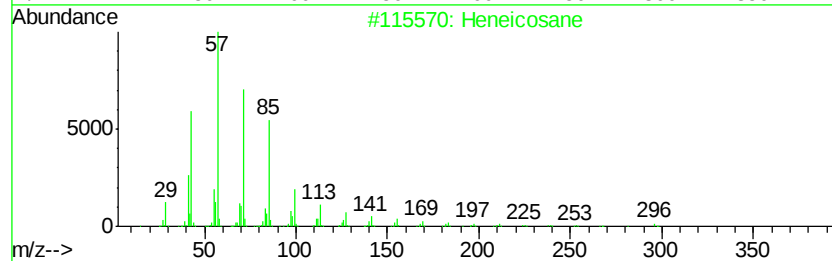
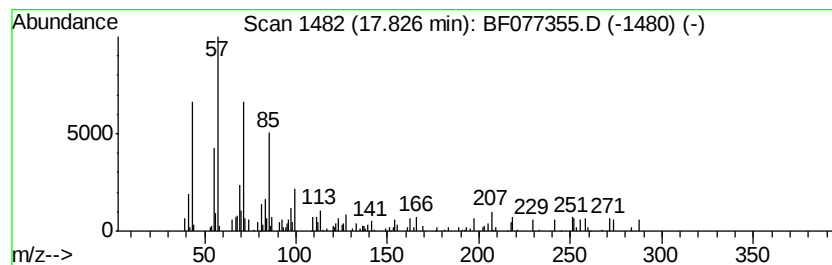
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heneicosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.83	3.80 ng	117981	Perylene-d12	18.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Heneicosane	296 C21H44	000629-94-7	76
2	Octadecane, 2-methyl-	268 C19H40	001560-88-9	64
3	Octadecane, 1-iodo-	380 C18H37I	000629-93-6	64
4	Eicosane, 2-methyl-	296 C21H44	001560-84-5	64
5	Nonadecane, 2-methyl-	282 C20H42	001560-86-7	64



Data Path : S:\HPCHEM1\BNA_F\DATA\BF021815\
Data File : BF077355.D
Acq On : 18 Feb 2015 16:40
Operator : TP/IZ
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Misc : S
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
FK-GF-04-005-B

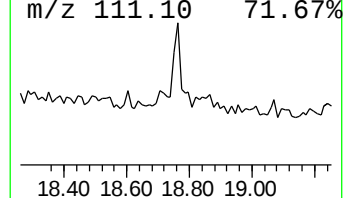
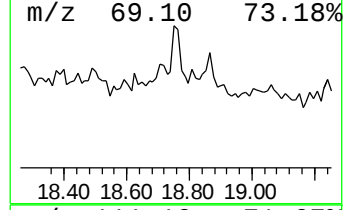
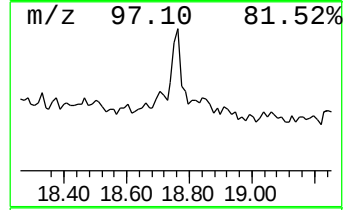
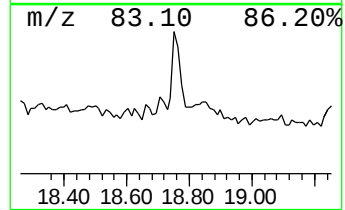
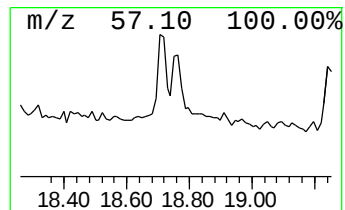
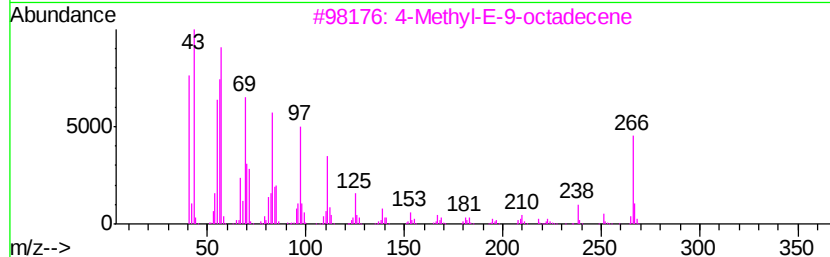
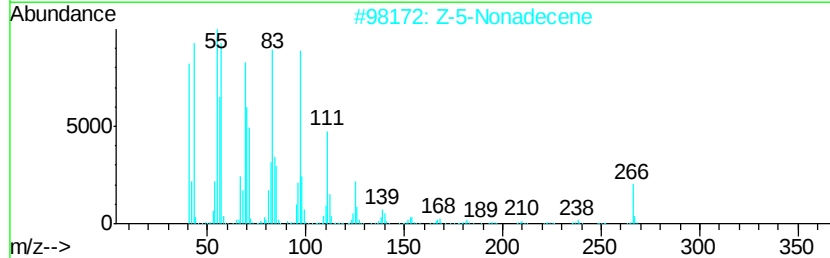
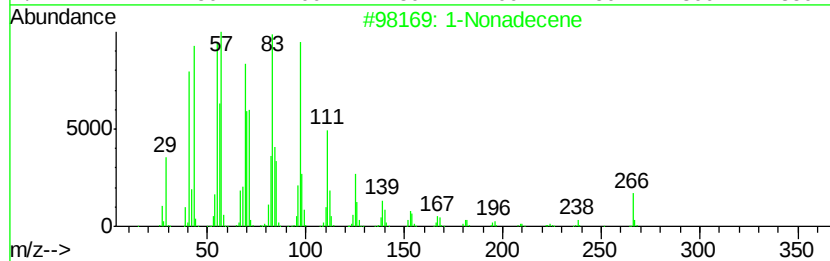
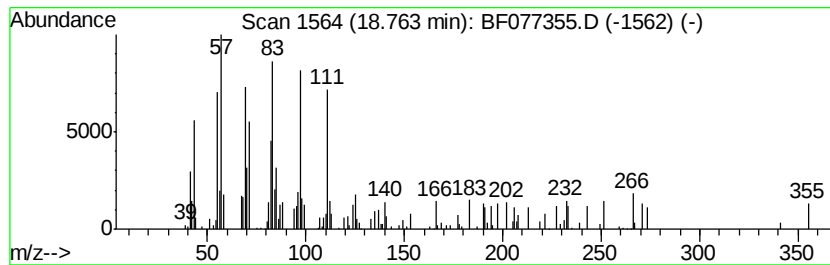
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 1-Nonadecene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.76	2.10 ng	65217	Perylene-d12	18.09

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Nonadecene	266	C19H38	018435-45-5	83
2			Z-5-Nonadecene	266	C19H38	1000131-11-8	83
3			4-Methyl-E-9-octadecene	266	C19H38	1000130-87-1	60
4			2-Methyl-E-7-octadecene	266	C19H38	1000130-92-3	49
5			2-Cyclohexen-1-ol, 1-methyl-	112	C7H12O	023758-27-2	46



Data Path : S:\HPCHEM1\BNA_F\DATA\BF021815\
Data File : BF077355.D
Acq On : 18 Feb 2015 16:40
Operator : TP/IZ
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Misc : S
ALS Vial : 3 Sample Multiplier: 1

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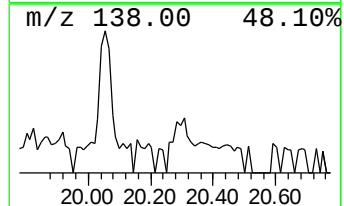
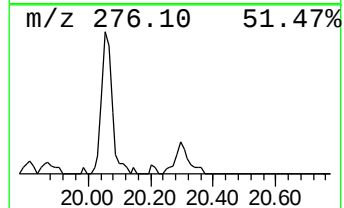
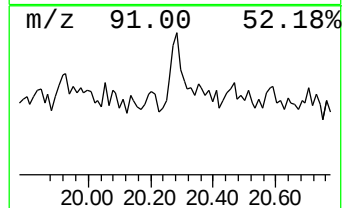
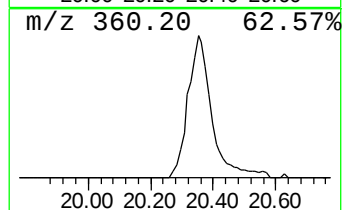
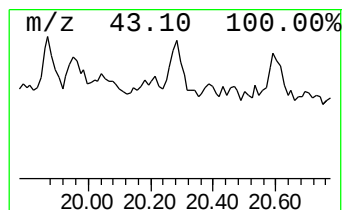
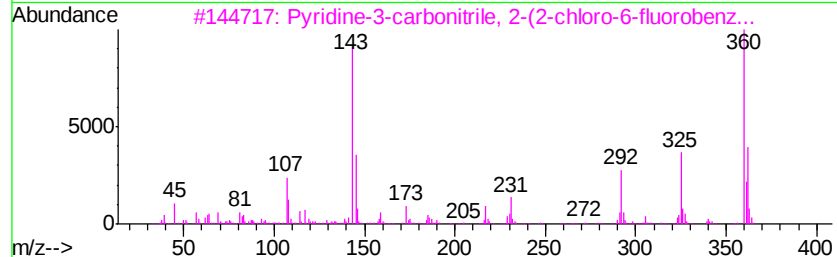
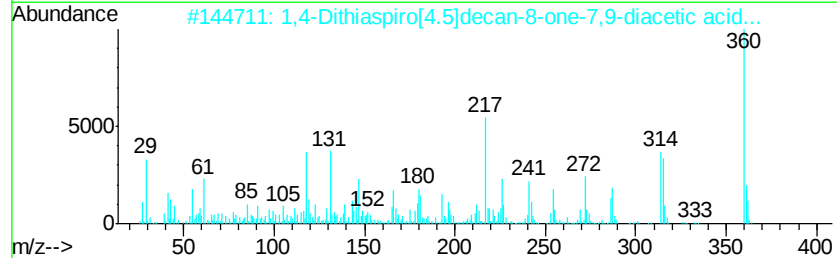
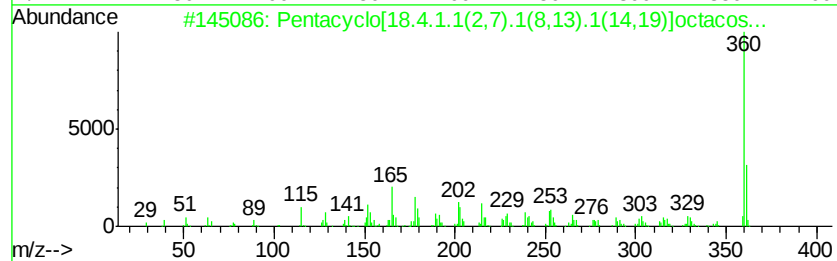
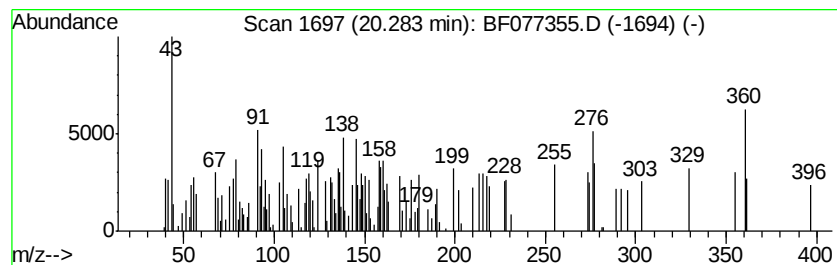
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown20.28 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.28	2.62 ng	81465	Perylene-d12	18.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentacyclo[18.4.1.1(2,7).1(8,13)...	360	C28H24	101077-60-5	14
2		1,4-Dithiaspiro[4.5]decan-8-one-...	360	C16H24O5S2	1000195-58-5	11
3		Pyridine-3-carbonitrile, 2-(2-ch...	360	C17H10ClFN2S2	298687-66-8	10
4		4-Chloro-4'-[2,4-diamino-5-pyrim...	360	C16H13ClN4O2S	052258-30-7	10
5		1H-Pyrrole-3-carboxylic acid, 4-...	360	C19H24N2O5	1000268-77-3	10



Data Path : S:\HPCHEM1\BNA_F\DATA\BF021815\
Data File : BF077355.D
Acq On : 18 Feb 2015 16:40
Operator : TP/IZ
Sample : G1300-03
Misc : S
ALS Vial : 3 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF021715.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
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Butane, 2-methoxy...	1.84	13.5	ng	187048	1	7.36	277872	20.0
2-Pentanone, 4-hy...	5.15	29.7	ng	412891	1	7.36	277872	20.0
unknown7.06	7.06	95.9	ng	1331850	1	7.36	277872	20.0
n-Hexadecanoic acid	13.67	4.4	ng	117501	4	12.96	528804	20.0
Tetradecanoic acid	14.65	3.9	ng	128364	5	16.27	653038	20.0
Trichloroacetic a...	16.13	11.8	ng	385209	5	16.27	653038	20.0
unknown17.00	17.00	2.9	ng	95602	5	16.27	653038	20.0
Heneicosane	17.83	3.8	ng	117981	6	18.09	621589	20.0
1-Nonadecene	18.76	2.1	ng	65217	6	18.09	621589	20.0
unknown20.28	20.28	2.6	ng	81465	6	18.09	621589	20.0