

Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
 Data File : BF079411.D
 Acq On : 21 May 2015 19:07
 Operator : TP/IZ
 Sample : G2325-03
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW15

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: BF079411.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.335	10	13	24	rVB	281977	704172	7.15%	2.255%
2	1.484	24	26	34	rVV	620923	1034142	10.50%	3.312%
3	1.610	36	37	46	rVV	477470	878397	8.92%	2.814%
4	1.724	46	47	87	rVB	4514641	9847848	100.00%	31.543%
5	5.347	362	364	372	rBV	640981	697872	7.09%	2.235%
6	6.639	475	477	482	rBV	305537	431938	4.39%	1.384%
7	6.776	487	489	492	rBV	1319539	1431071	14.53%	4.584%
8	7.073	513	515	517	rBV	313125	349800	3.55%	1.120%
9	7.256	528	531	533	rBV	1651693	1505916	15.29%	4.824%
10	7.770	573	576	578	rBV	1044999	1177212	11.95%	3.771%
11	8.650	650	653	655	rBV	432002	478072	4.85%	1.531%
12	9.999	768	771	773	rBV	2237466	2424347	24.62%	7.765%
13	10.811	840	842	845	rBV	417202	535577	5.44%	1.715%
14	11.794	925	928	931	rBV	1496968	1534499	15.58%	4.915%
15	11.999	944	946	948	rBV	109362	120178	1.22%	0.385%
16	12.057	949	951	953	rVV	323909	328742	3.34%	1.053%
17	12.102	953	955	957	rVV2	391302	510638	5.19%	1.636%
18	12.148	957	959	963	rVV2	234486	525091	5.33%	1.682%
19	12.217	963	965	968	rVV2	138850	254701	2.59%	0.816%
20	12.274	968	970	973	rVV2	244681	324605	3.30%	1.040%
21	12.331	973	975	977	rVV	302548	407512	4.14%	1.305%
22	12.377	977	979	981	rVB	180988	195686	1.99%	0.627%
23	12.651	1001	1003	1006	rBV	596623	575967	5.85%	1.845%
24	14.651	1175	1178	1180	rBV	2442284	2587481	26.27%	8.288%
25	15.943	1289	1291	1294	rBV	586242	623343	6.33%	1.997%
26	16.948	1377	1379	1384	rBV3	113264	223799	2.27%	0.717%
27	17.360	1412	1415	1420	rBV3	106600	208421	2.12%	0.668%
28	17.726	1444	1447	1455	rBV	526443	967099	9.82%	3.098%
29	18.160	1482	1485	1490	rBV2	102695	178221	1.81%	0.571%
30	18.606	1521	1524	1530	rBV2	80761	157898	1.60%	0.506%

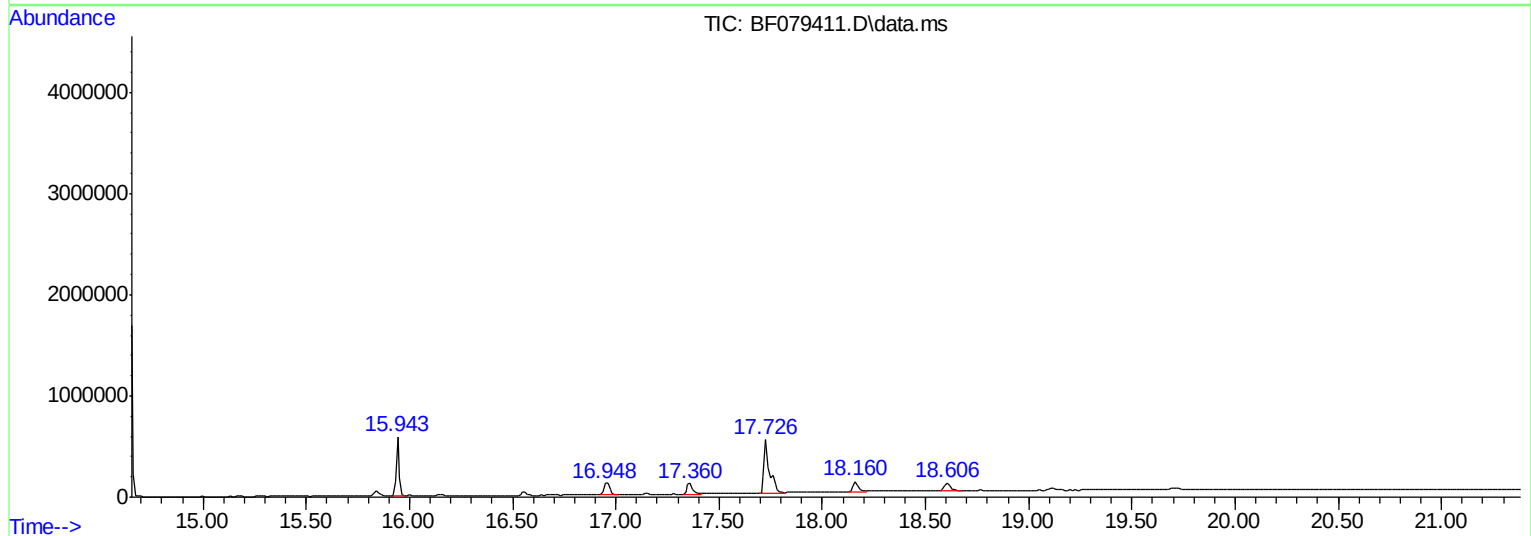
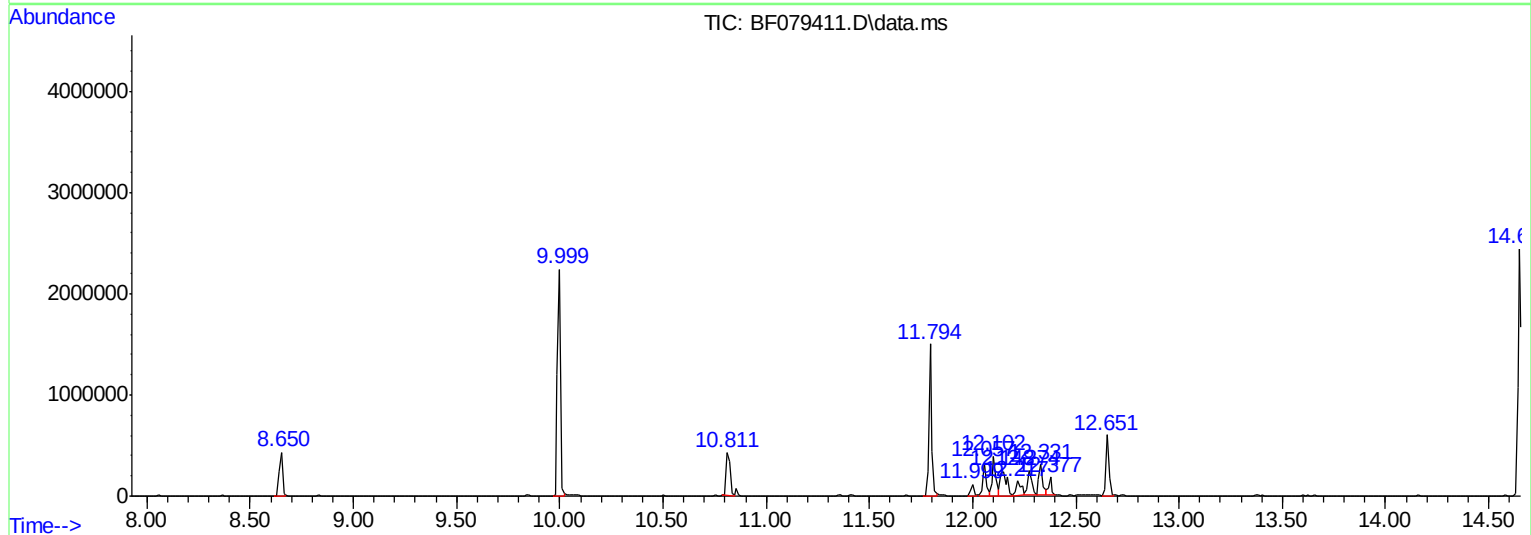
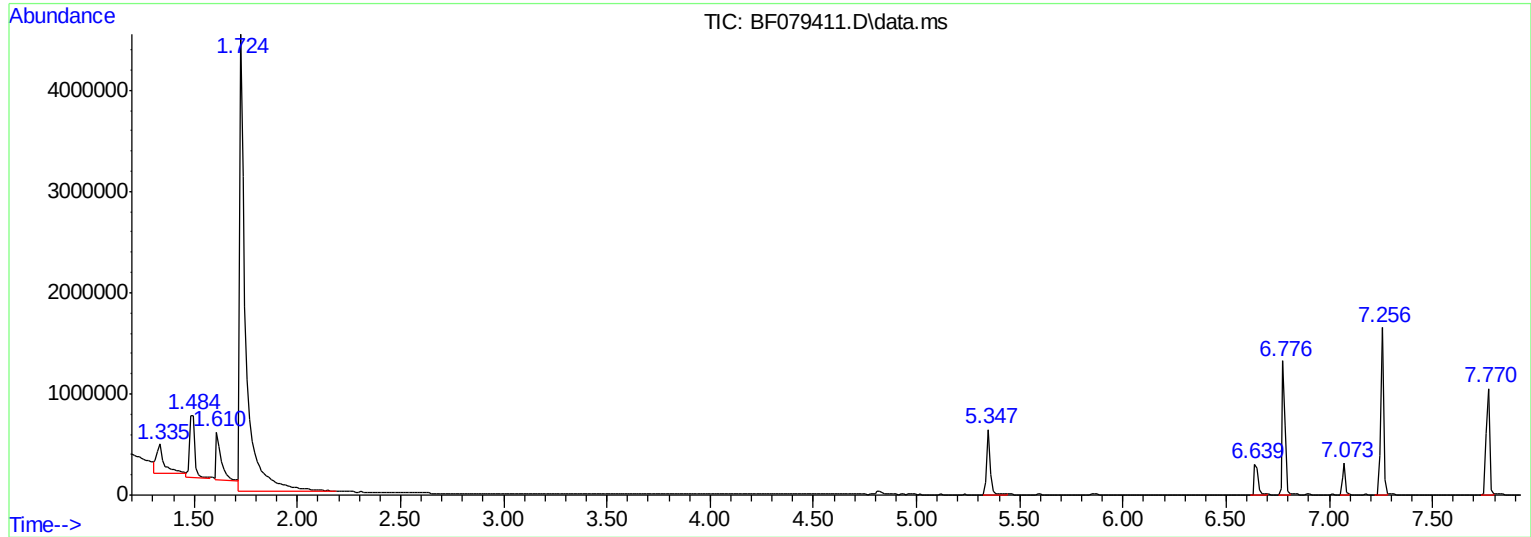
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Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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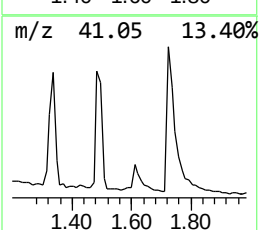
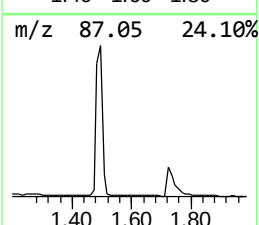
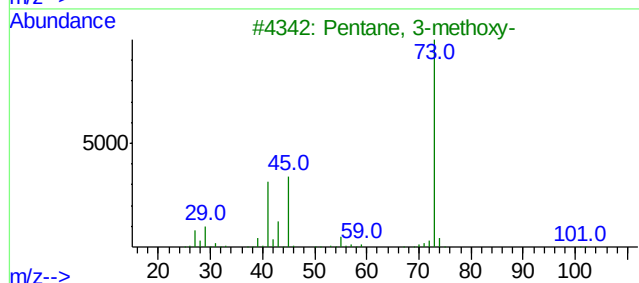
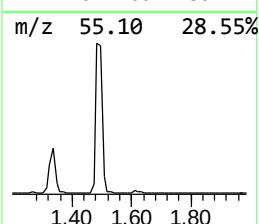
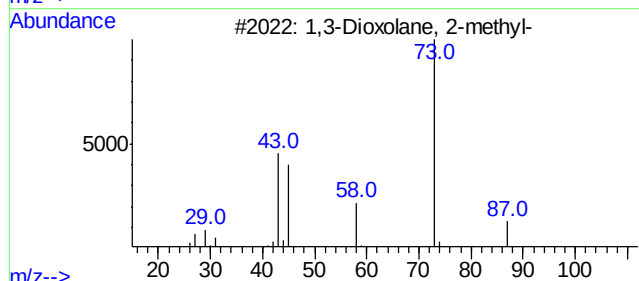
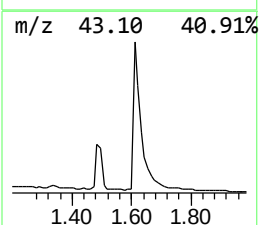
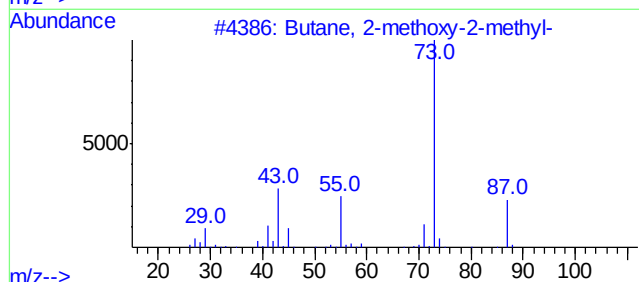
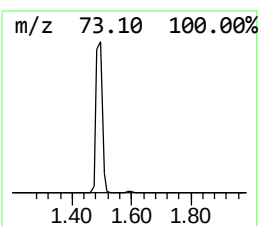
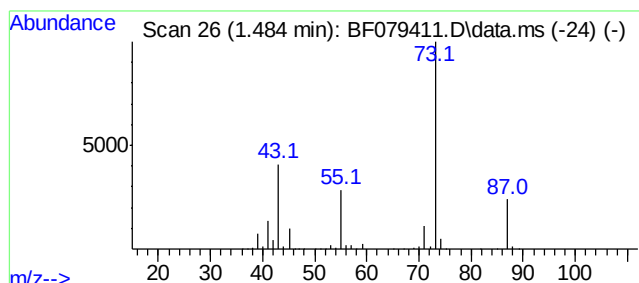
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Butane, 2-methoxy-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.484	59.13 ng	1034140	1,4-Dichlorobenzene-d4	7.073

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
4	Silane, tetramethyl-	88	C4H12Si	000075-76-3	10
5	Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9



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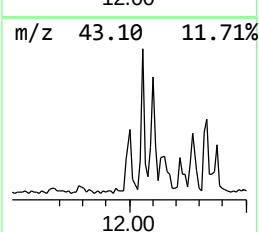
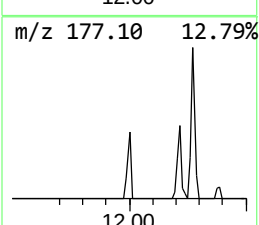
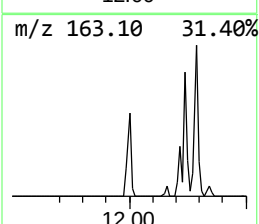
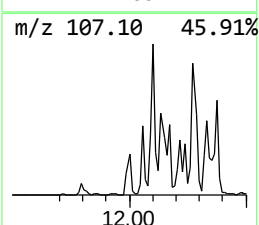
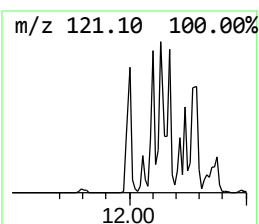
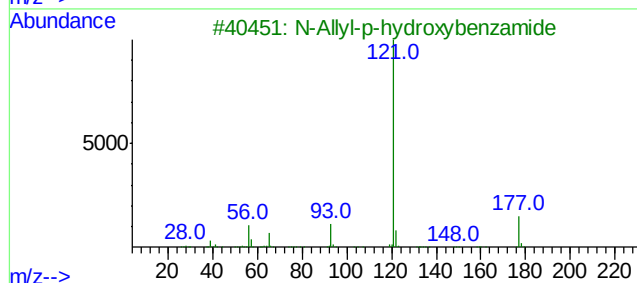
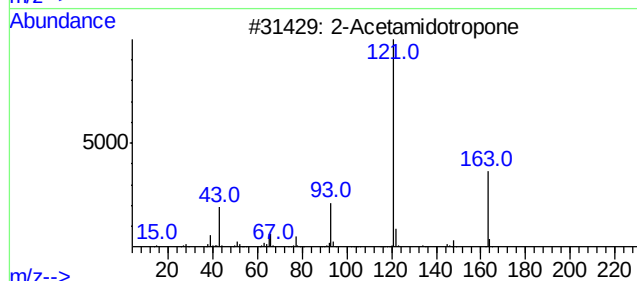
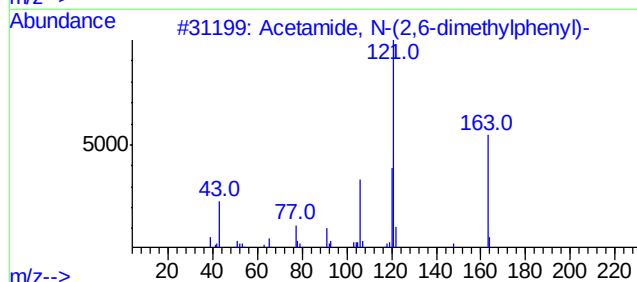
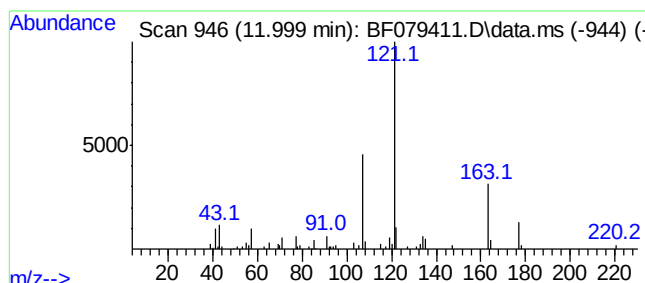
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Peak Number 7 Acetamide, N-(2,6-dimethylp... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.999	4.17 ng	120178	Phenanthrene-d10	12.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(2,6-dimethylphenyl)-	163	C10H13NO	002198-53-0	50
2	2-Acetamidotropone	163	C9H9NO2	006422-12-4	47
3	N-Allyl-p-hydroxybenzamide	177	C10H11NO2	090921-72-5	43
4	2-Propanone, 1-(4-methoxyphenyl)-	164	C10H12O2	000122-84-9	43
5	p-Sec-butylphenyl acetate	192	C12H16O2	003245-24-7	43



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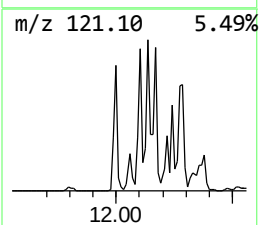
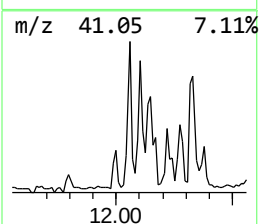
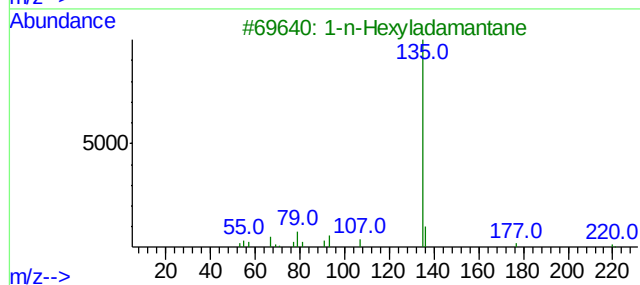
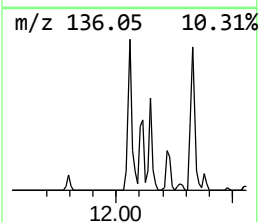
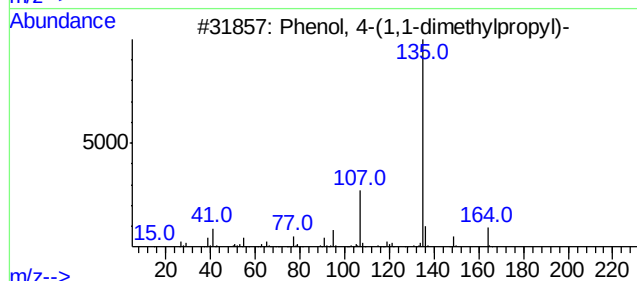
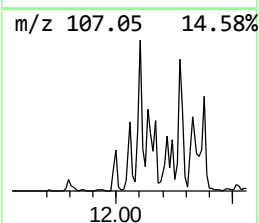
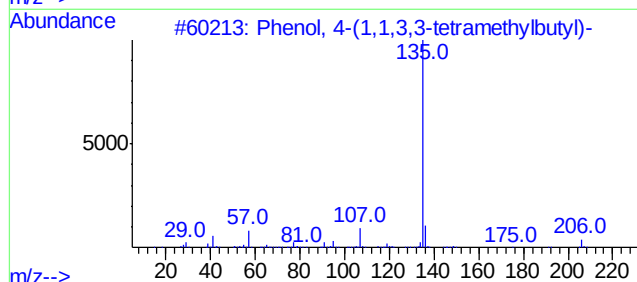
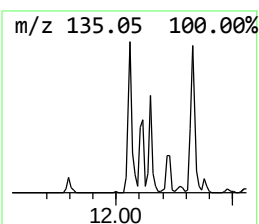
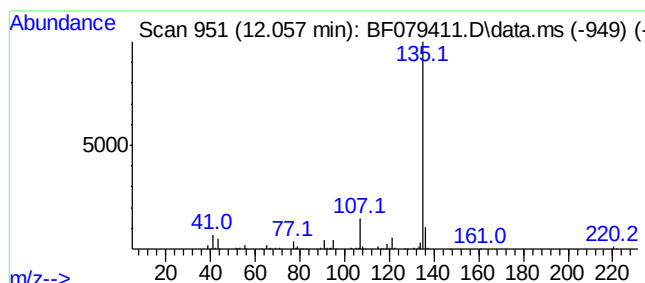
Instrument :
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.057	11.42 ng	328742	Phenanthrene-d10	12.651	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
2	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
3	1-n-Hexyladamantane	220	C16H28	022458-75-9	59
4	1-Adamantanecarboxylic acid, 2-a...	314	C21H30O2	1000282-94-3	50
5	m-Anisic acid, cyclobutyl ester	206	C12H14O3	1000292-25-0	50



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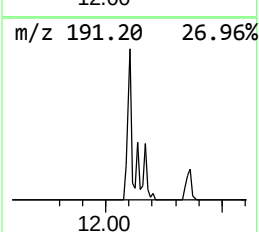
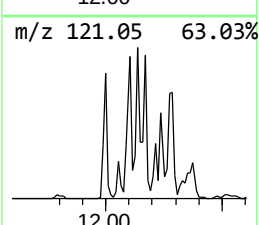
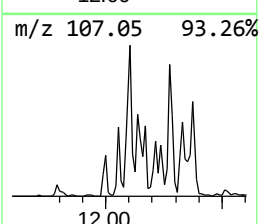
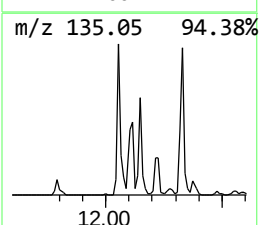
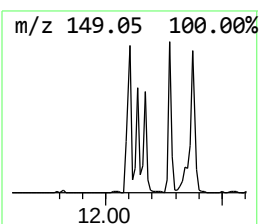
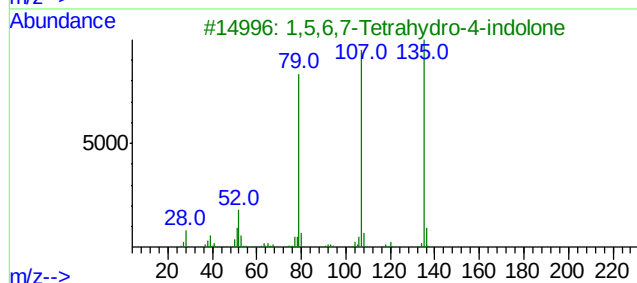
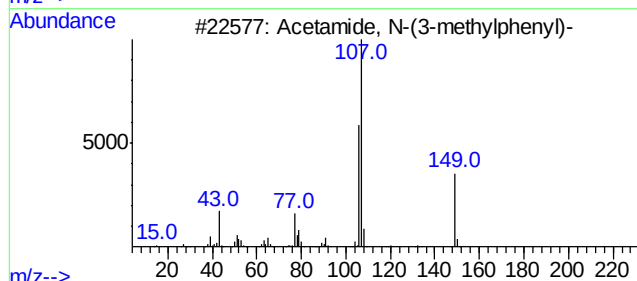
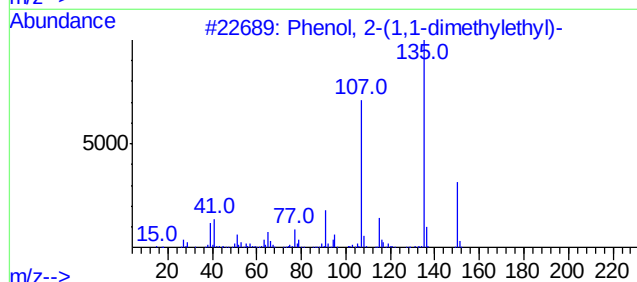
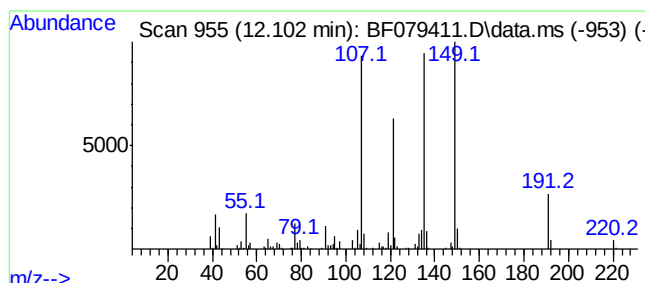
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 9 Phenol, 2-(1,1-dimethylethyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.102	17.73 ng	510638	Phenanthrene-d10	12.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	55
2	Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	49
3	1,5,6,7-Tetrahydro-4-indolone	135	C8H9NO	013754-86-4	46
4	2,5-Cyclohexadiene, 1,4-diethyl-...	164	C12H20	1000150-21-6	43
5	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	35



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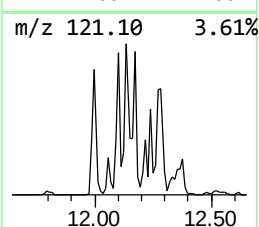
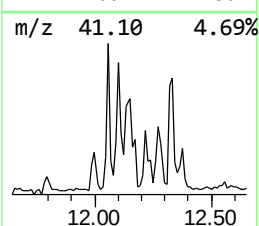
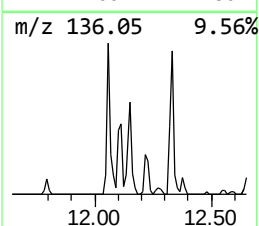
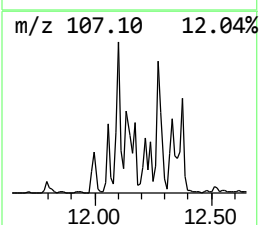
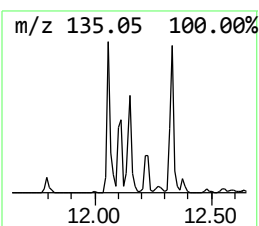
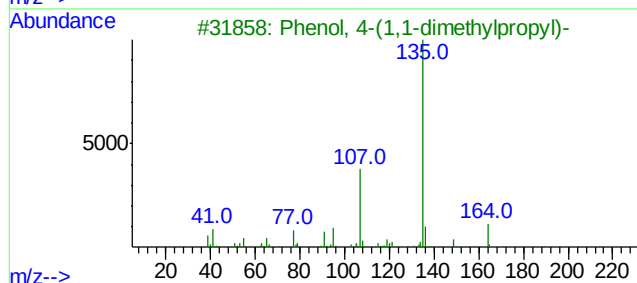
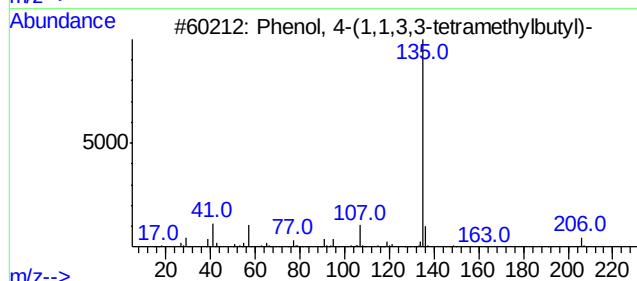
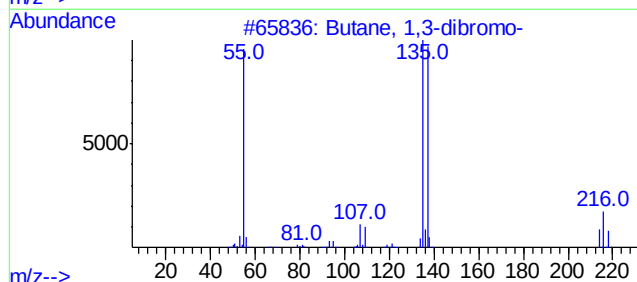
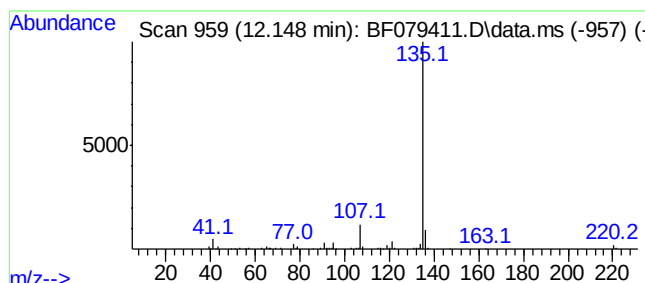
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Peak Number 10 Butane, 1,3-dibromo- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.148	18.23 ng	525091	Phenanthrene-d10	12.651	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 1,3-dibromo-	214	C4H8Br2	000107-80-2	78
2	Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
3	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
4	2H-1,4-Benzoxazine, 3,4-dihydro-	135	C8H9NO	005735-53-5	56
5	Carbamic acid, N-[1,1-bis(triflu...	413	C19H25F6NO2	296242-69-8	56



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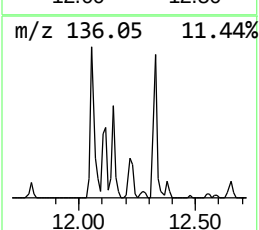
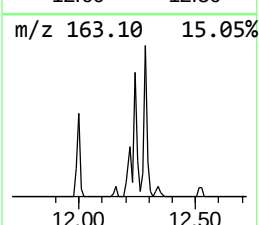
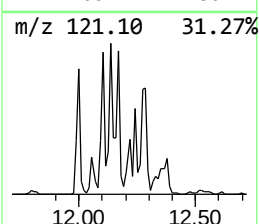
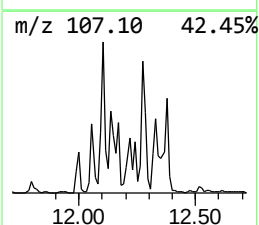
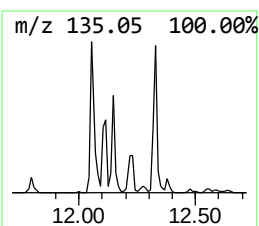
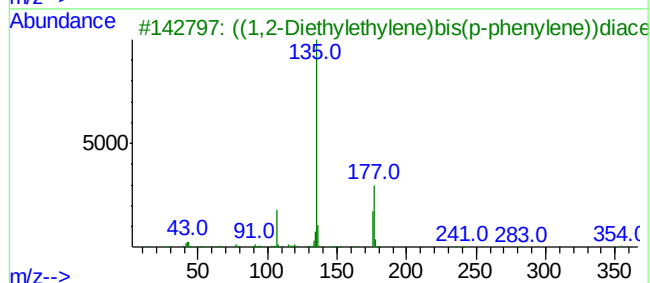
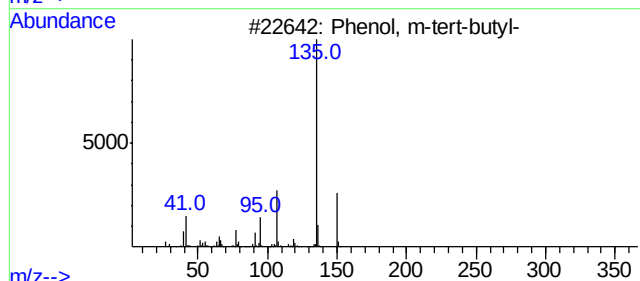
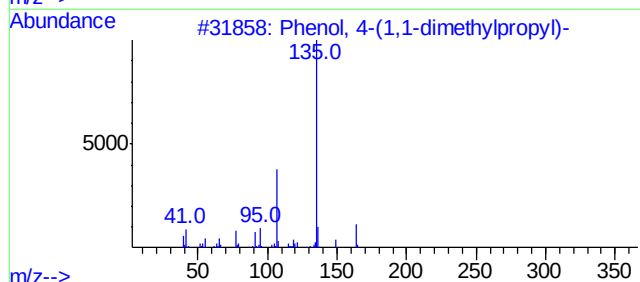
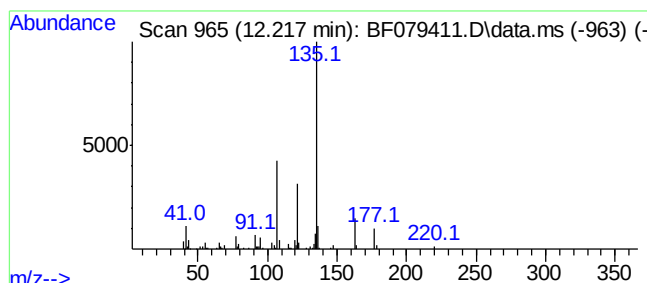
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Phenol, 4-(1,1-dimethylprop... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.217	8.84 ng	254701	Phenanthrene-d10	12.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	72
2	Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	64
3	((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	64
4	Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	50
5	S-[1,2,3-Benzotriazin-4-yl]-N-me...	220	C9H8N4OS	1000254-70-9	47



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079411.D
Acq On : 21 May 2015 19:07
Operator : TP/IZ
Sample : G2325-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

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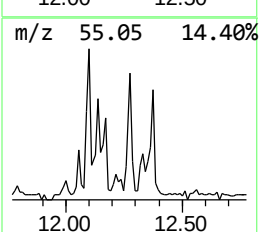
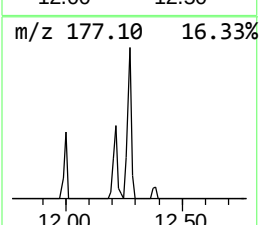
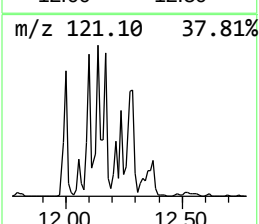
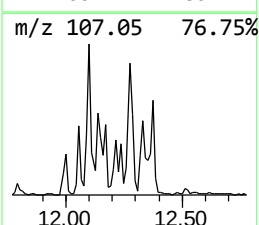
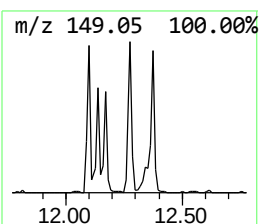
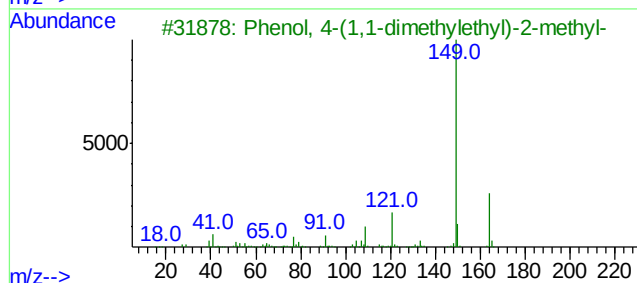
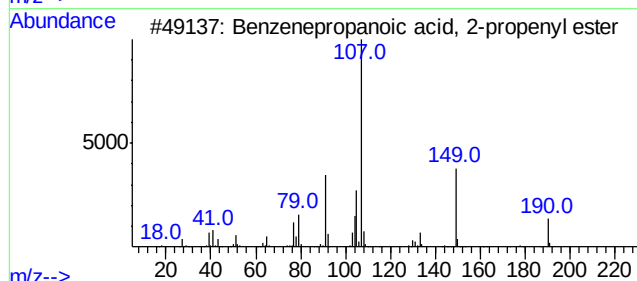
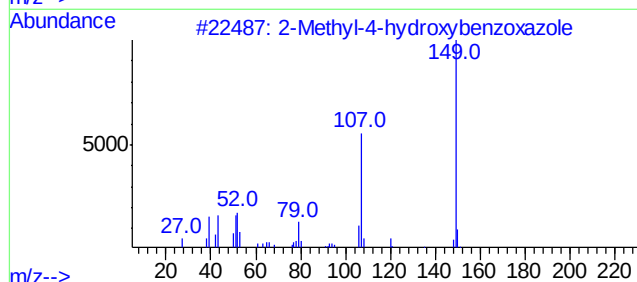
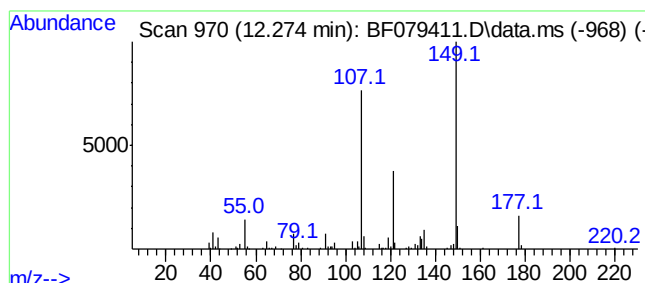
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2-Methyl-4-hydroxybenzoxazole Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.274	11.27 ng	324605	Phenanthrene-d10	12.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-4-hydroxybenzoxazole	149	C8H7NO2	051110-60-2	64
2	Benzenepropanoic acid, 2-propenyl...	190	C12H14O2	015814-45-6	50
3	Phenol, 4-(1,1-dimethylethyl)-2-...	164	C11H16O	000098-27-1	38
4	Phenol, 4-butyl-	150	C10H14O	001638-22-8	38
5	4-Methyl-2-tert-octylphenol	220	C15H24O	004979-46-8	35



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079411.D
Acq On : 21 May 2015 19:07
Operator : TP/IZ
Sample : G2325-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW15

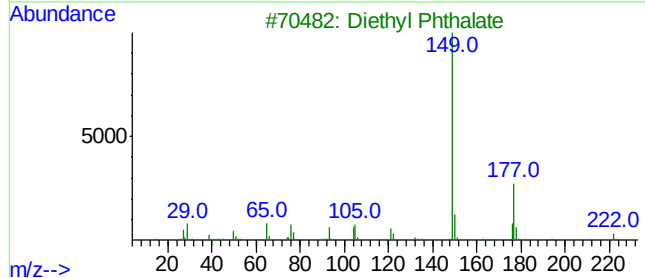
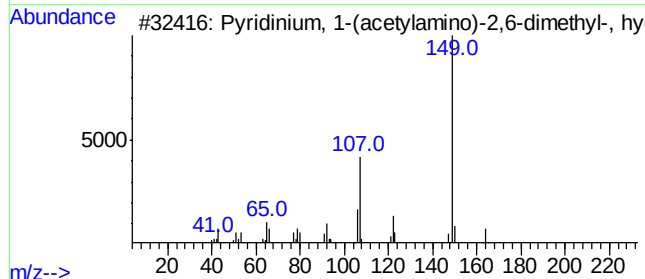
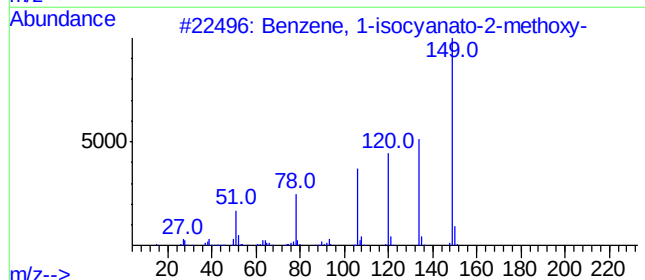
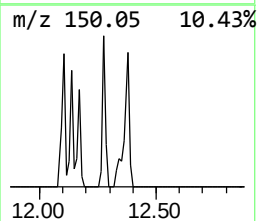
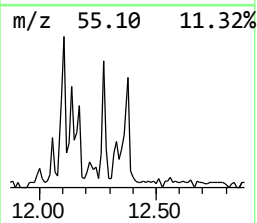
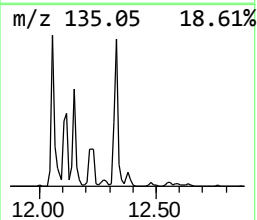
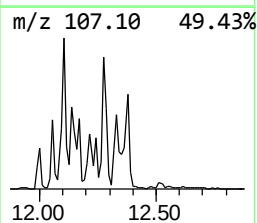
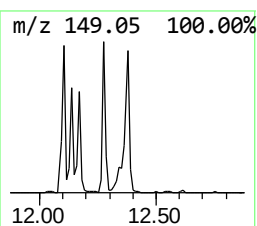
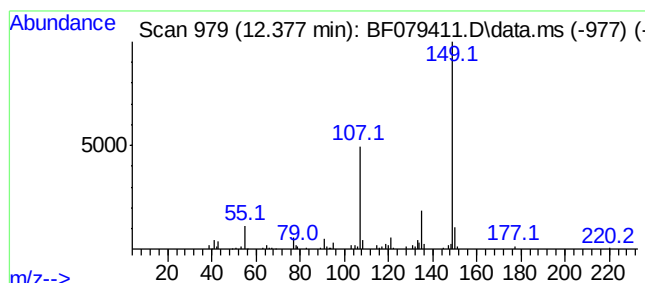
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Benzene, 1-isocyanato-2-met... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.377	6.80 ng	195686	Phenanthrene-d10	12.651

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-isocyanato-2-methoxy-	149	C8H7NO2	000700-87-8	50
2	Pyridinium, 1-(acetylamino)-2,6-...	164	C9H12N2O	031020-35-6	40
3	Diethyl Phthalate	222	C12H14O4	000084-66-2	38
4	Phenol, 4-(1,1-dimethylethyl)-2-...	164	C11H16O	000098-27-1	38
5	Phenol, 2-methyl-4-(1,1,3,3-tetr...	220	C15H24O	002219-84-3	38



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079411.D
Acq On : 21 May 2015 19:07
Operator : TP/IZ
Sample : G2325-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

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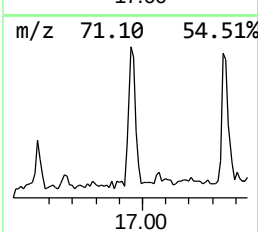
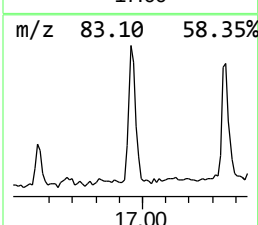
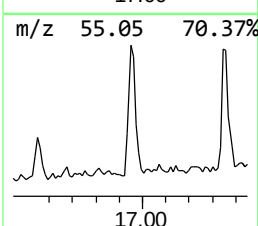
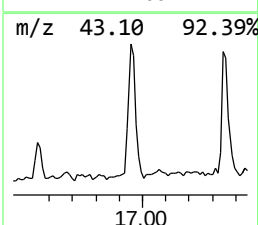
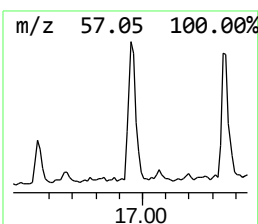
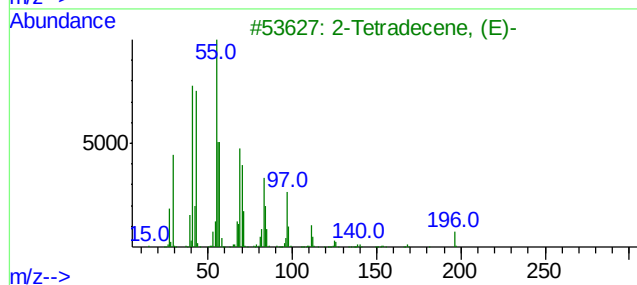
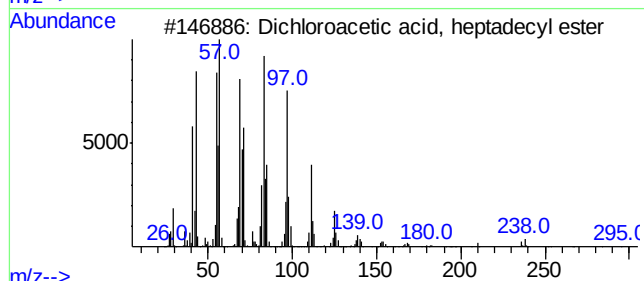
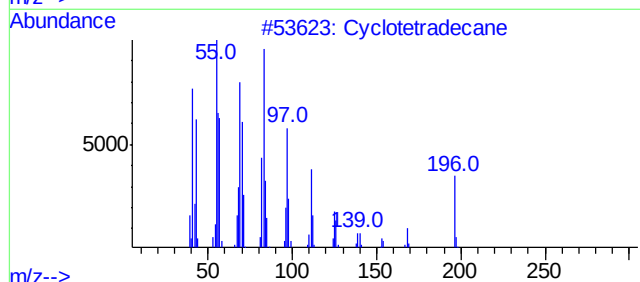
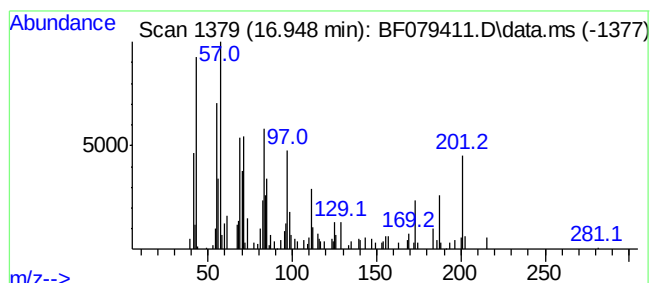
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Cyclotetradecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.948	4.63 ng	223799	Perylene-d12	17.726

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetradecane	196	C14H28	000295-17-0	94
2	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	90
3	2-Tetradecene, (E)-	196	C14H28	035953-53-8	89
4	1-Hexadecanol, 2-methyl-	256	C17H36O	002490-48-4	70
5	Trifluoroacetic acid, n-octadecy...	366	C20H37F3O2	079392-43-1	64



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
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Acq On : 21 May 2015 19:07
Operator : TP/IZ
Sample : G2325-03
Misc :
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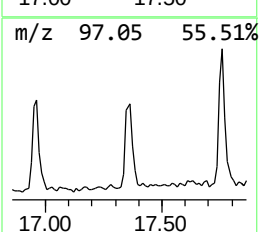
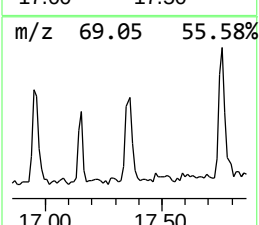
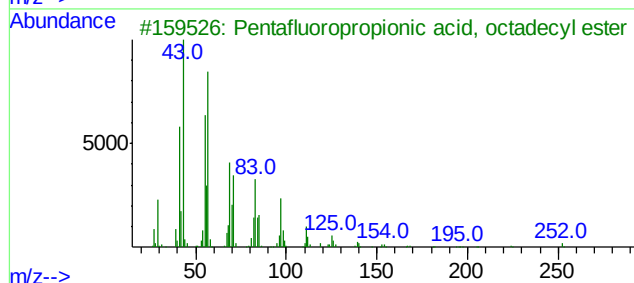
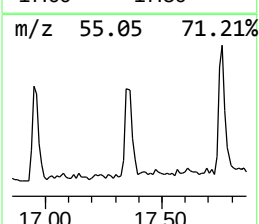
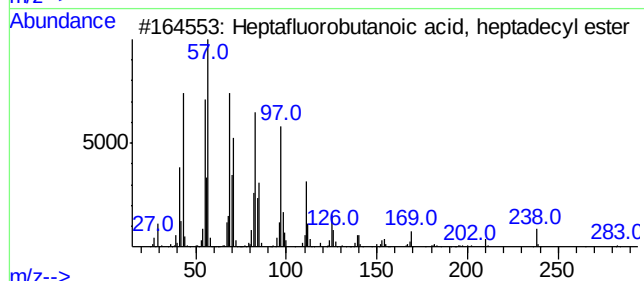
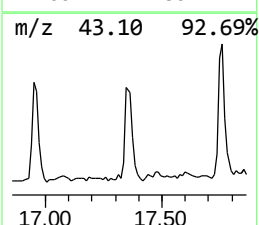
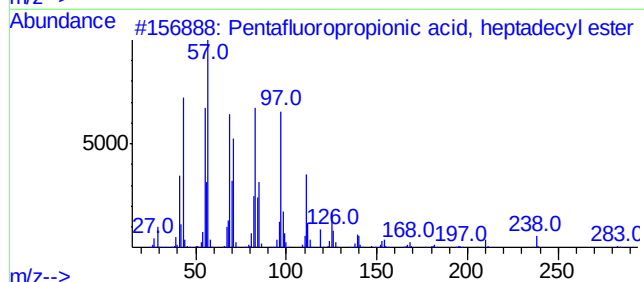
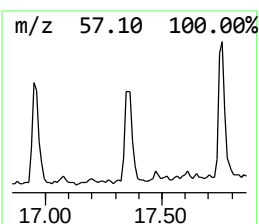
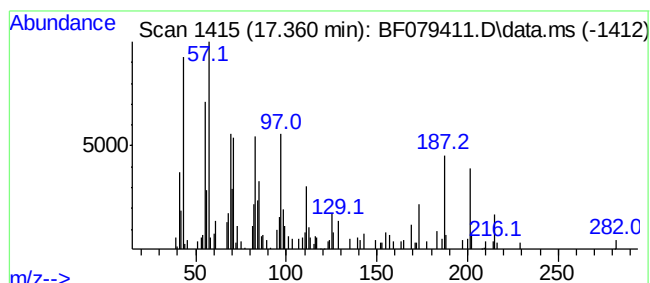
Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 Pentafluoropropionic acid, ... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.360	4.31 ng	208421	Perylene-d12	17.726	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	91
2	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	90
3	Pentafluoropropionic acid, octad...	416	C21H37F5O2	1000280-07-7	64
4	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	60
5	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	60



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079411.D
Acq On : 21 May 2015 19:07
Operator : TP/IZ
Sample : G2325-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW15

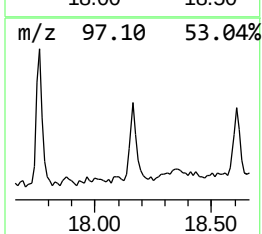
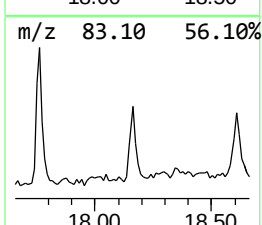
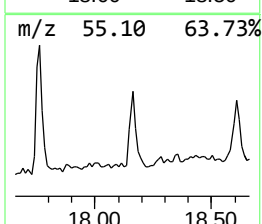
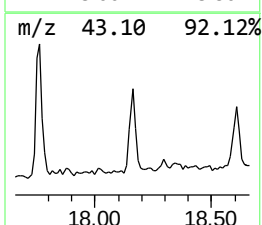
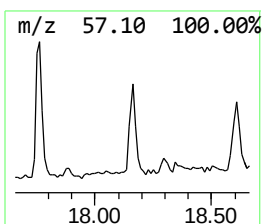
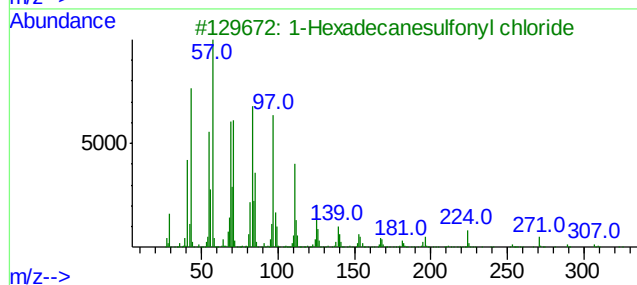
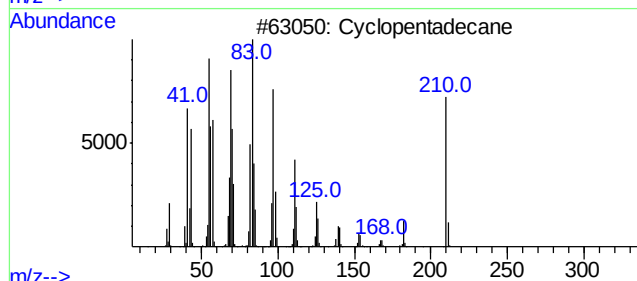
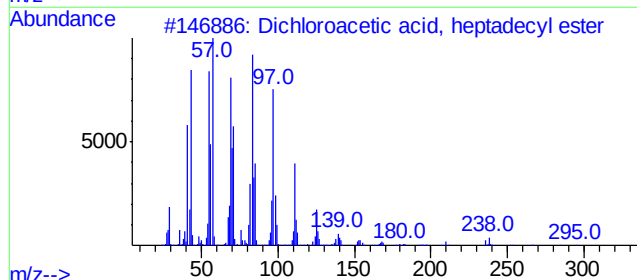
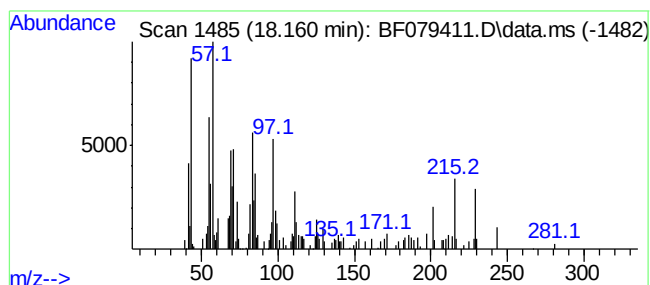
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Dichloroacetic acid, heptad... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.160	3.69 ng	178221	Perylene-d12	17.726

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	92
2	Cyclopentadecane	210	C15H30	000295-48-7	89
3	1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	70
4	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	70
5	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	64



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
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Acq On : 21 May 2015 19:07
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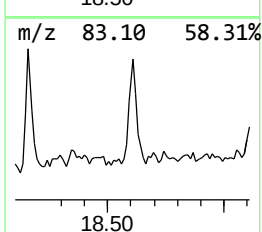
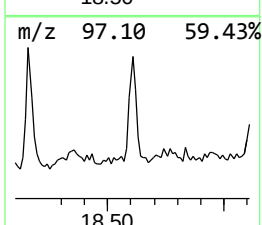
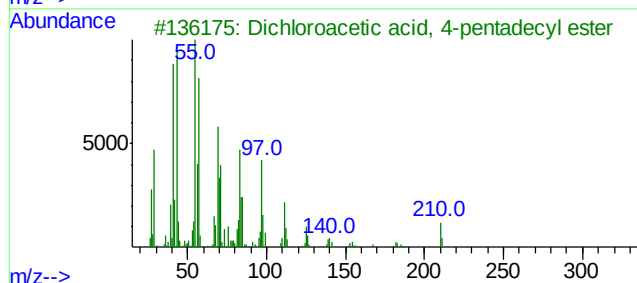
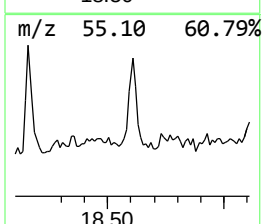
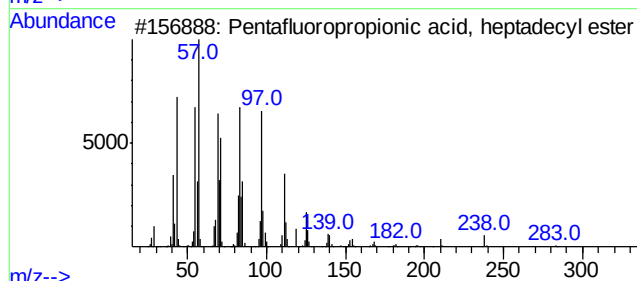
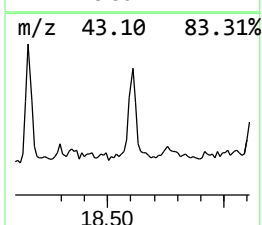
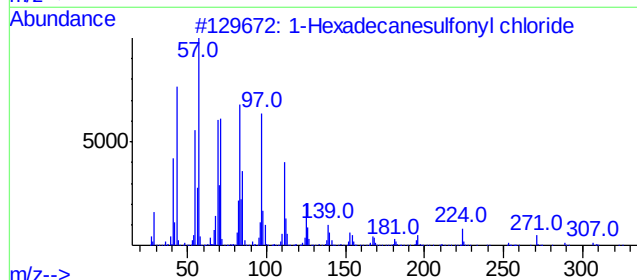
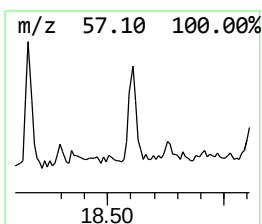
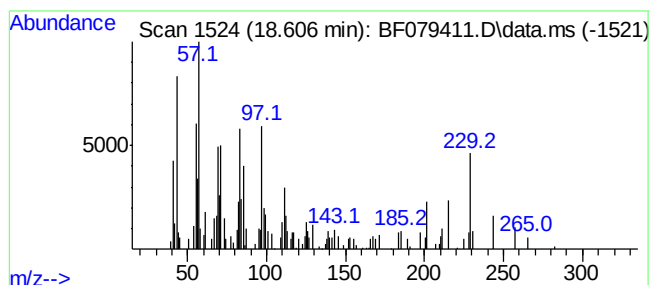
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 1-Hexadecanesulfonyl chloride Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.606	3.27 ng	157898	Perylene-d12	17.726	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexadecanesulfonyl chloride	324	C16H33ClO2S	038775-38-1	62
2	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	1000283-04-2	60
3	Dichloroacetic acid, 4-pentadecy...	338	C17H32Cl2O2	1000280-64-9	58
4	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	53
5	Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	53



Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079411.D
Acq On : 21 May 2015 19:07
Operator : TP/IZ
Sample : G2325-03
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW15

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
Butane, 2-metho...	1.484	59.1	ng	1034140	1	7.073	349800	20.0
Acetamide, N-(2...	11.999	4.2	ng	120178	4	12.651	575967	20.0
Phenol, 4-(1,1,...	12.057	11.4	ng	328742	4	12.651	575967	20.0
Phenol, 2-(1,1-...	12.102	17.7	ng	510638	4	12.651	575967	20.0
Butane, 1,3-dib...	12.148	18.2	ng	525091	4	12.651	575967	20.0
Phenol, 4-(1,1-...	12.217	8.8	ng	254701	4	12.651	575967	20.0
2-Methyl-4-hydr...	12.274	11.3	ng	324605	4	12.651	575967	20.0
Benzene, 1-isoc...	12.377	6.8	ng	195686	4	12.651	575967	20.0
Cyclotetradecane	16.948	4.6	ng	223799	6	17.726	967099	20.0
Pentafluoroprop...	17.360	4.3	ng	208421	6	17.726	967099	20.0
Dichloroacetic ...	18.160	3.7	ng	178221	6	17.726	967099	20.0
1-Hexadecanesul...	18.606	3.3	ng	157898	6	17.726	967099	20.0