

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
 Data File : BF095321.D
 Acq On : 22 May 2017 18:58
 Operator : SJ/MA
 Sample : I3232-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-PS-01-003

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-BF050217.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.119	538	541	561	rBV	130612	345623	13.00%	1.510%
2	5.719	640	643	675	rBV	975583	2235715	84.06%	9.765%
3	7.301	907	912	928	rBV	1316861	2242670	84.32%	9.795%
4	7.419	928	932	950	rVB	1608185	2474690	93.05%	10.809%
5	7.760	986	990	1007	rVB	304473	389566	14.65%	1.701%
6	7.995	1025	1030	1044	rBV	1310634	1641669	61.73%	7.170%
7	8.660	1138	1143	1161	rBV	798420	1360590	51.16%	5.943%
8	9.778	1329	1333	1352	rBV	316132	620844	23.34%	2.712%
9	11.548	1628	1634	1651	rBV	1753964	2659595	100.00%	11.616%
10	12.248	1749	1753	1771	rBV	242796	394806	14.84%	1.724%
11	12.613	1810	1815	1839	rBV2	450464	747402	28.10%	3.264%
12	13.442	1952	1956	1961	rVB	31477	36383	1.37%	0.159%
13	13.907	2030	2035	2054	rBV	887125	1568036	58.96%	6.849%
14	14.213	2083	2087	2090	rBV	29748	35474	1.33%	0.155%
15	15.013	2218	2223	2243	rVB	377441	642116	24.14%	2.805%
16	15.965	2381	2385	2396	rBV3	52066	101298	3.81%	0.442%
17	17.330	2612	2617	2636	rBV	1502401	1811808	68.12%	7.913%
18	18.571	2824	2828	2832	rBV2	76502	102285	3.85%	0.447%
19	18.683	2843	2847	2861	rBV	290113	447771	16.84%	1.956%
20	19.359	2958	2962	2964	rBV	46340	58981	2.22%	0.258%
21	19.877	3046	3050	3055	rBV	245401	263973	9.93%	1.153%
22	20.083	3081	3085	3090	rBV	61645	70964	2.67%	0.310%
23	20.318	3121	3125	3128	rVB5	20803	34471	1.30%	0.151%
24	20.359	3128	3132	3142	rBV	296490	463583	17.43%	2.025%
25	20.600	3168	3173	3177	rBV2	149346	194799	7.32%	0.851%
26	20.695	3186	3189	3197	rVB2	21114	33468	1.26%	0.146%
27	20.806	3205	3208	3213	rVB2	33958	42405	1.59%	0.185%
28	21.024	3241	3245	3251	rBV3	28929	52015	1.96%	0.227%
29	21.483	3318	3323	3332	rVB6	51678	93777	3.53%	0.410%
30	21.736	3361	3366	3370	rBV7	20715	40768	1.53%	0.178%
31	22.147	3428	3436	3445	rBV9	90223	309670	11.64%	1.353%
32	22.241	3448	3452	3454	rBV4	23867	37369	1.41%	0.163%
33	22.383	3471	3476	3483	rBV9	38201	92472	3.48%	0.404%
34	22.571	3503	3508	3519	rVB8	69088	217639	8.18%	0.951%

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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 >
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35	22.988	3574	3579	3590	rVB5	67381	162552	6.11%	0.710%
36	23.688	3688	3698	3714	rBV7	125164	524216	19.71%	2.290%
37	23.835	3716	3723	3736	rVB6	105440	344085	12.94%	1.503%

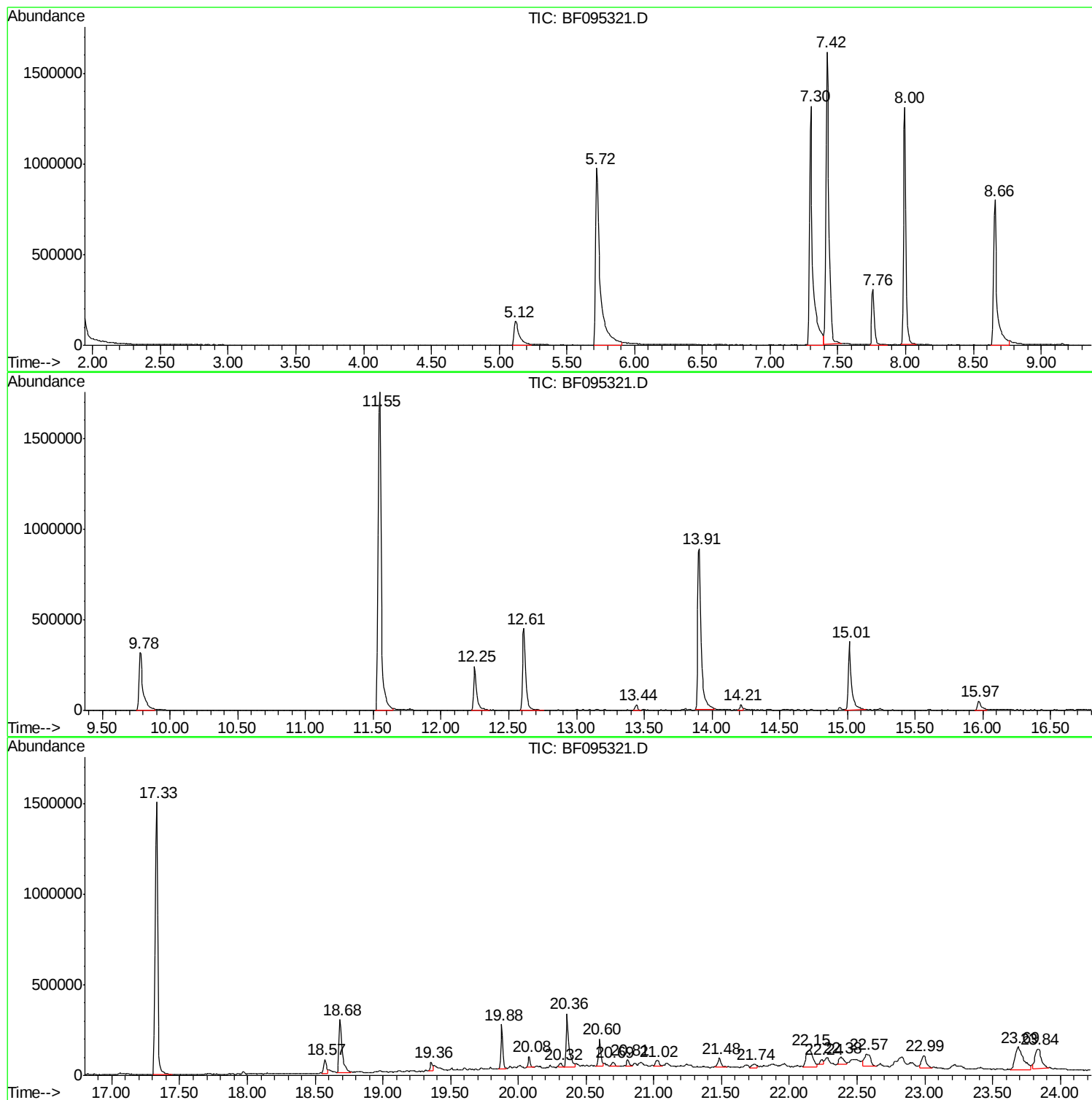
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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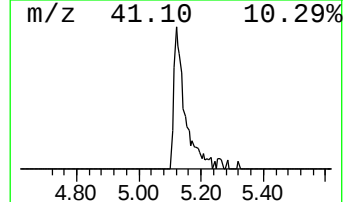
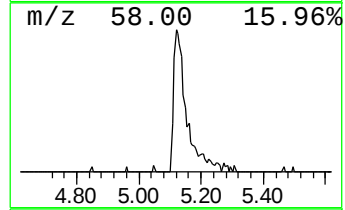
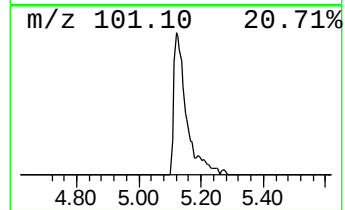
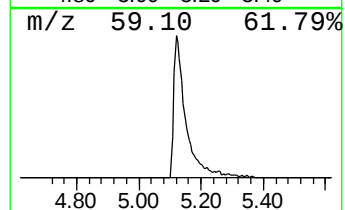
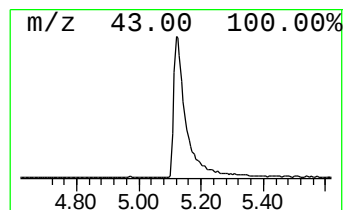
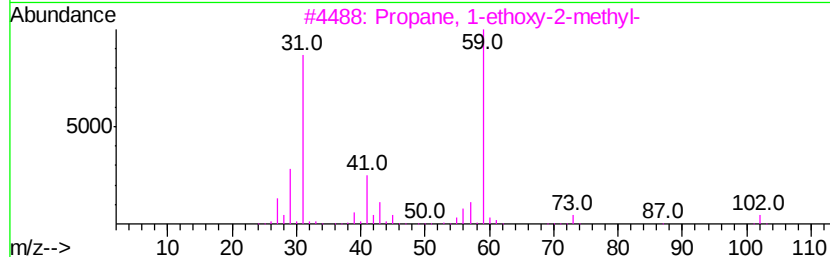
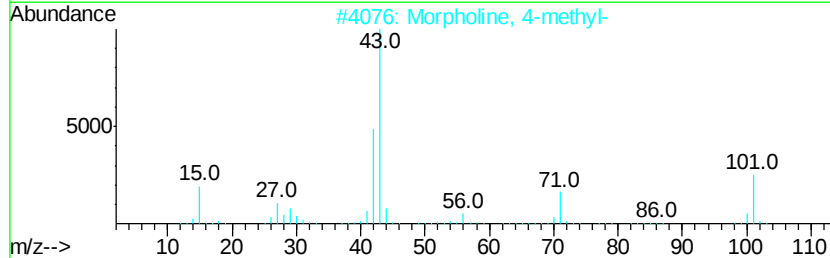
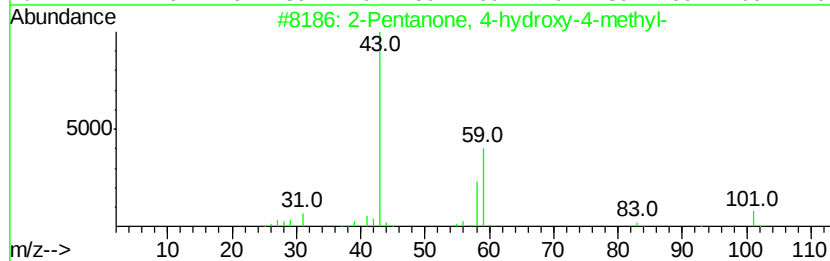
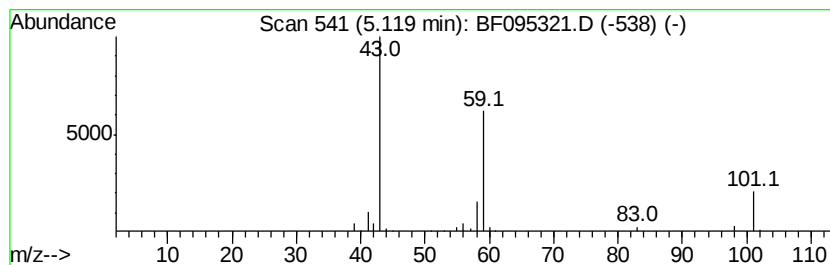
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	17.74 ng	345623	1,4-Dichlorobenzene-d4	7.76

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
3			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9
4			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	9
5			Propanoic acid, 2-nitro-, methyl...	133	C4H7NO4	006118-50-9	9



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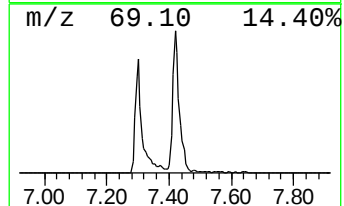
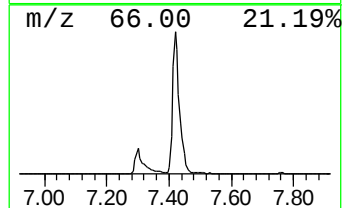
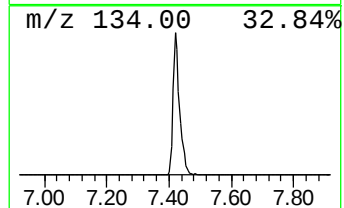
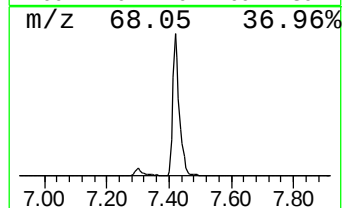
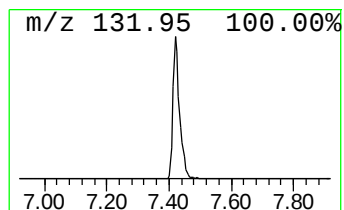
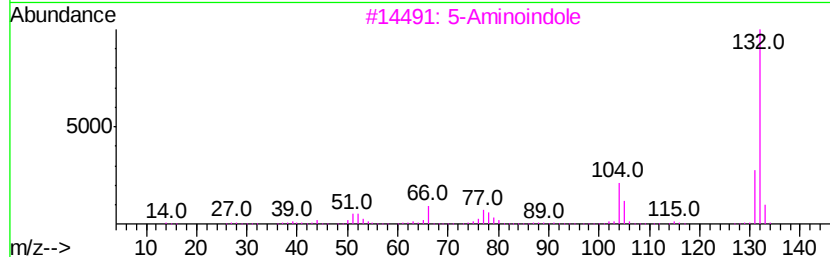
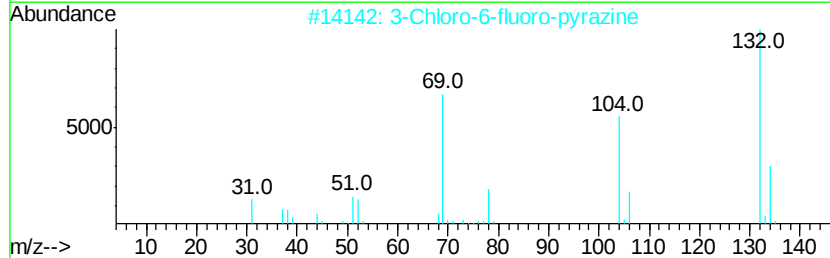
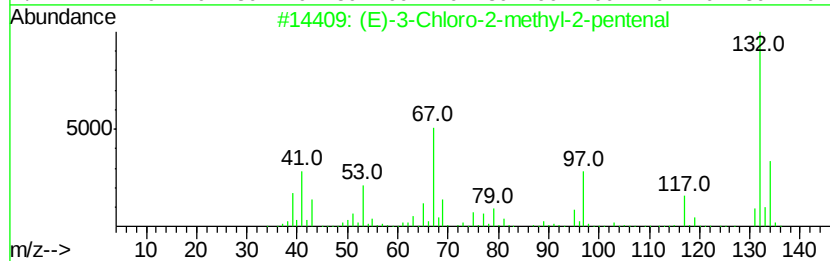
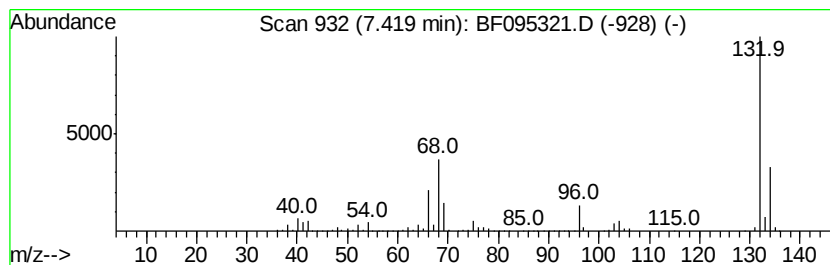
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown7.42 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.42	127.05 ng	2474690	1,4-Dichlorobenzene-d4	7.76

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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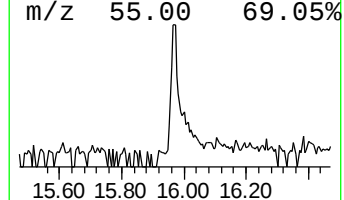
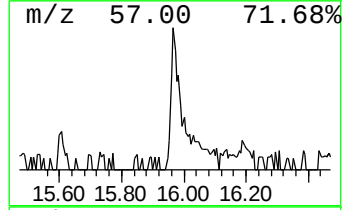
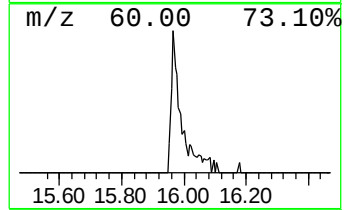
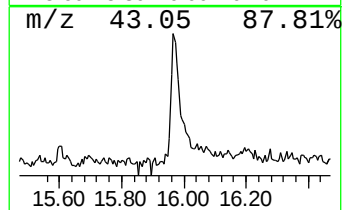
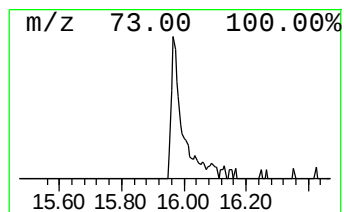
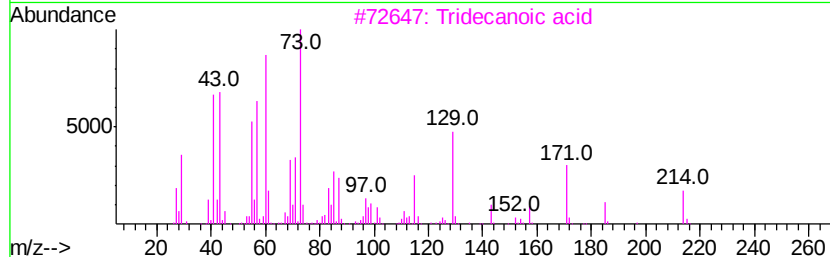
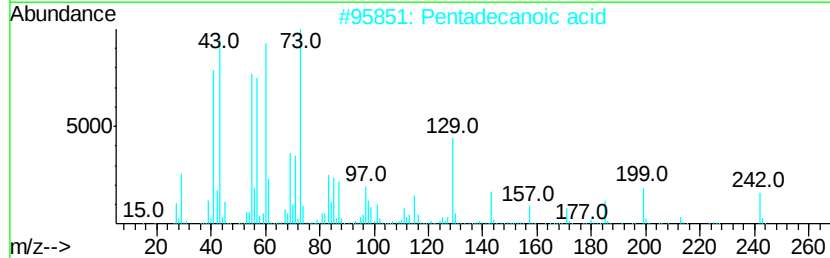
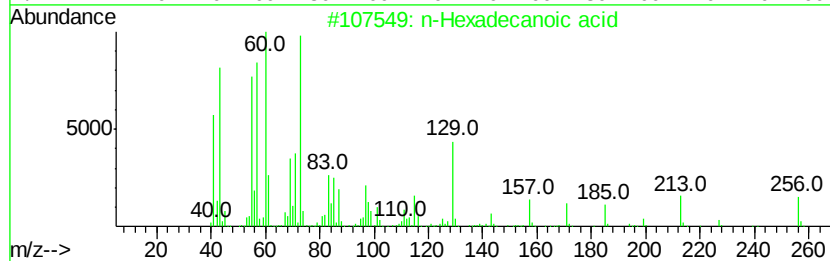
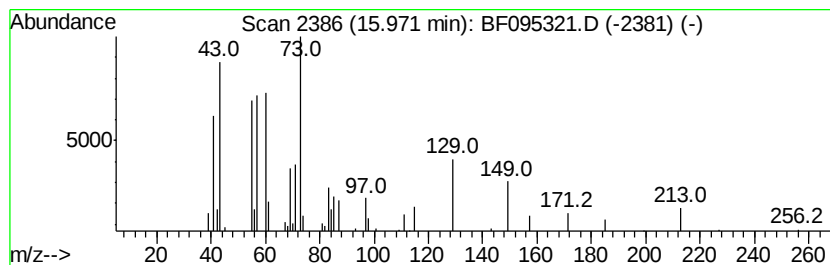
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	3.16 ng	101298	Phenanthrene-d10	15.01

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	64
3		Tridecanoic acid	214	C13H26O2	000638-53-9	64
4		n-Decanoic acid	172	C10H20O2	000334-48-5	50
5		Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	25



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ClientSampleId :
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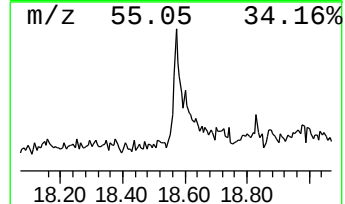
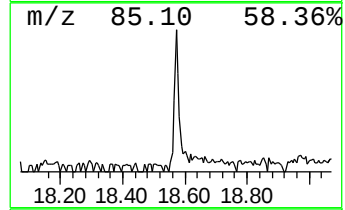
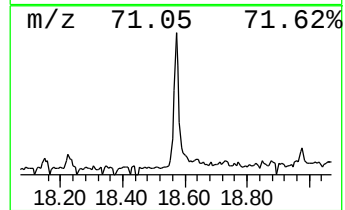
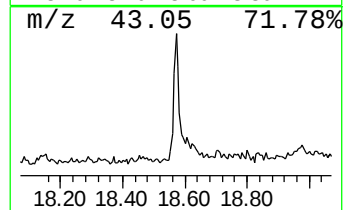
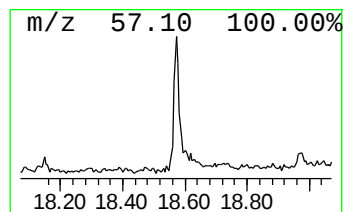
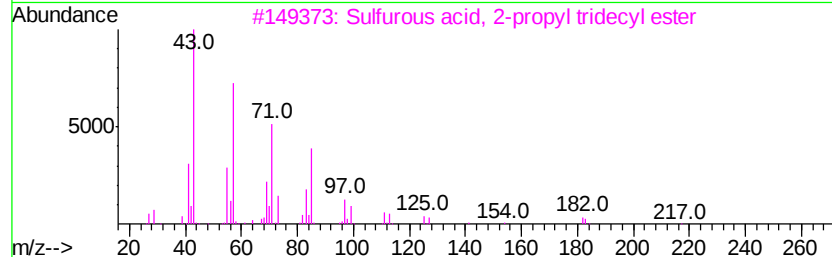
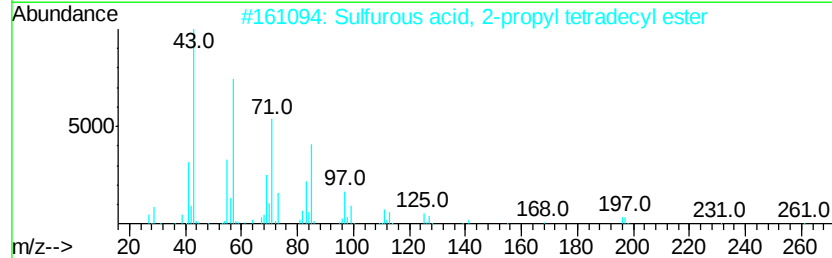
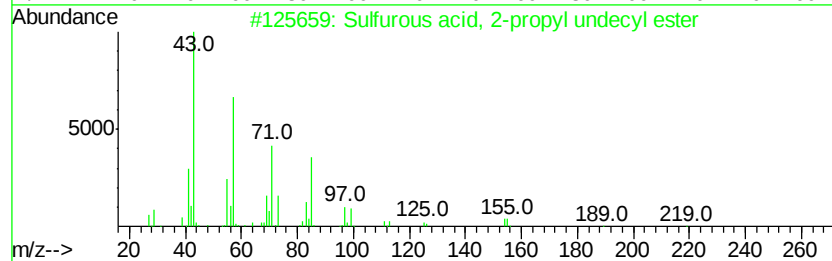
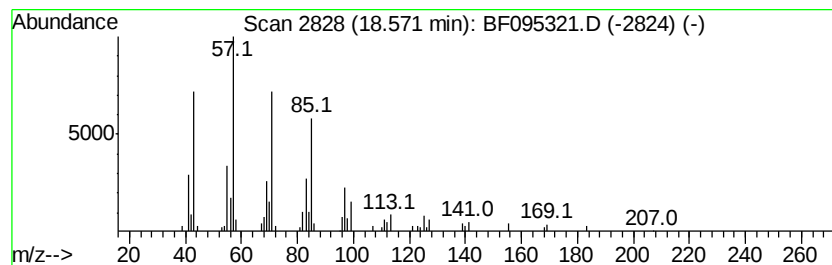
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Sulfurous acid, 2-propyl un... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.57	4.57 ng	102285	Chrysene-d12	18.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Sulfurous acid, 2-propyl undecyl...	278	C14H30O3S	1000309-12-2	87
2		Sulfurous acid, 2-propyl tetradecyl...	320	C17H36O3S	1000309-12-5	86
3		Sulfurous acid, 2-propyl tridecyl...	306	C16H34O3S	1000309-12-4	86
4		2-methyloctacosane	408	C29H60	1000376-72-8	80
5		Heneicosane	296	C21H44	000629-94-7	74



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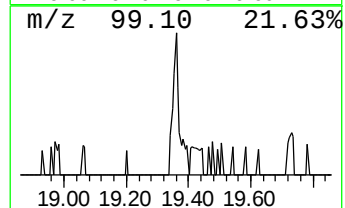
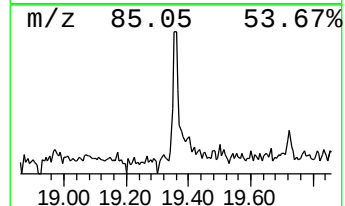
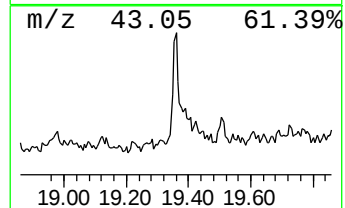
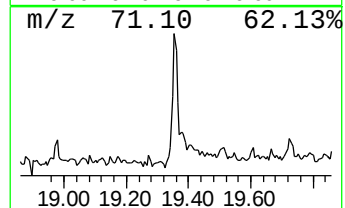
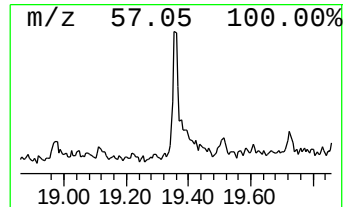
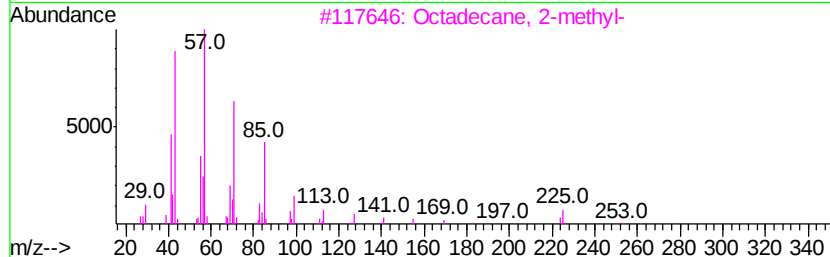
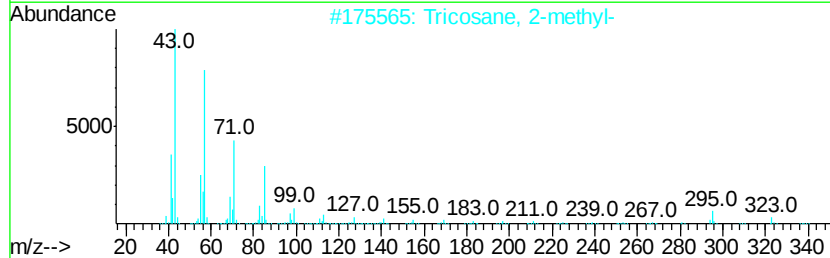
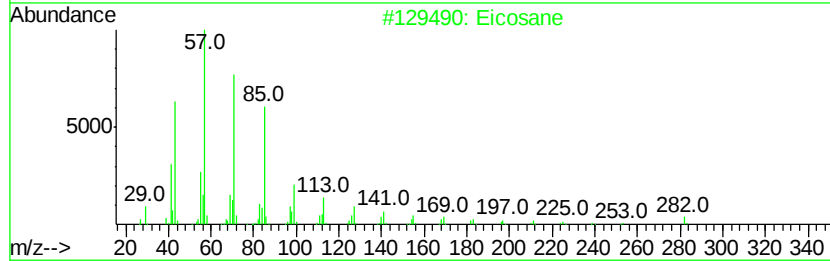
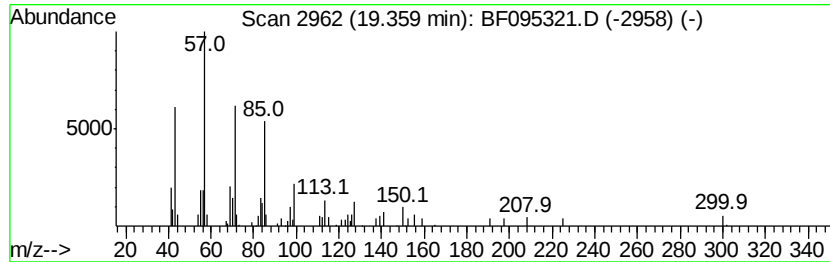
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Eicosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.36	2.63 ng	58981	Chrysene-d12	18.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	83
2		Tricosane, 2-methyl-	338	C24H50	001928-30-9	80
3		Octadecane, 2-methyl-	268	C19H40	001560-88-9	80
4		Octadecane	254	C18H38	000593-45-3	80
5		Heptadecane, 2-methyl-	254	C18H38	001560-89-0	80



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
Data File : BF095321.D
Acq On : 22 May 2017 18:58
Operator : SJ/MA
Sample : I3232-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-PS-01-003

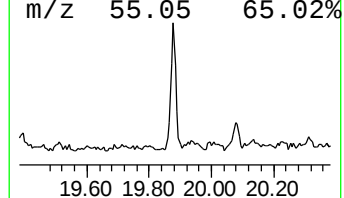
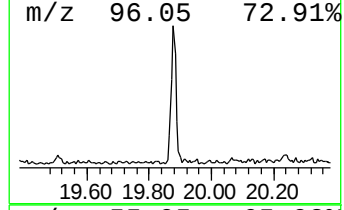
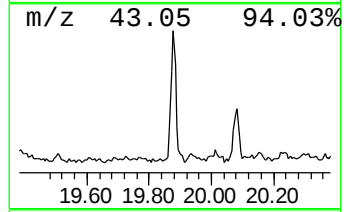
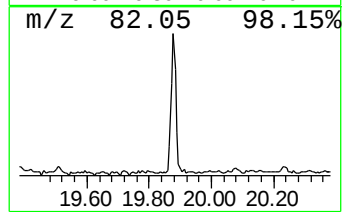
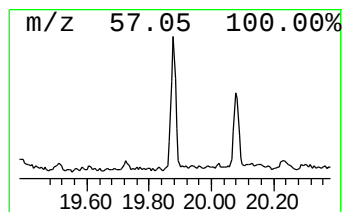
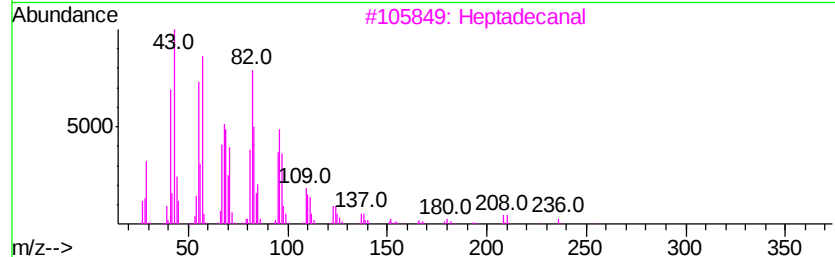
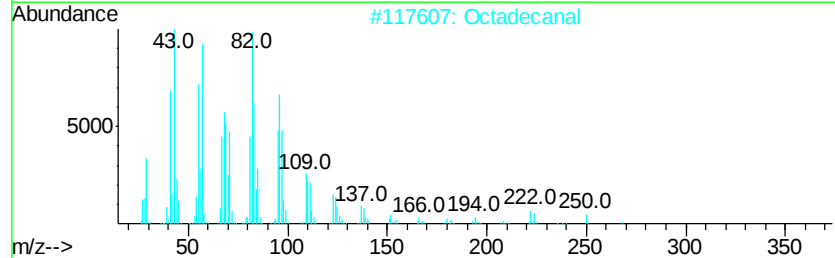
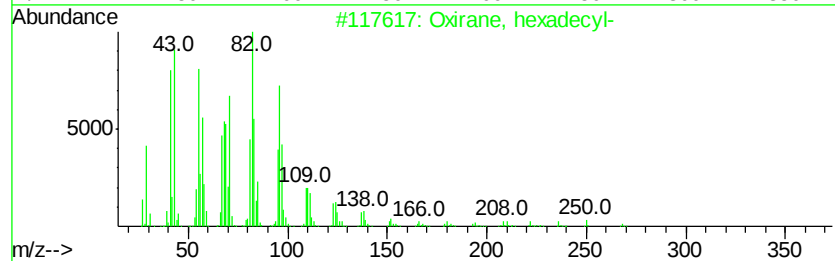
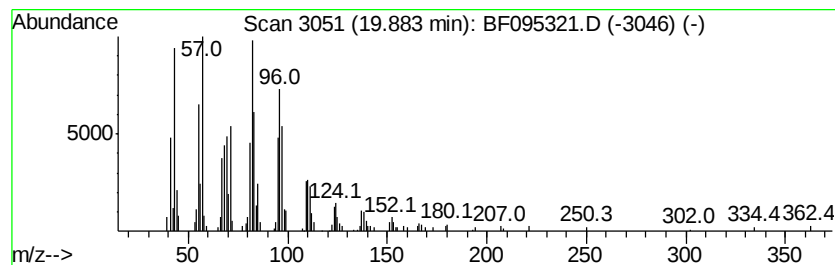
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Oxirane, hexadecyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.88	11.39 ng	263973	Perylene-d12	20.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
2		Octadecanal	268	C18H36O	000638-66-4	91
3		Heptadecanal	254	C17H34O	1000376-70-0	91
4		(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	87
5		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
Data File : BF095321.D
Acq On : 22 May 2017 18:58
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ClientSampleId :
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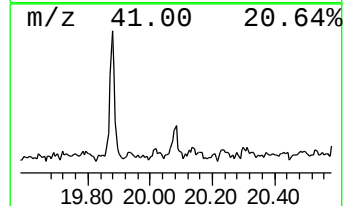
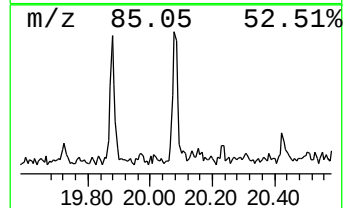
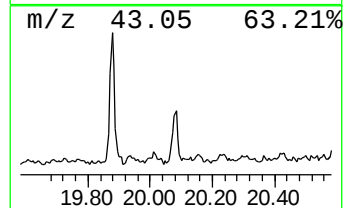
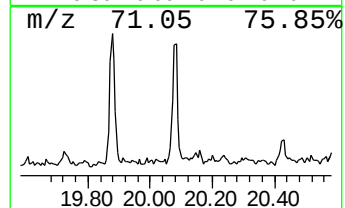
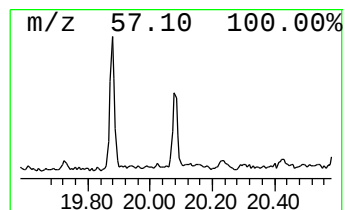
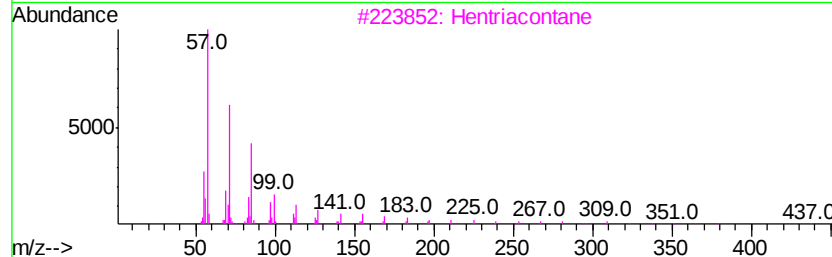
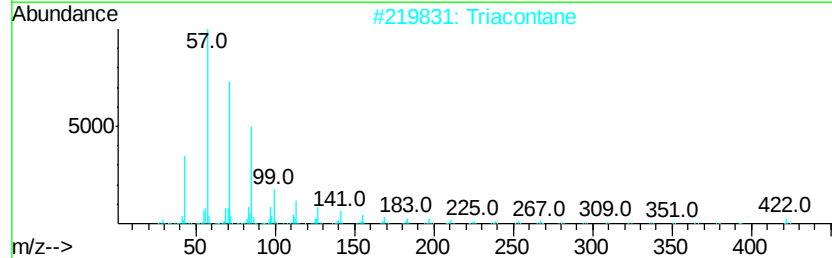
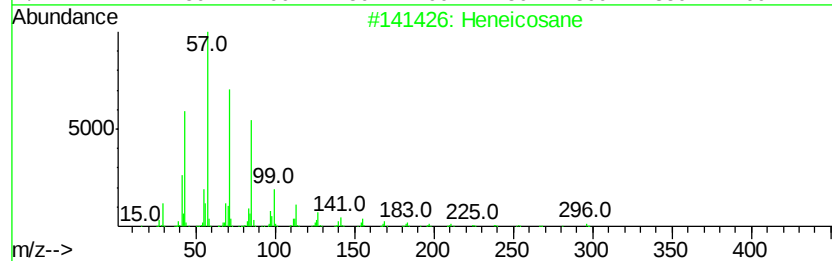
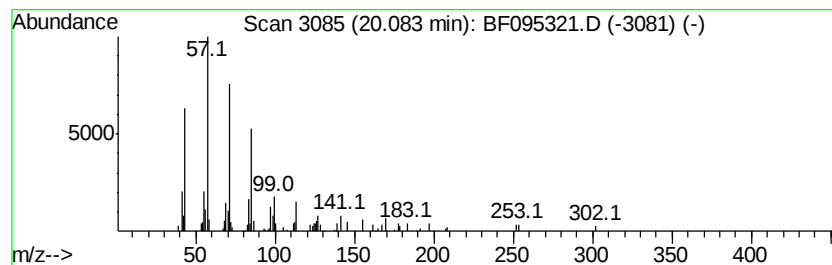
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Heneicosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.08	3.06 ng	70964	Perylene-d12	20.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heneicosane	296	C21H44	000629-94-7	90
2		Triacontane	422	C30H62	000638-68-6	90
3		Hentriacontane	437	C31H64	000630-04-6	87
4		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	87
5		Octacosane	394	C28H58	000630-02-4	87



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
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Operator : SJ/MA
Sample : I3232-01
Misc :
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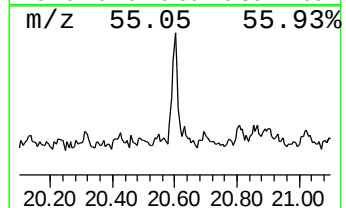
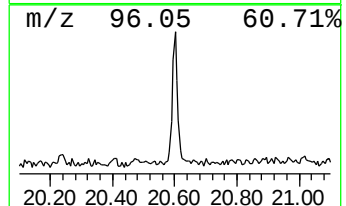
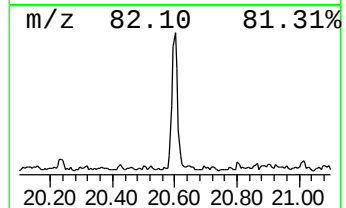
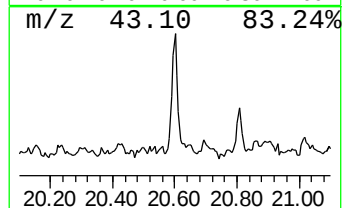
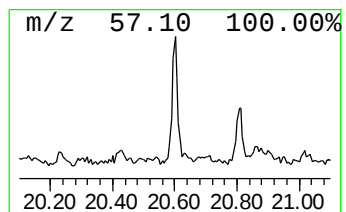
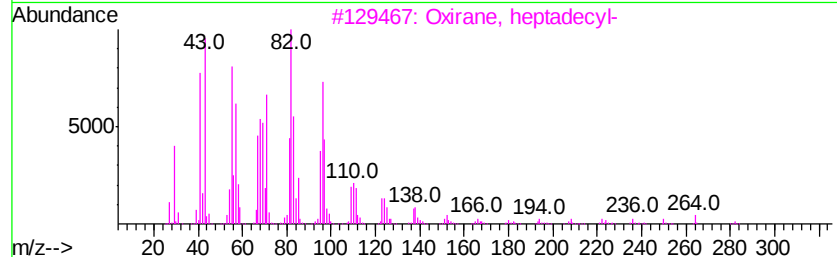
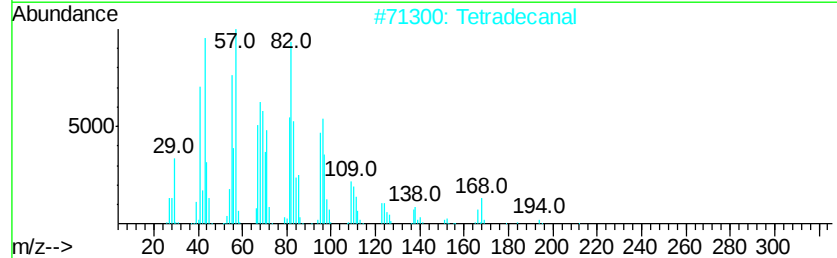
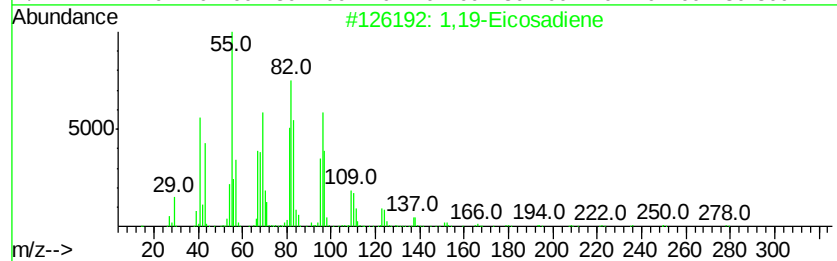
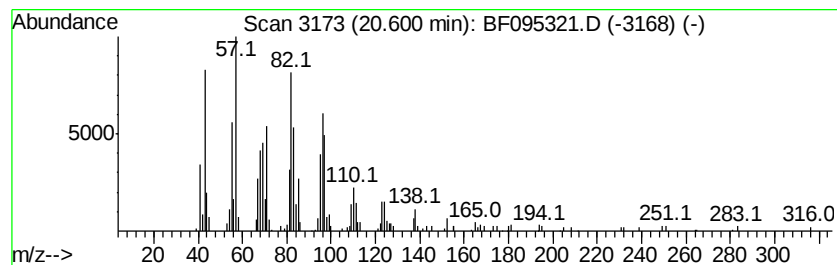
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 1,19-Eicosadiene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.60	8.40 ng	194799	Perylene-d12	20.36
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1,19-Eicosadiene	278 C20H38	014811-95-1	80
2	Tetradecanal	212 C14H28O	000124-25-4	72
3	Oxirane, heptadecyl-	282 C19H38O	067860-04-2	68
4	Oxirane, hexadecyl-	268 C18H36O	007390-81-0	68
5	Hexadecanal	240 C16H32O	000629-80-1	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
Data File : BF095321.D
Acq On : 22 May 2017 18:58
Operator : SJ/MA
Sample : I3232-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

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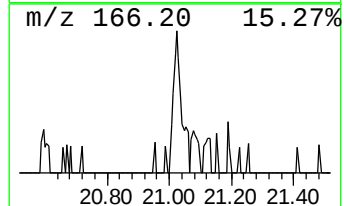
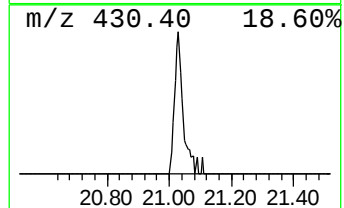
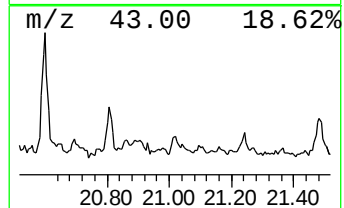
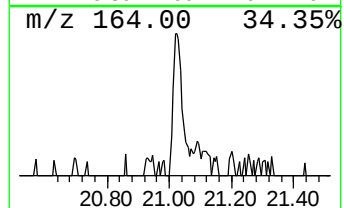
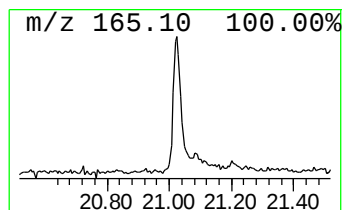
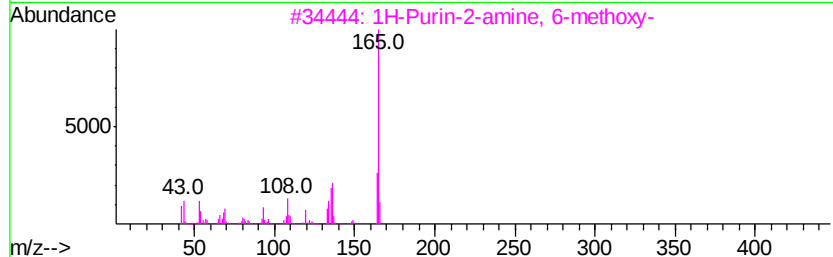
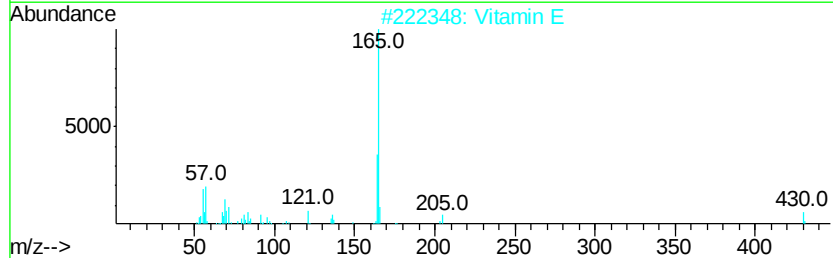
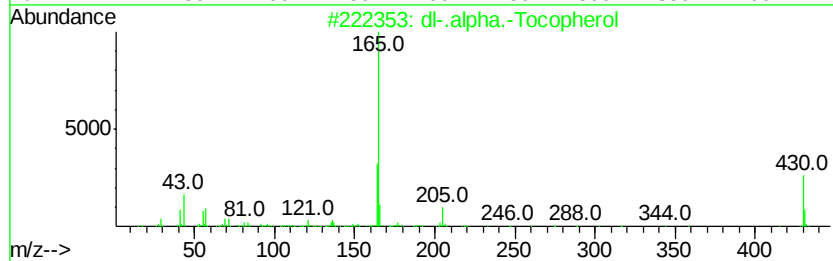
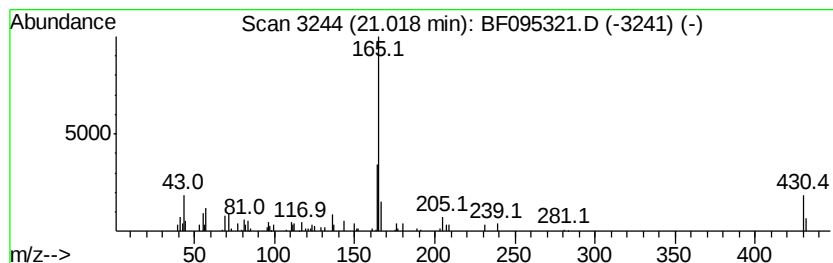
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 dl-.alpha.-Tocopherol Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.02	2.24 ng	52015	Perylene-d12	20.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	dl-.alpha.-Tocopherol	430	C29H50O2	010191-41-0	83
2		Vitamin E	430	C29H50O2	000059-02-9	80
3		1H-Purin-2-amine, 6-methoxy-	165	C6H7N5O	020535-83-5	50
4		Furo[3,4-c]pyridine-3,4(1H,5H)-d...	165	C8H7N03	007472-18-6	40
5		6H-Purin-6-one, 1,7-dihydro-2-(m...	165	C6H7N5O	010030-78-1	37



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
Data File : BF095321.D
Acq On : 22 May 2017 18:58
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ClientSampleId :
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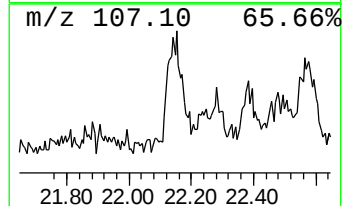
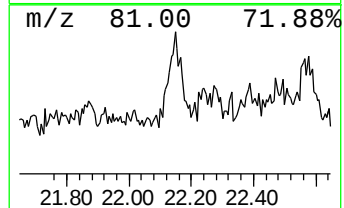
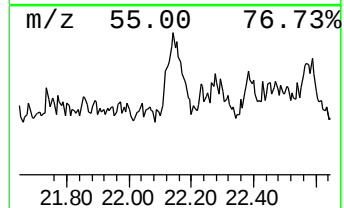
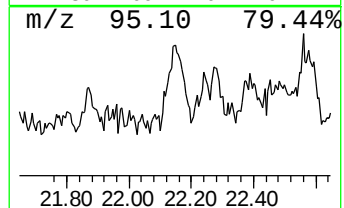
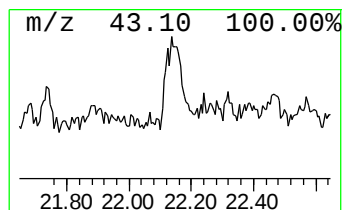
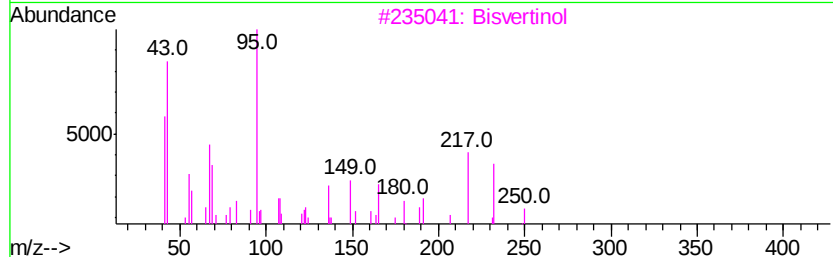
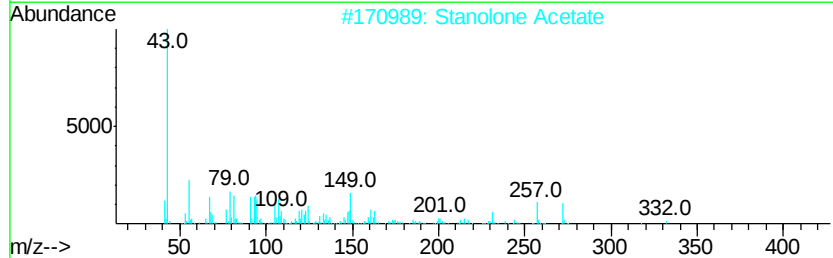
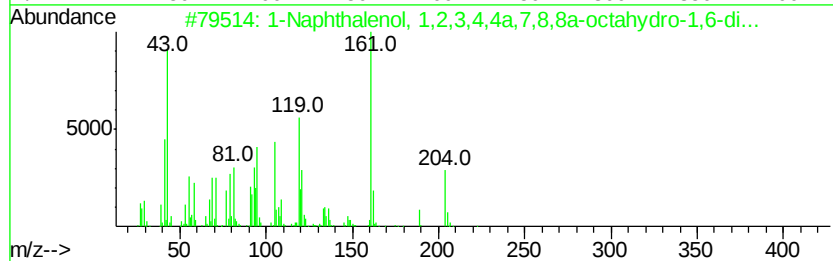
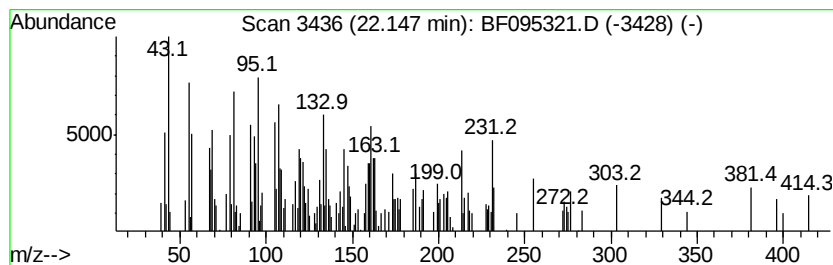
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown22.15 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.15	13.36 ng	309670	Perylene-d12	20.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Naphthalenol, 1,2,3,4,4a,7,8,8...	222	C15H26O	019435-97-3	22
2		Stanolone Acetate	332	C21H32O3	1000379-14-9	16
3		Bisvertinol	498	C28H34O8	103804-06-4	14
4		5-(4-Nitro-phenylcarbamoyl)-1H-i...	276	C11H8N4O5	1000286-29-9	14
5		Androst-5-en-7-one, 3-(acetyloxy...	358	C23H34O3	054549-89-2	10



Data Path : Z:\HPCHEM1\BNA F\DATA\BF052217\
Data File : BF095321.D
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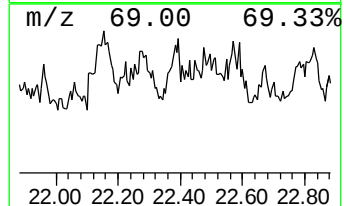
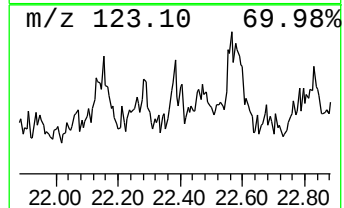
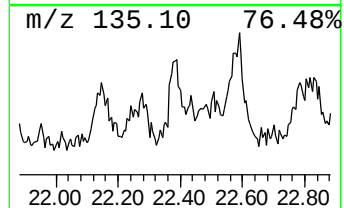
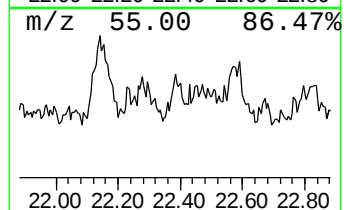
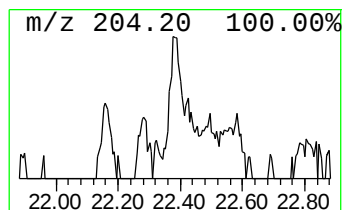
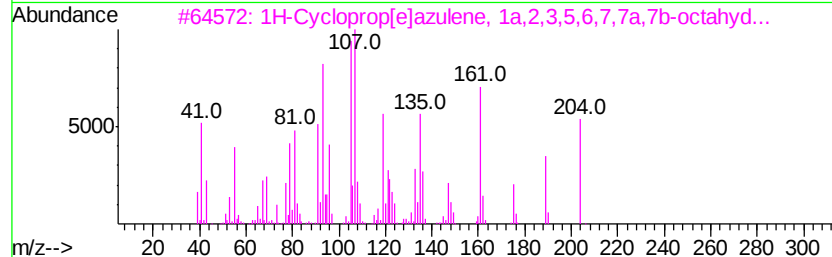
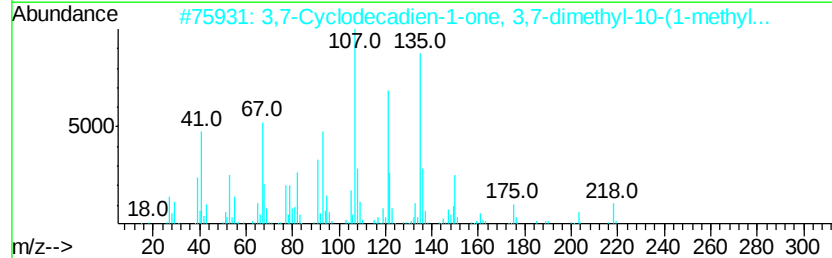
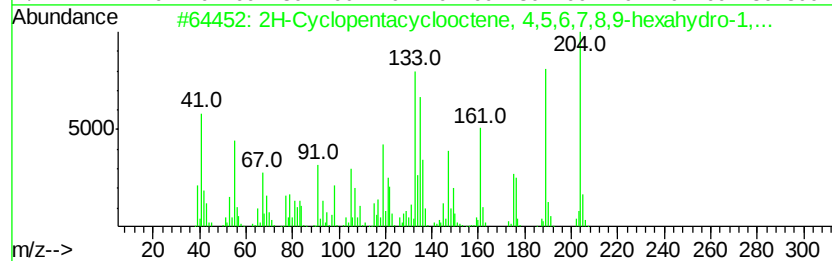
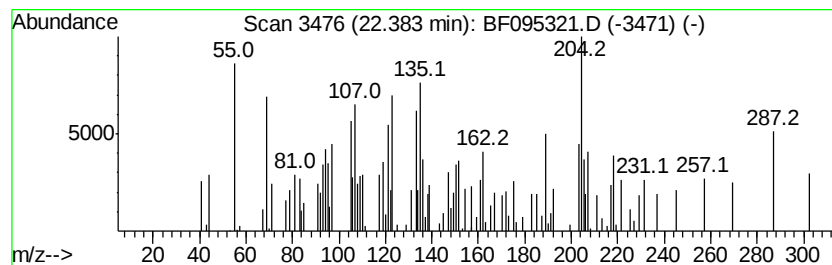
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown22.38 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.38	3.99 ng	92472	Perylene-d12	20.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-Cyclopentacyclooctene, 4,5,6,...	204	C15H24	1000221-85-8	49
2		3,7-Cyclodecadien-1-one, 3,7-dim...	218	C15H22O	006902-91-6	20
3		1H-Cycloprop[elazulene, 1a,2,3,5...	204	C15H24	021747-46-6	20
4		Boron, 1,5-cyclooctanedivl(2,4-p...	218	C13H23BN2	136704-99-9	18
5		.delta.-Selinene	204	C15H24	028624-23-9	15



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
Data File : BF095321.D
Acq On : 22 May 2017 18:58
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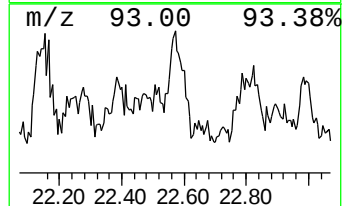
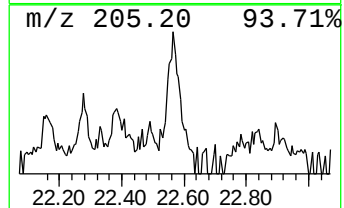
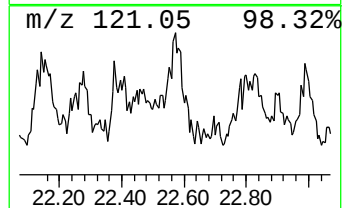
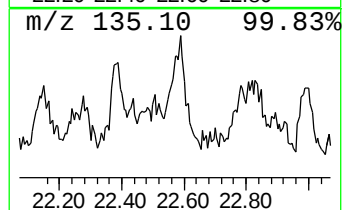
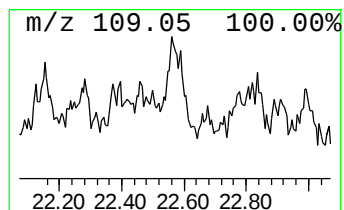
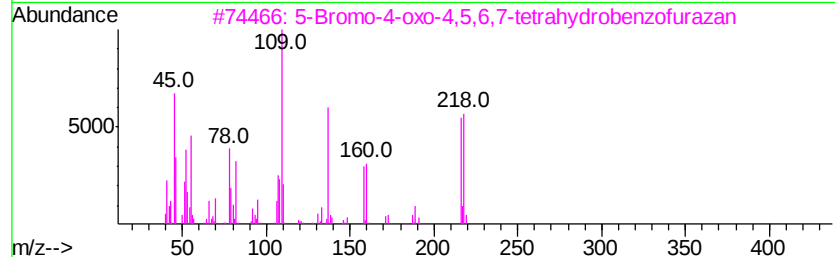
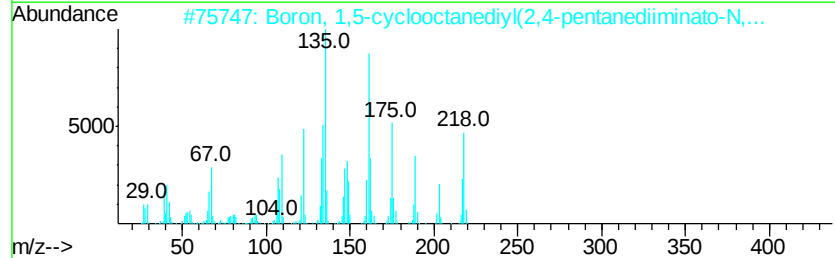
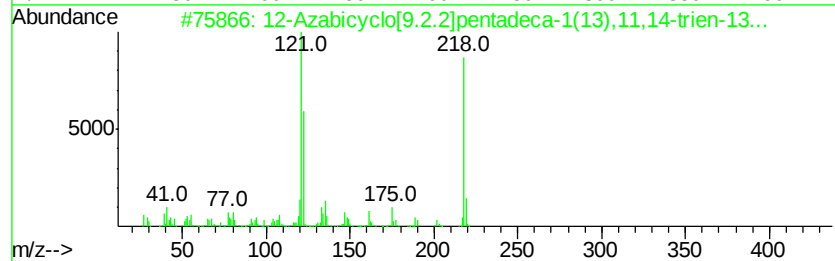
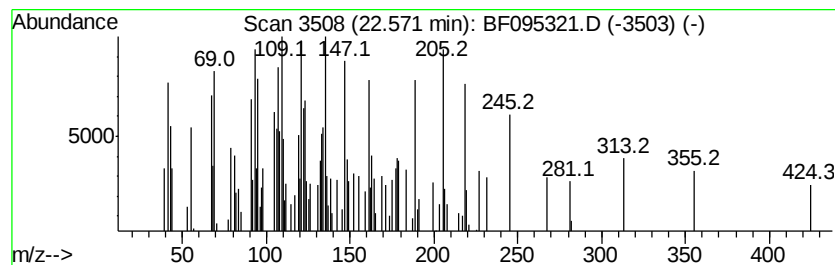
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 12-Azabicyclo[9.2.2]pentade... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.57	9.39 ng	217639	Perylene-d12	20.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	12-Azabicyclo[9.2.2]pentadeca-1(...	218	C14H22N2	1000185-25-0	62
2		Boron, 1,5-cyclooctanediyl(2,4-p...	218	C13H23BN2	136704-99-9	38
3		5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	30
4		2(3H)-Naphthalenone, 4,4a,5,6,7,...	218	C15H22O	019598-45-9	30
5		4,8-Methanoazulen-9-ol, decahydr...	222	C15H26O	004586-22-5	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
Data File : BF095321.D
Acq On : 22 May 2017 18:58
Operator : SJ/MA
Sample : I3232-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-PS-01-003

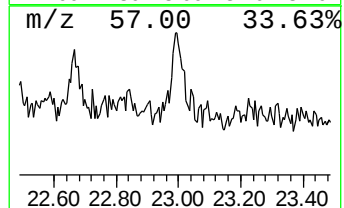
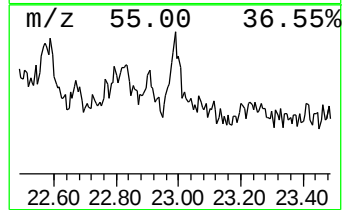
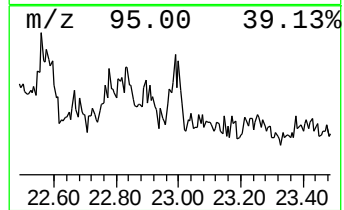
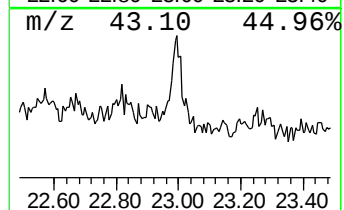
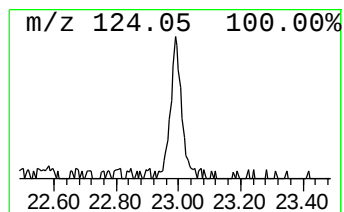
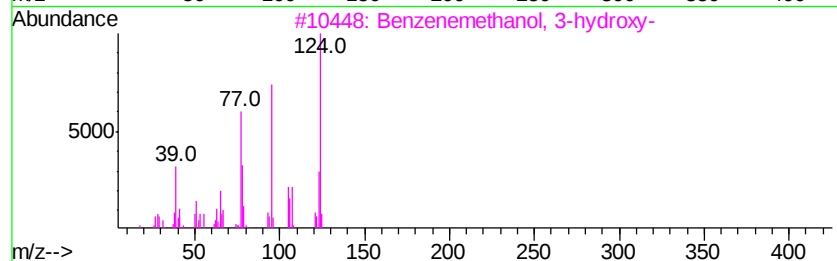
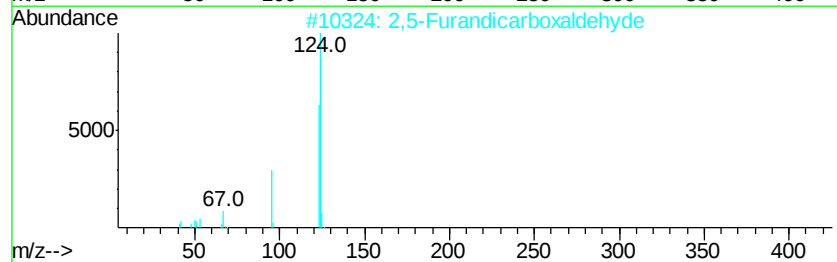
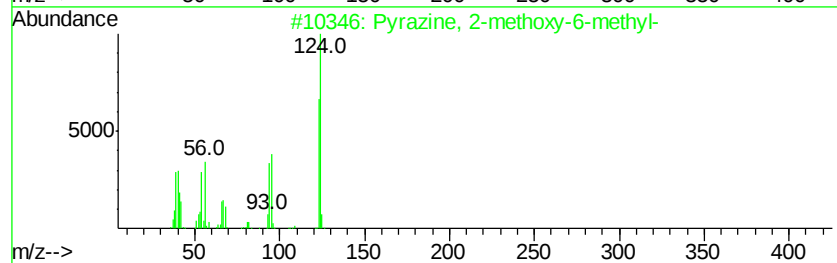
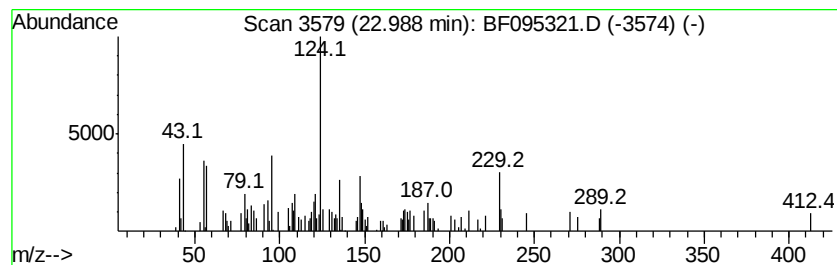
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 unknown22.99 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.99	7.01 ng	162552	Perylene-d12	20.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrazine, 2-methoxy-6-methyl-	124	C6H8N2O	002882-21-5	38
2		2,5-Furandicarboxaldehyde	124	C6H4O3	000823-82-5	38
3		Benzenemethanol, 3-hydroxy-	124	C7H8O2	000620-24-6	38
4		5-Ethylcyclopent-1-enecarboxalde...	124	C8H12O	036431-60-4	37
5		Testosterone	288	C19H28O2	000058-22-0	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
Data File : BF095321.D
Acq On : 22 May 2017 18:58
Operator : SJ/MA
Sample : I3232-01
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
FK-PS-01-003

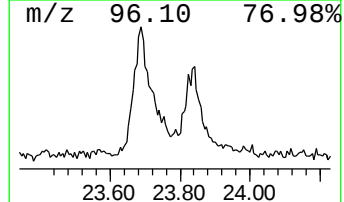
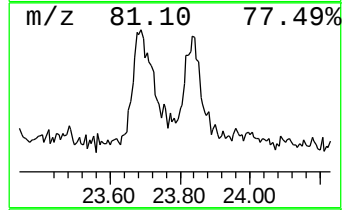
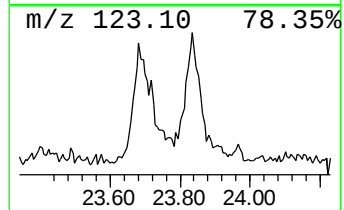
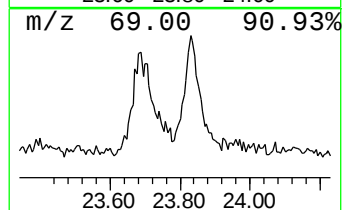
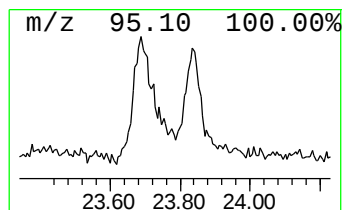
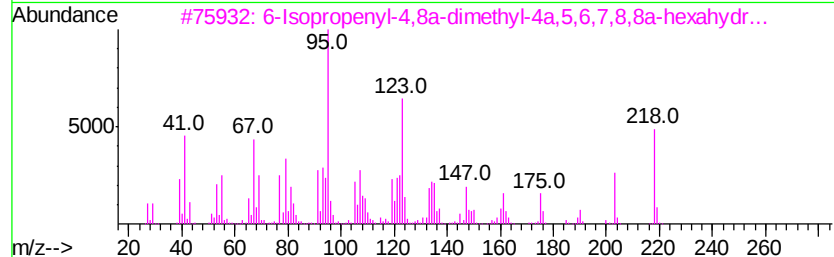
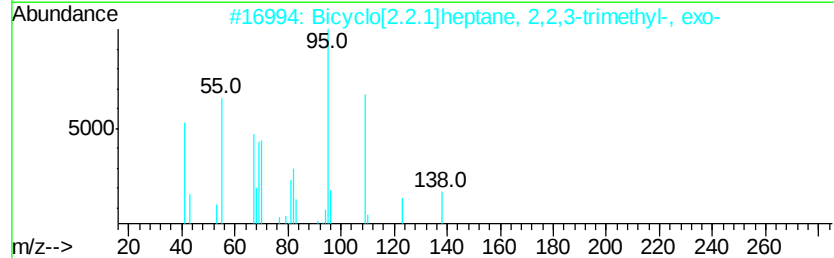
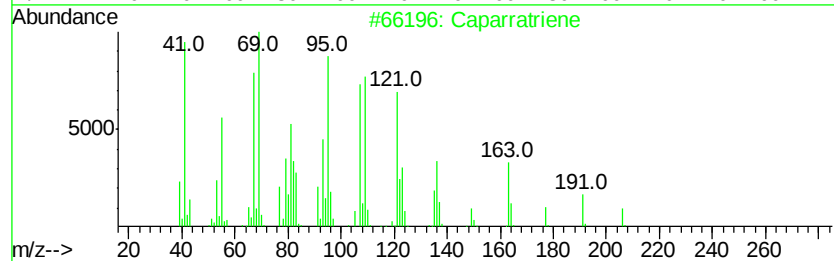
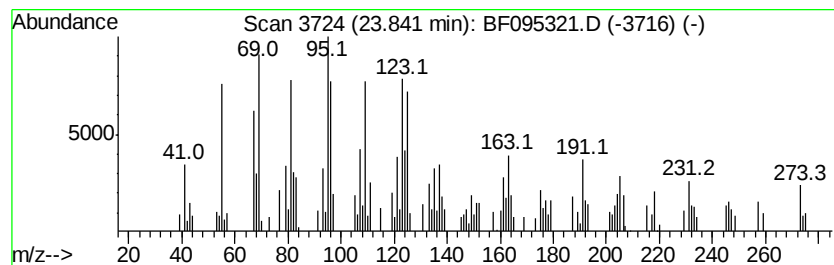
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 unknown23.84 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.84	14.84 ng	344085	Perylene-d12	20.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Caparratriene	206	C15H26	1000374-08-7	43
2		Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-41-8	38
3		6-Isopropenyl-4,8a-dimethyl-4a,5...	218	C15H22O	086917-79-5	38
4		Bicyclo[2.2.1]heptane, 2,2,3-tri...	138	C10H18	020536-40-7	38
5		Cyclopentane, 1-methyl-1-(2-meth...	138	C10H18	074764-47-9	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF052217\
 Data File : BF095321.D
 Acq On : 22 May 2017 18:58
 Operator : SJ/MA
 Sample : I3232-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-PS-01-003

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.12	17.7	ng	345623	1	7.76	389566	20.0
unknown7.42	7.42	127.1	ng	2474690	1	7.76	389566	20.0
n-Hexadecanoic acid	15.97	3.2	ng	101298	4	15.01	642116	20.0
Sulfurous acid, 2...	18.57	4.6	ng	102285	5	18.68	447771	20.0
Eicosane	19.36	2.6	ng	58981	5	18.68	447771	20.0
Oxirane, hexadecyl-	19.88	11.4	ng	263973	6	20.36	463583	20.0
Heneicosane	20.08	3.1	ng	70964	6	20.36	463583	20.0
1,19-Eicosadiene	20.60	8.4	ng	194799	6	20.36	463583	20.0
dl-.alpha.-Tocoph...	21.02	2.2	ng	52015	6	20.36	463583	20.0
unknown22.15	22.15	13.4	ng	309670	6	20.36	463583	20.0
unknown22.38	22.38	4.0	ng	92472	6	20.36	463583	20.0
12-Azabicyclo[9.2...	22.57	9.4	ng	217639	6	20.36	463583	20.0
unknown22.99	22.99	7.0	ng	162552	6	20.36	463583	20.0
unknown23.84	23.84	14.8	ng	344085	6	20.36	463583	20.0