

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
 Data File : BF082297.D
 Acq On : 15 Oct 2015 9:09
 Operator : UM/IZ
 Sample : G4016-04
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CS-4

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-BF101015.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	4.308	187	190	193	rBV	1397313	2324860	32.04%	9.421%
2	5.040	246	254	256	rBV	2426752	7255320	100.00%	29.402%
3	6.457	375	378	388	rBV	209897	263281	3.63%	1.067%
4	6.823	407	410	412	rBV	472050	427310	5.89%	1.732%
5	6.983	421	424	426	rBV	1022907	1371359	18.90%	5.557%
6	7.394	457	460	462	rBV	1149633	1106329	15.25%	4.483%
7	8.114	520	523	525	rBV	581274	586430	8.08%	2.376%
8	9.189	615	617	620	rBV	2073030	2157173	29.73%	8.742%
9	9.589	650	652	653	rBV	347010	279013	3.85%	1.131%
10	9.875	674	677	679	rBV	533629	706935	9.74%	2.865%
11	10.766	754	755	760	rVB2	132583	189962	2.62%	0.770%
12	11.361	805	807	809	rBV	730529	725165	9.99%	2.939%
13	11.578	824	826	831	rBV	52221	73993	1.02%	0.300%
14	12.138	873	875	878	rBV	97614	84492	1.16%	0.342%
15	12.961	944	947	949	rBV	2418086	2461474	33.93%	9.975%
16	13.452	988	990	993	rBV	49259	89409	1.23%	0.362%
17	13.898	1027	1029	1033	rVB	77944	130214	1.79%	0.528%
18	14.069	1042	1044	1049	rBV	541789	622609	8.58%	2.523%
19	14.607	1089	1091	1093	rBV2	49462	88294	1.22%	0.358%
20	14.938	1117	1120	1125	rBV	1216424	1482511	20.43%	6.008%
21	15.578	1174	1176	1178	rVB	94409	115883	1.60%	0.470%
22	15.635	1178	1181	1185	rBV	291503	491634	6.78%	1.992%
23	15.761	1190	1192	1194	rVB	94832	131134	1.81%	0.531%
24	16.310	1237	1240	1250	rVB	219492	586840	8.09%	2.378%
25	16.652	1268	1270	1274	rBV3	45603	132093	1.82%	0.535%
26	16.732	1274	1277	1285	rVB	256443	511576	7.05%	2.073%
27	17.293	1323	1326	1329	rVV	70614	145609	2.01%	0.590%
28	17.361	1330	1332	1336	rVB	62822	135330	1.87%	0.548%

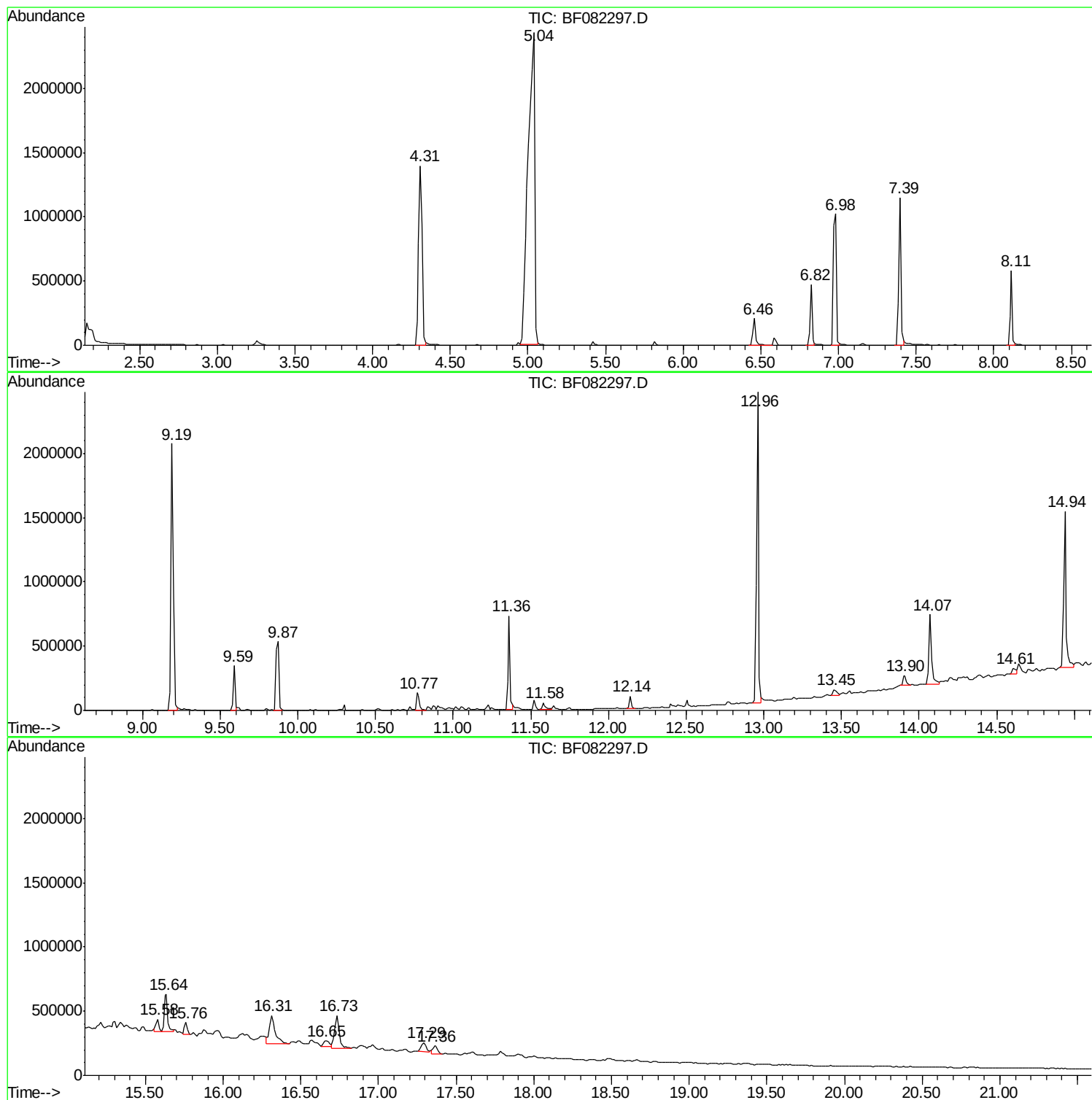
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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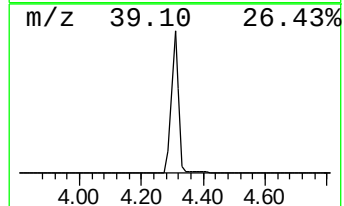
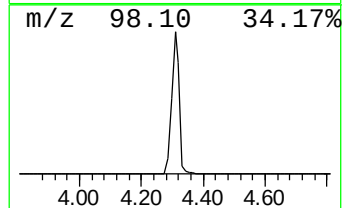
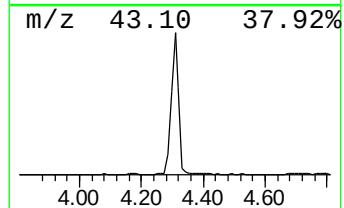
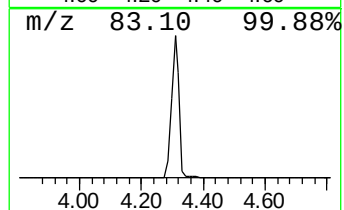
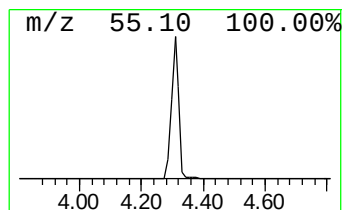
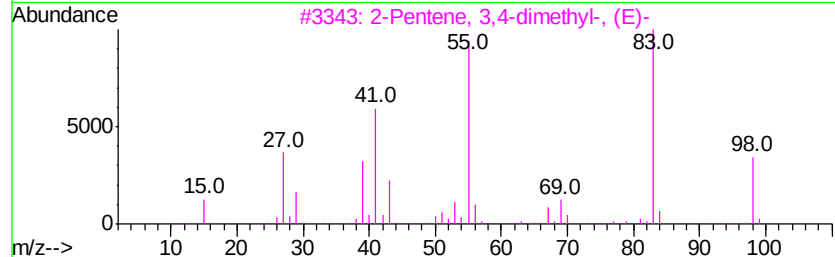
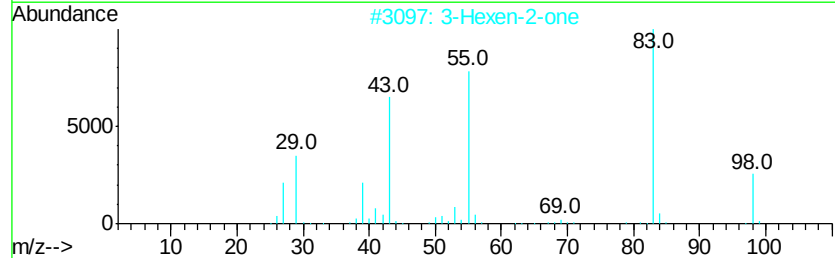
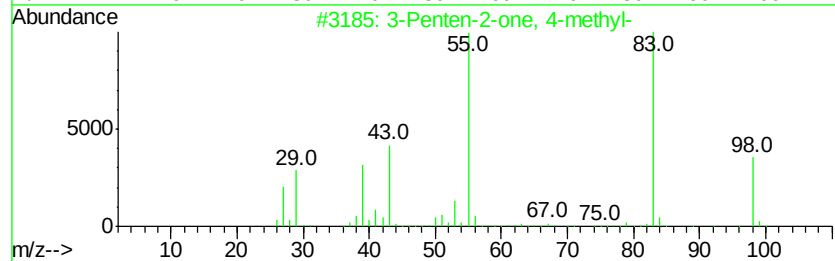
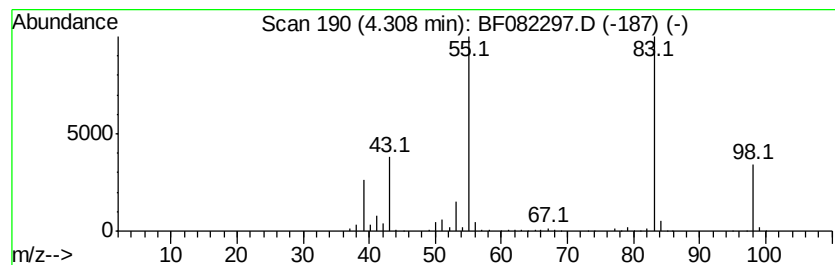
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.31	108.81 ng	2324860	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	94
2		3-Hexen-2-one	98	C6H10O	000763-93-9	91
3		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90
4		2-Pentene, 3,4-dimethyl-, (Z)-	98	C7H14	004914-91-4	86
5		2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	86



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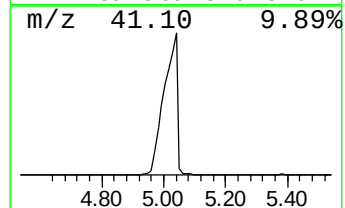
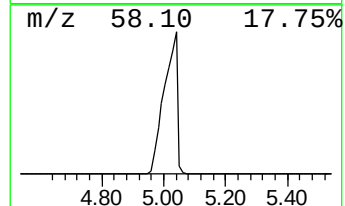
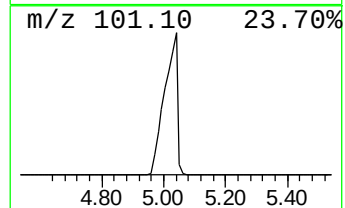
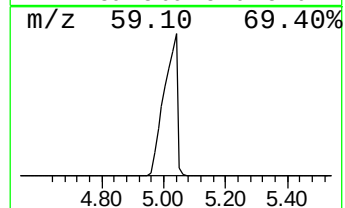
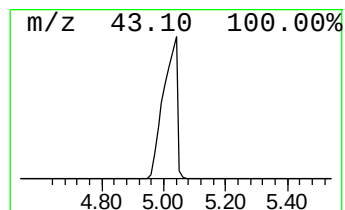
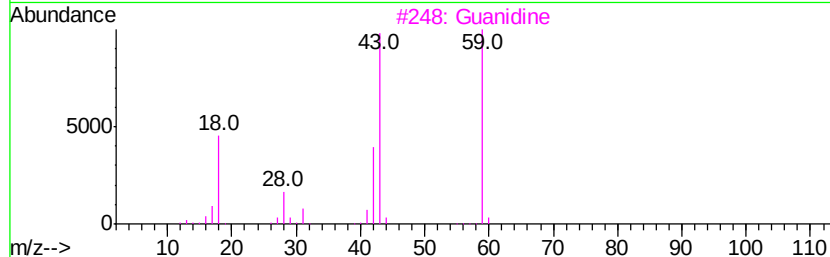
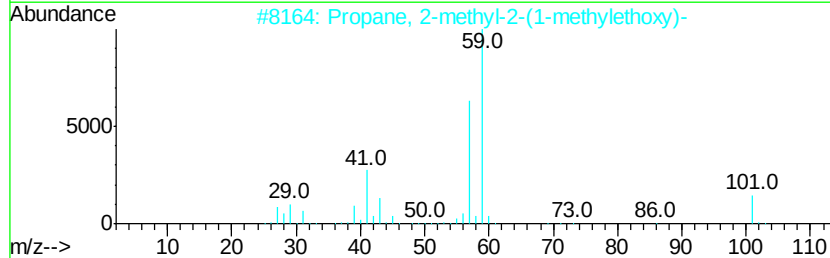
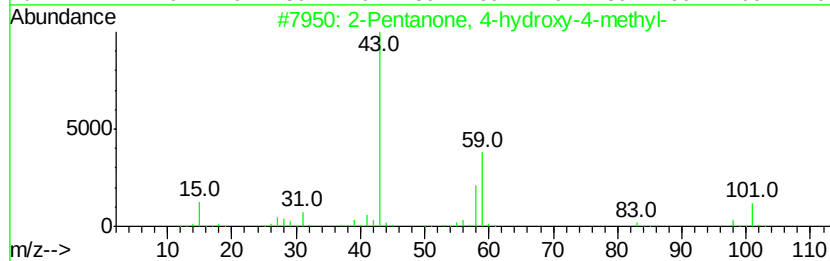
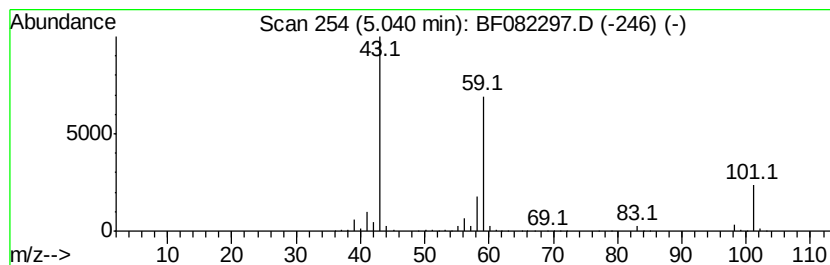
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
5.04	339.58 ng	7255320	1,4-Dichlorobenzene-d4	6.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	53
3	Guanidine	59	CH5N3	000113-00-8	43
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



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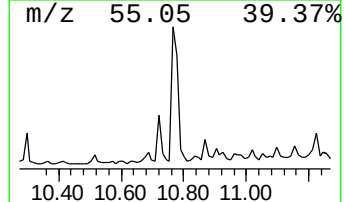
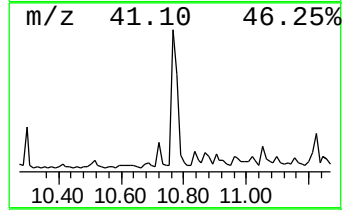
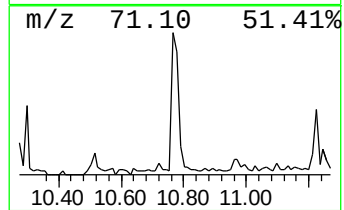
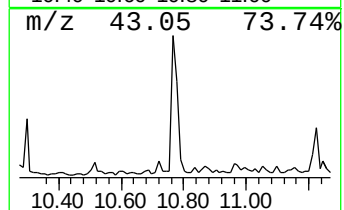
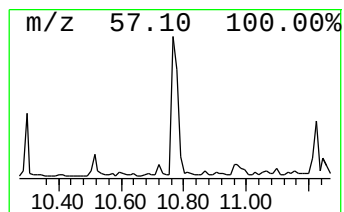
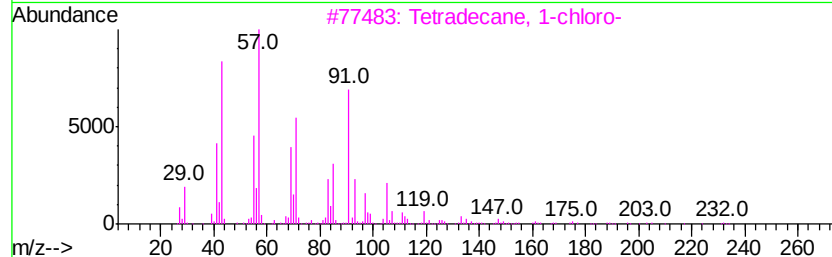
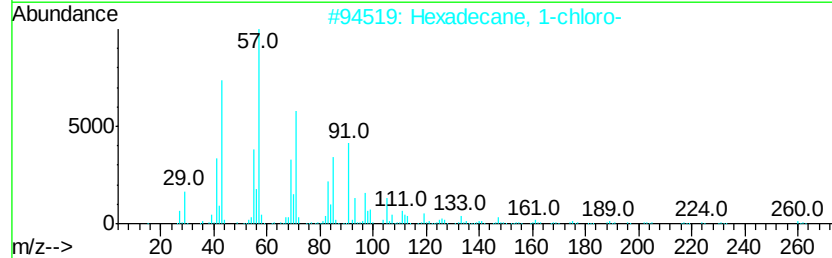
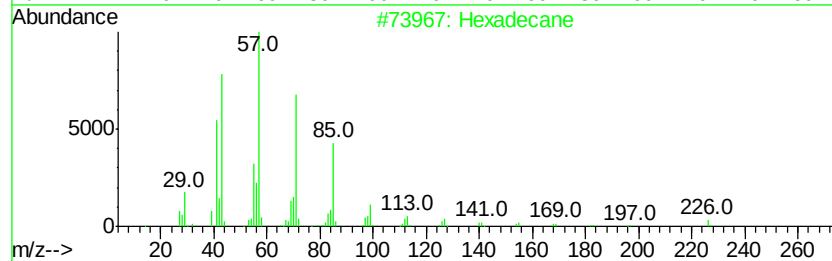
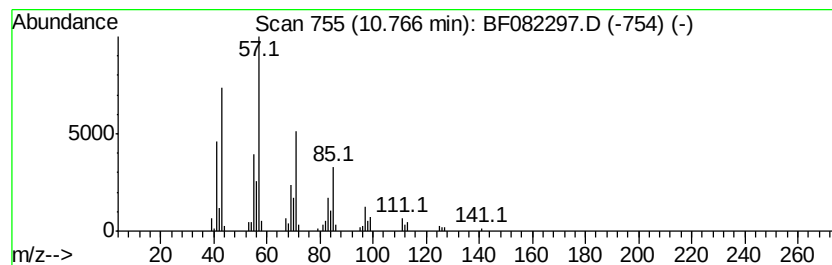
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Peak Number 4 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.77	5.24 ng	189962	Phenanthrene-d10	11.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	86
2	Hexadecane, 1-chloro-		260	C16H33Cl	004860-03-1	80
3	Tetradecane, 1-chloro-		232	C14H29Cl	002425-54-9	72
4	Tetradecane		198	C14H30	000629-59-4	72
5	Hexatriacontane		507	C36H74	000630-06-8	72



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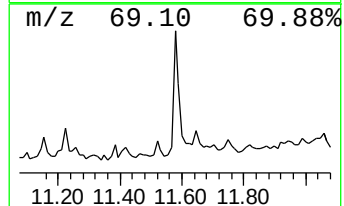
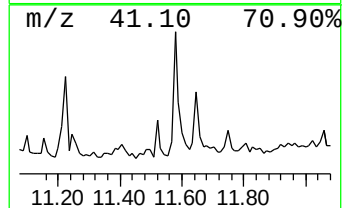
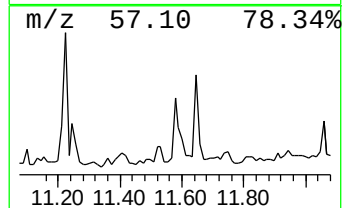
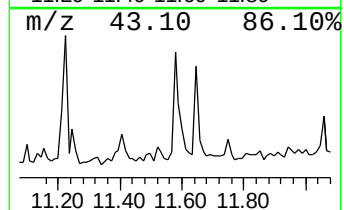
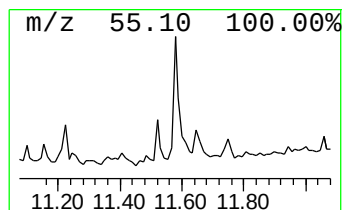
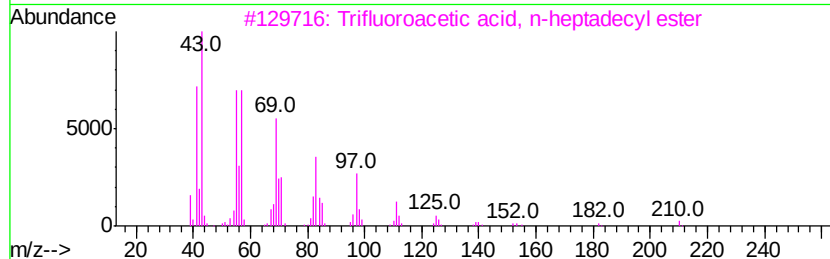
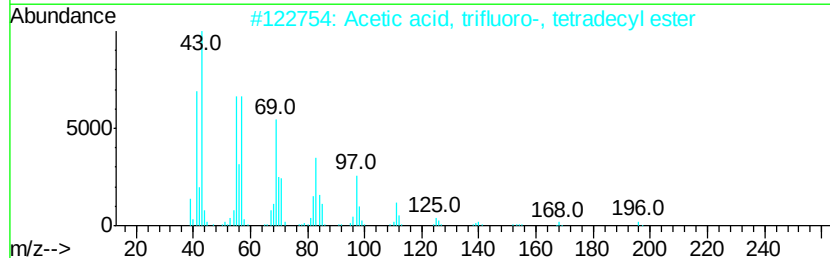
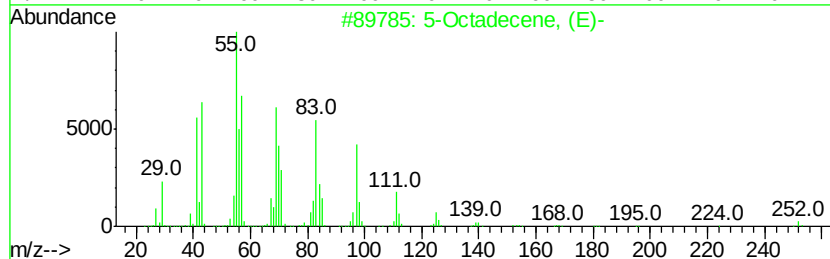
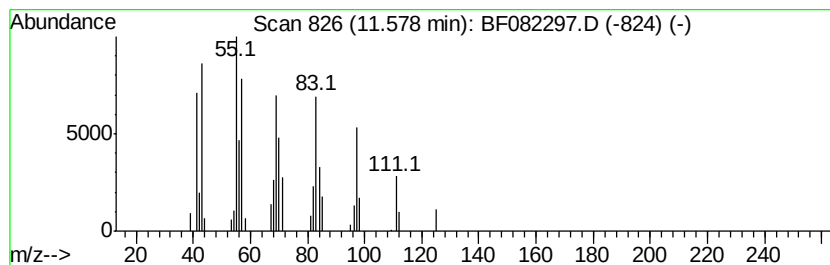
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Peak Number 5 5-Octadecene, (E)- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.58	2.04 ng	73993	Phenanthrene-d10	11.36		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Octadecene, (E)-	252	C18H36	007206-21-5	91
2		Acetic acid, trifluoro-, tetradec...	310	C16H29F3O2	006222-02-2	90
3		Trifluoroacetic acid, n-heptadec...	324	C17H31F3O2	1000216-79-2	83
4		Pentafluoropropionic acid, tride...	346	C16H27F5O2	1000280-07-3	80
5		Cyclotetradecane	196	C14H28	000295-17-0	72



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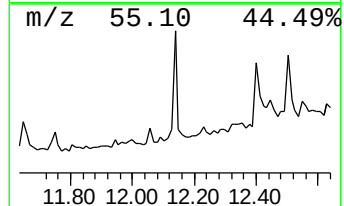
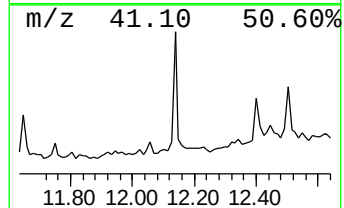
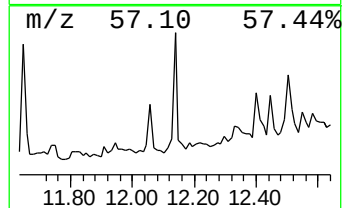
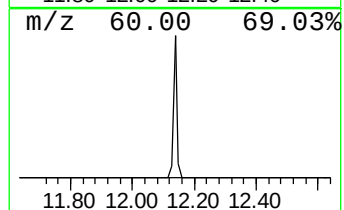
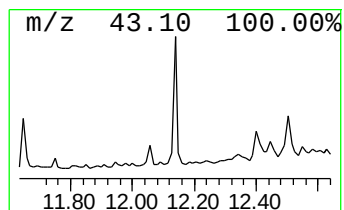
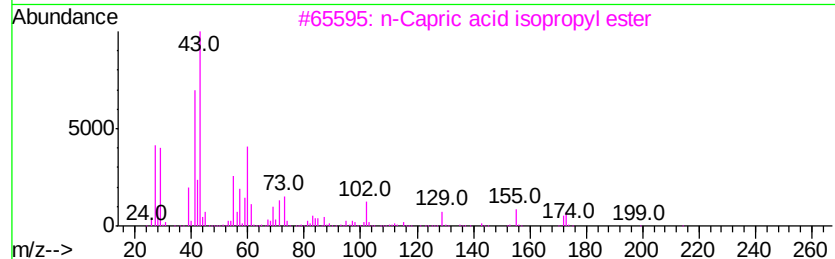
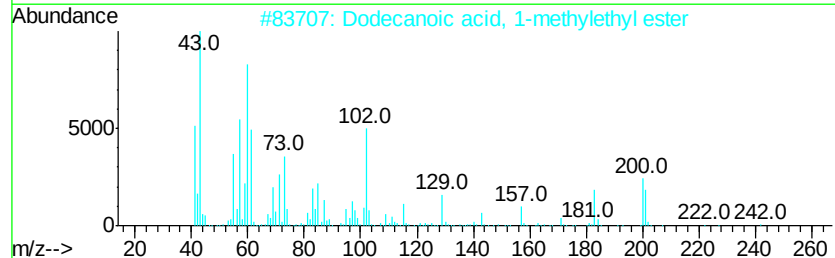
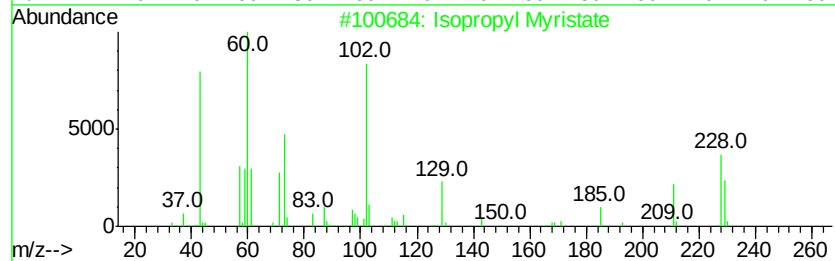
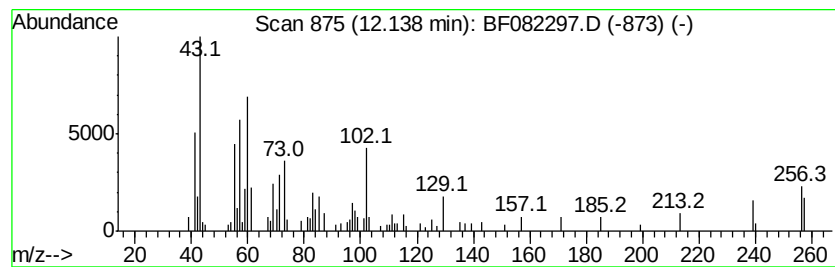
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Peak Number 6 Isopropyl Myristate Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.14	2.33 ng	84492	Phenanthrene-d10	11.36
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Isopropyl Myristate	270 C17H34O2	000110-27-0	53
2	Dodecanoic acid, 1-methylethyl e...	242 C15H30O2	010233-13-3	49
3	n-Capric acid isopropyl ester	214 C13H26O2	002311-59-3	47
4	n-Hexadecanoic acid	256 C16H32O2	000057-10-3	42
5	.alpha.-D-Galactopyranoside, methyl	194 C7H14O6	003396-99-4	27



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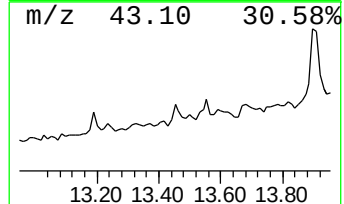
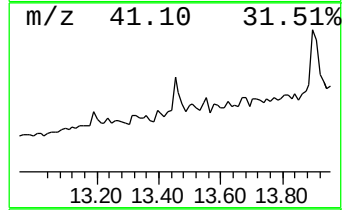
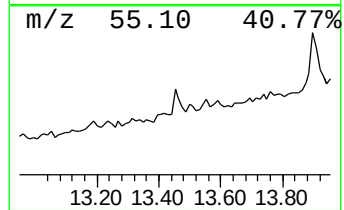
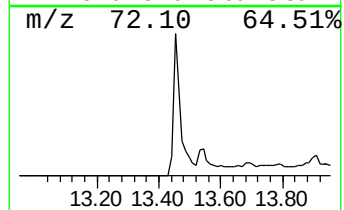
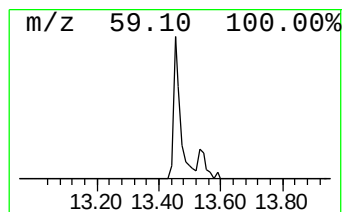
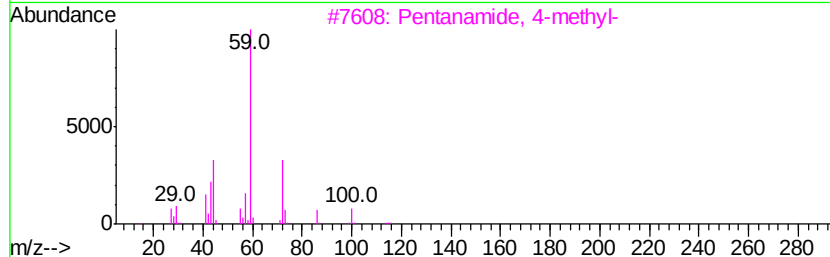
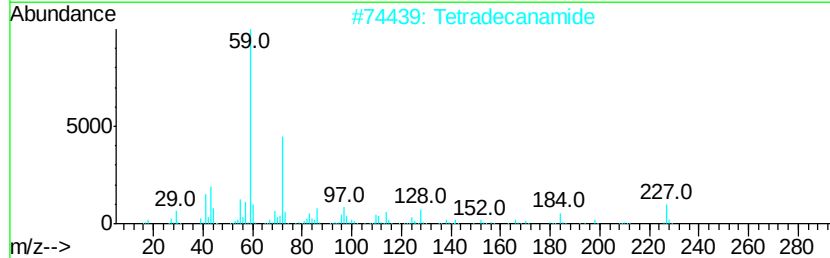
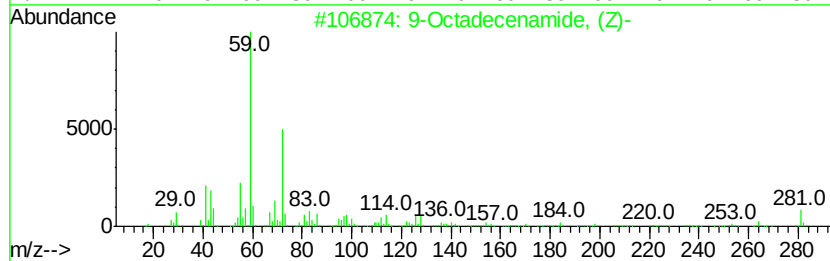
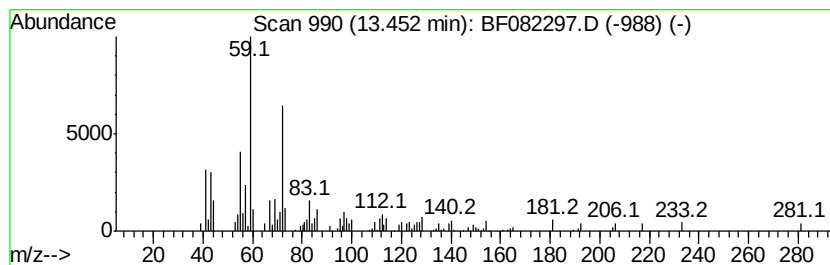
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 9-Octadecenamide, (Z)- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.45	2.87 ng	89409	Chrysene-d12	14.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	83
2		Tetradecanamide	227	C14H29NO	000638-58-4	72
3		Pentanamide, 4-methyl-	115	C6H13NO	001119-29-5	64
4		7-Nonenamide	155	C9H17NO	090949-53-4	59
5		Octanamide	143	C8H17NO	000629-01-6	53



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082297.D
Acq On : 15 Oct 2015 9:09
Operator : UM/IZ
Sample : G4016-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

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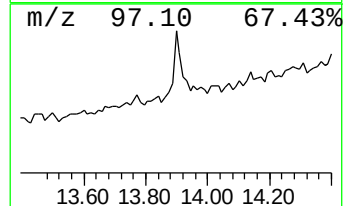
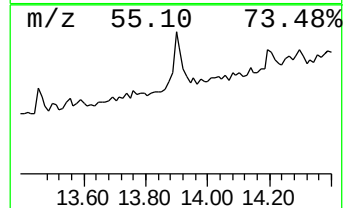
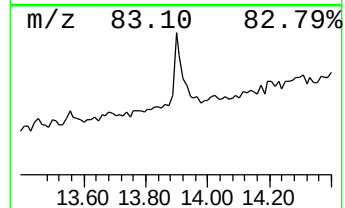
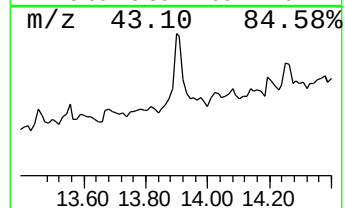
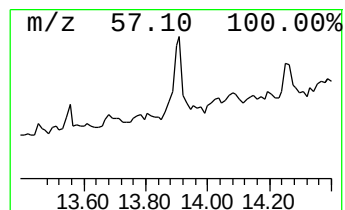
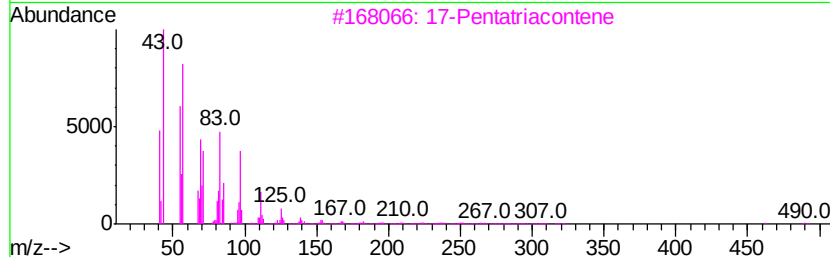
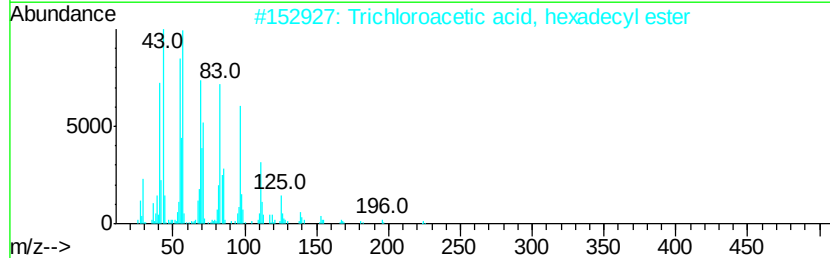
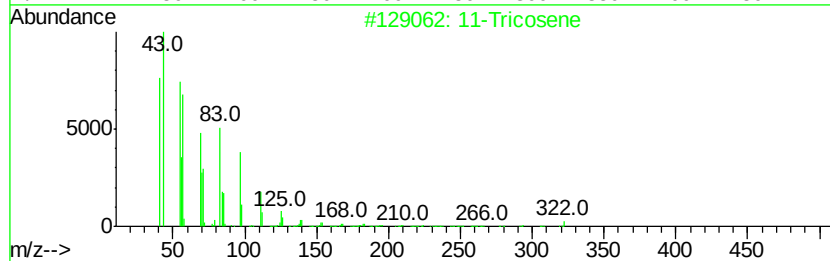
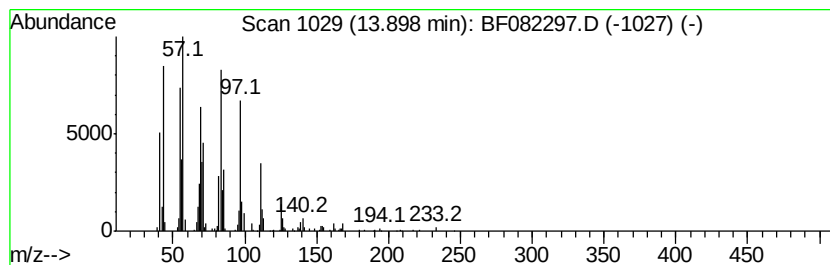
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 11-Tricosene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	4.18 ng	130214	Chrysene-d12	14.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11-Tricosene	322	C23H46	052078-56-5	91
2		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	91
3		17-Pentatriacontene	491	C35H70	006971-40-0	91
4		3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	87
5		Heptafluorobutyric acid, n-penta...	424	C19H31F7O2	1000216-79-3	81



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082297.D
Acq On : 15 Oct 2015 9:09
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Sample : G4016-04
Misc :
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Instrument :
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ClientSampleId :
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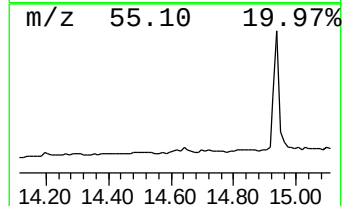
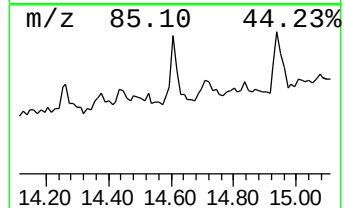
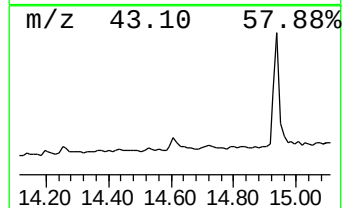
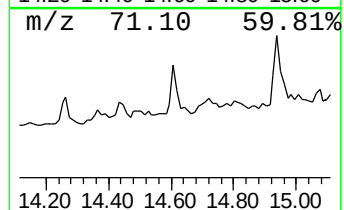
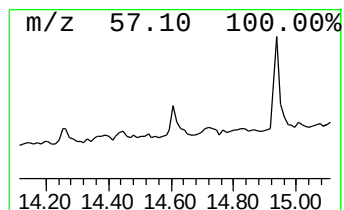
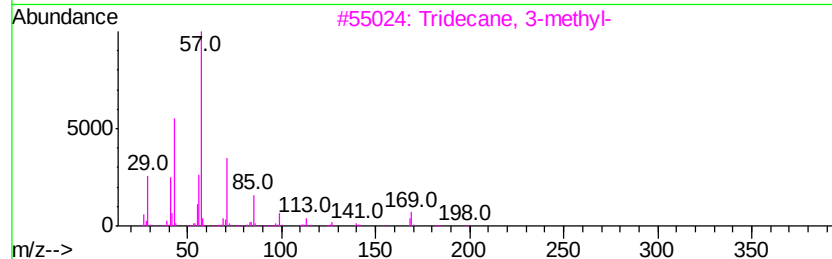
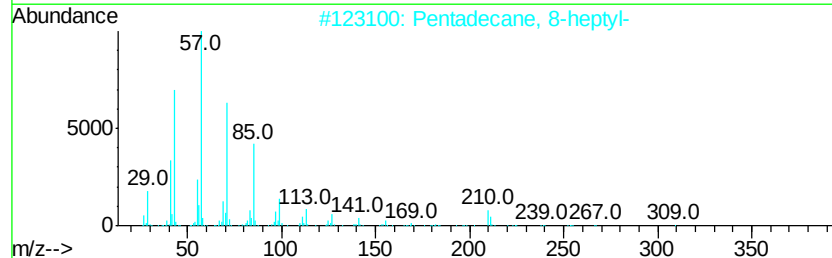
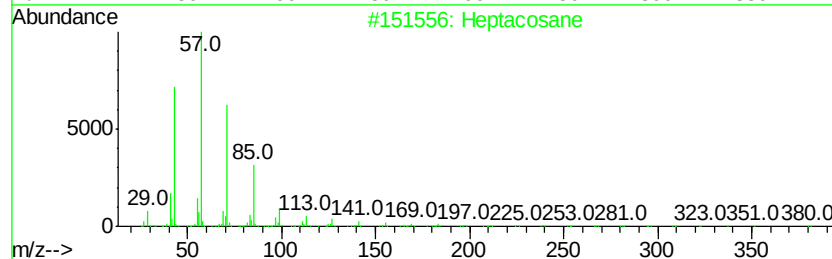
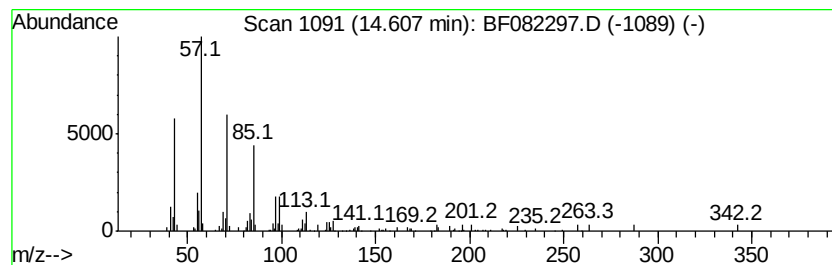
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Heptacosane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.61	2.84 ng	88294	Chrysene-d12	14.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptacosane	380	C27H56	000593-49-7	86
2		Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	86
3		Tridecane, 3-methyl-	198	C14H30	006418-41-3	86
4		Hexadecane	226	C16H34	000544-76-3	80
5		Eicosane, 10-methyl-	296	C21H44	054833-23-7	78



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082297.D
Acq On : 15 Oct 2015 9:09
Operator : UM/IZ
Sample : G4016-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

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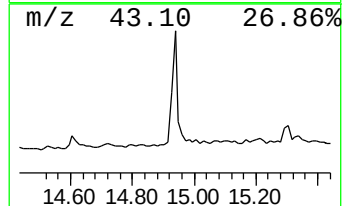
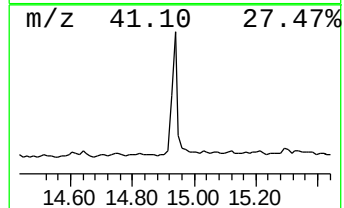
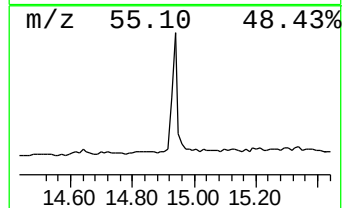
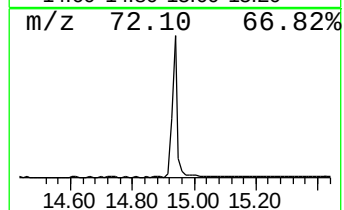
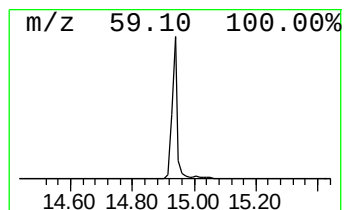
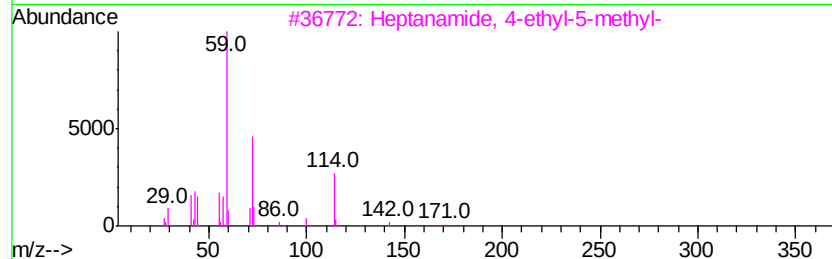
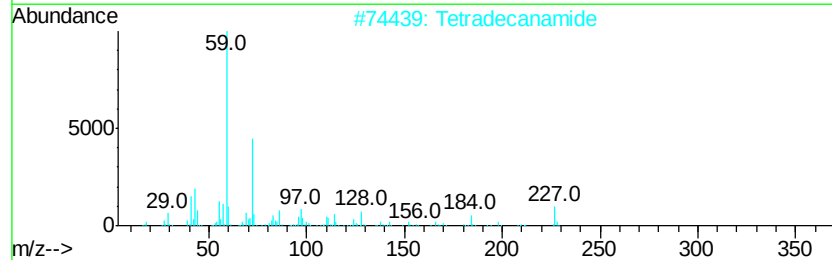
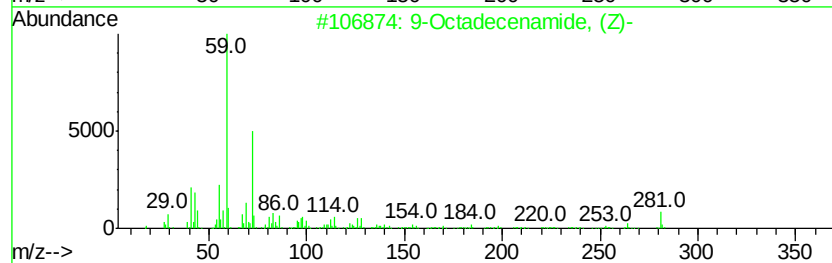
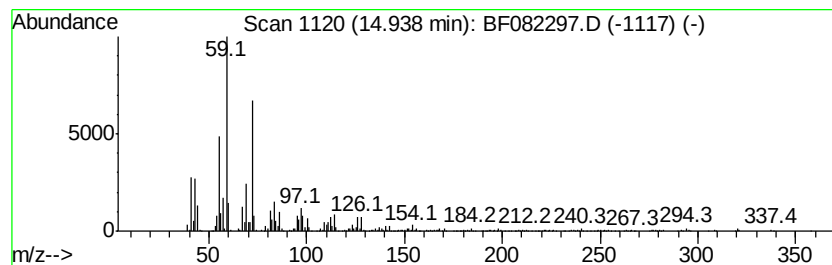
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetradecanamide Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.94	60.31 ng	1482510	Perylene-d12	15.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
2		Tetradecanamide	227	C14H29NO	000638-58-4	78
3		Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	64
4		Decanamide-	171	C10H21NO	002319-29-1	59
5		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	52



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
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Sample : G4016-04
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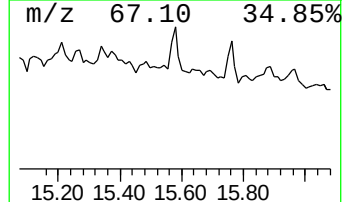
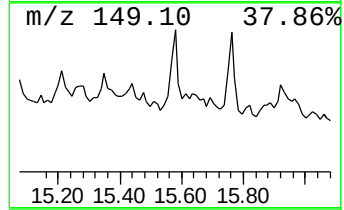
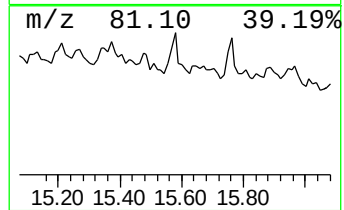
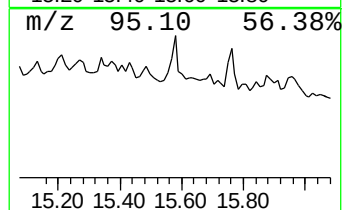
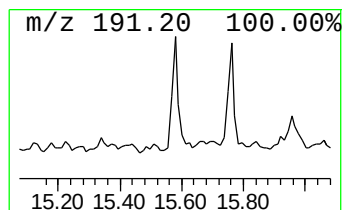
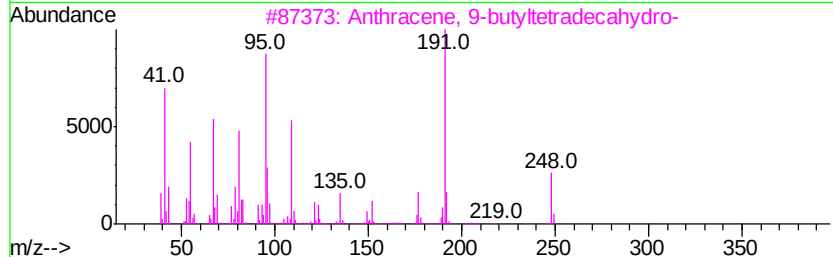
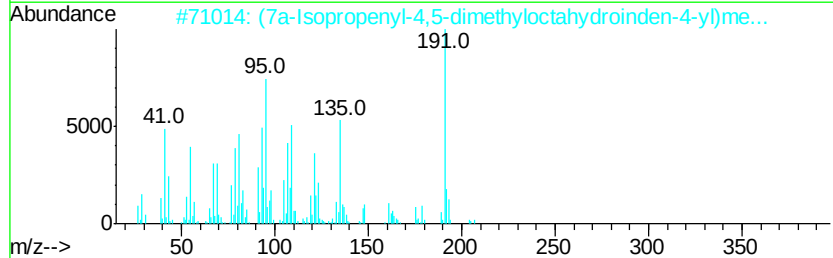
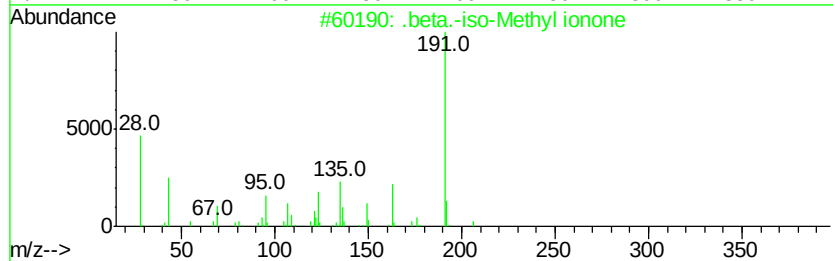
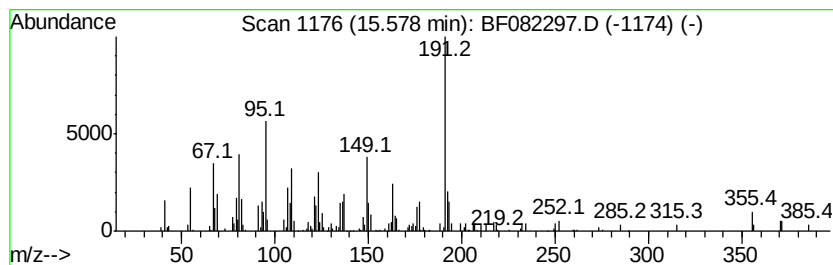
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 .beta.-iso-Methyl ionone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.58	4.71 ng	115883	Perylene-d12	15.64	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	58
2	(7a-Isopropenyl-4,5-dimethylocta...	222	C15H26O	1000187-02-9	42
3	Anthracene, 9-butyltetradecahydro-	248	C18H32	055133-89-6	38
4	Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	38
5	1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	35



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082297.D
Acq On : 15 Oct 2015 9:09
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Misc :
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ClientSampleId :
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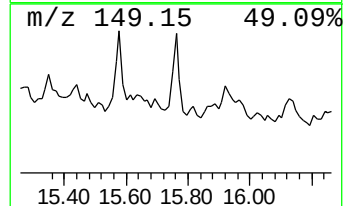
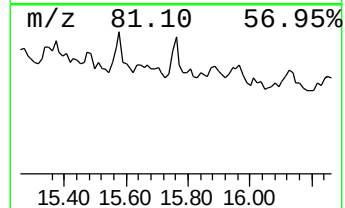
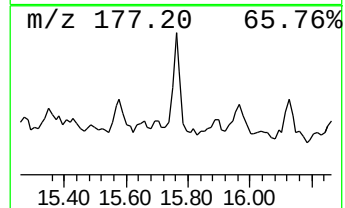
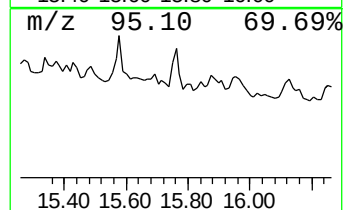
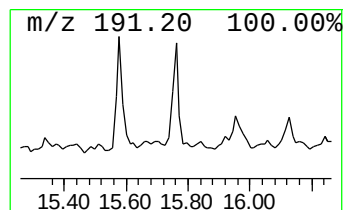
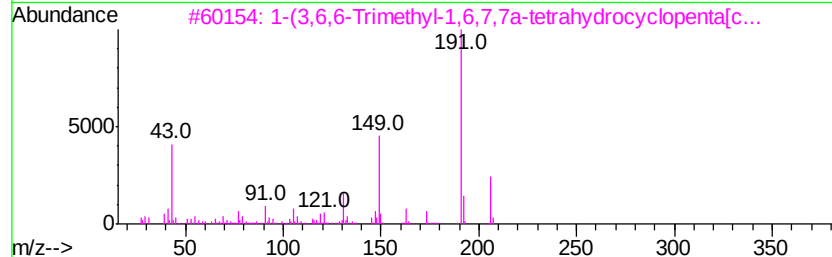
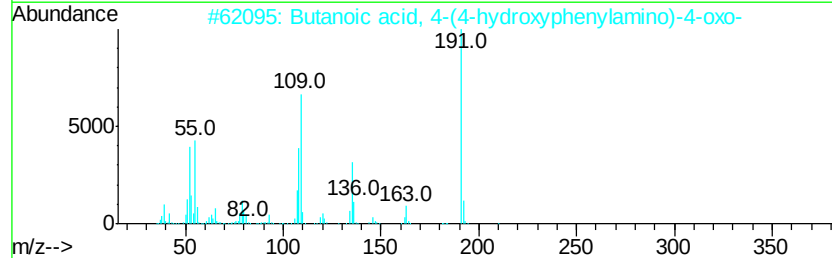
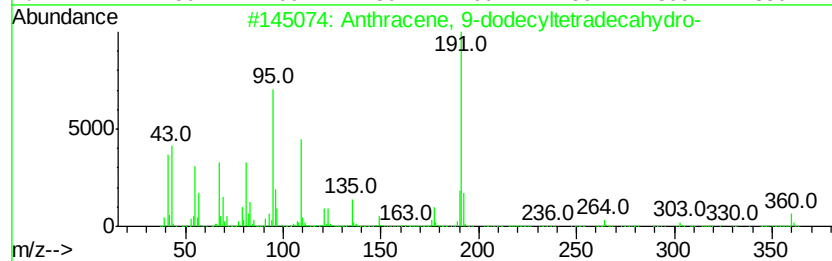
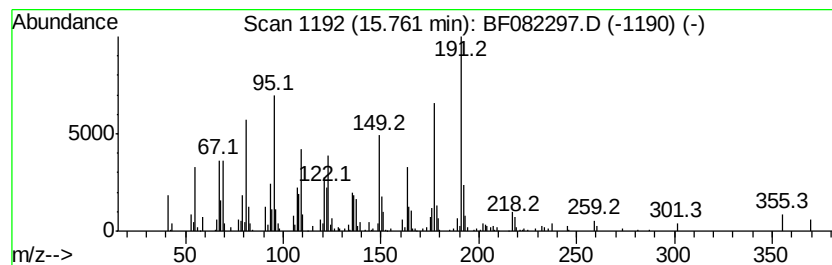
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown15.76 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.76	5.33 ng	131134	Perylene-d12	15.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	47
2		Butanoic acid, 4-(4-hydroxypheny...	209	C10H11N04	062558-67-2	27
3		1-(3,6,6-Trimethyl-1,6,7,7a-tetr...	206	C13H18O2	1000194-97-2	22
4		(7a-Isopropenyl-4,5-dimethylocta...	222	C15H26O	1000187-02-9	14
5		Silane, [4-(1,1-dimethylethyl)ph...	206	C13H22Si	018412-68-5	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
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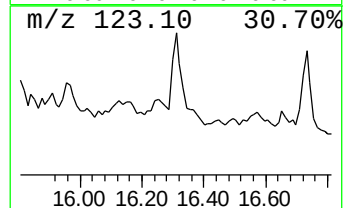
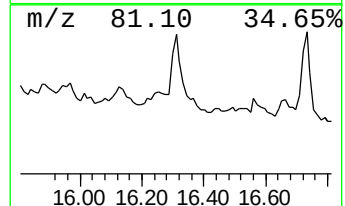
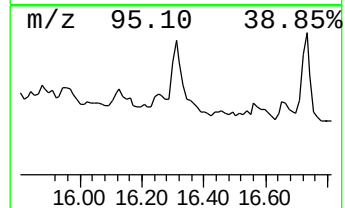
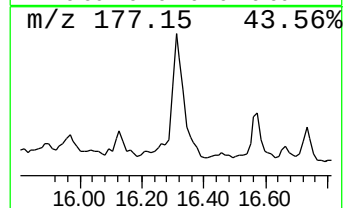
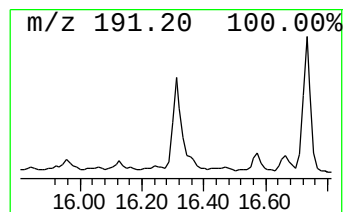
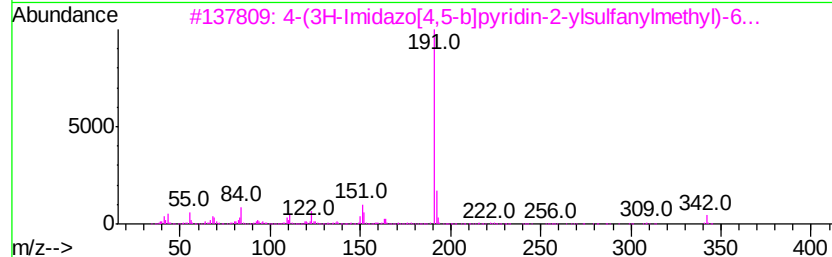
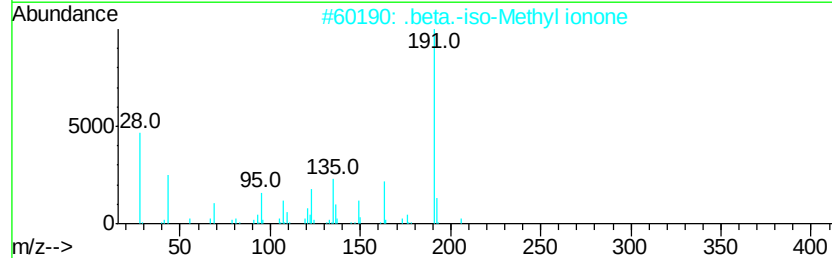
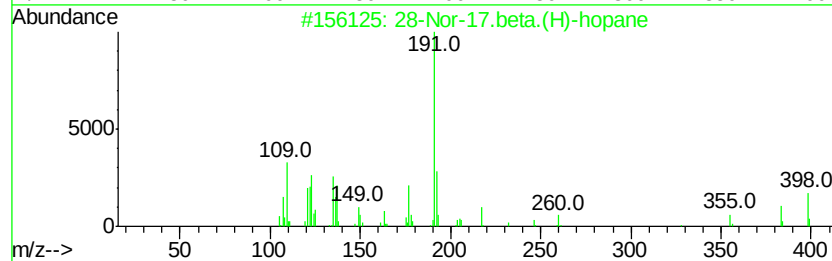
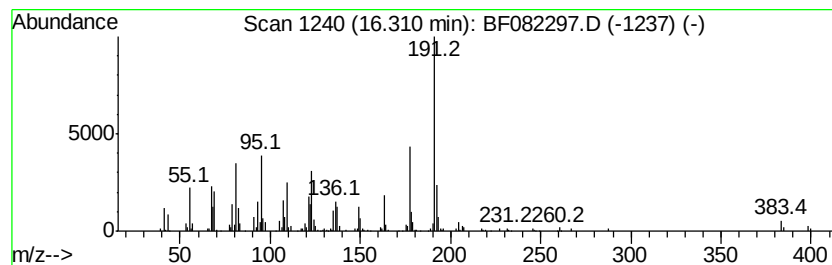
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 28-Nor-17.beta.(H)-hopane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.31	23.87 ng	586840	Perylene-d12	15.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	58
2		.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	38
3		4-(3H-Imidazo[4,5-b]pyridin-2-yl...)	342	C15H18N8S	1000276-08-2	35
4		1-Penten-3-one, 1-(2,6,6-trimeth...)	206	C14H22O	000127-43-5	35
5		2,4,7-Pteridinetriamine, 6-methyl-	191	C7H9N7	017539-50-3	27



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082297.D
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Operator : UM/IZ
Sample : G4016-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

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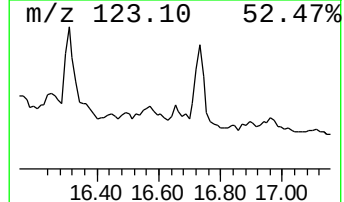
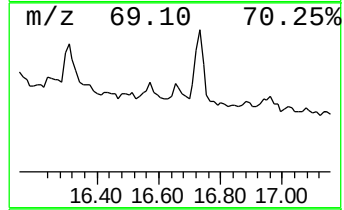
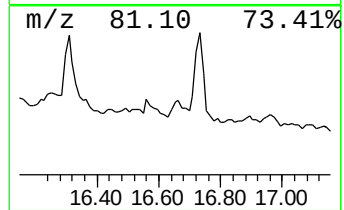
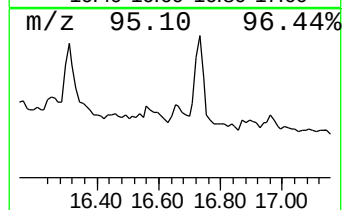
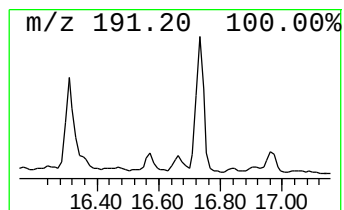
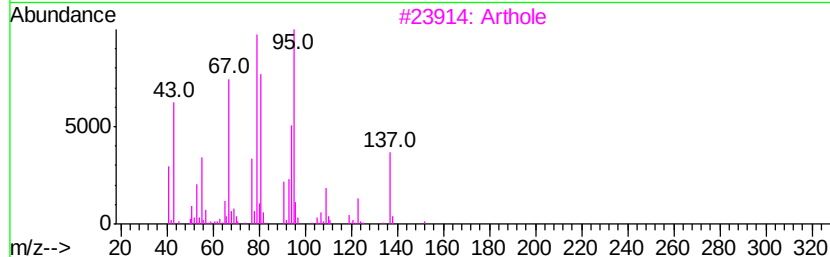
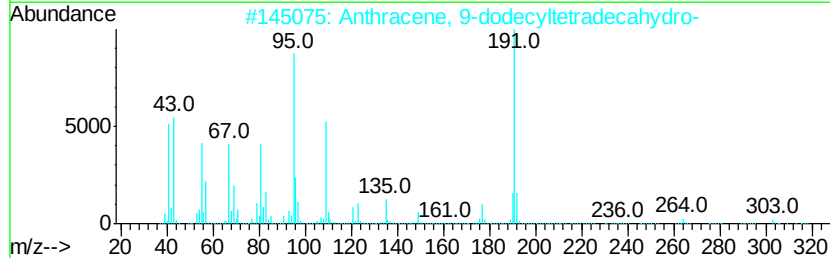
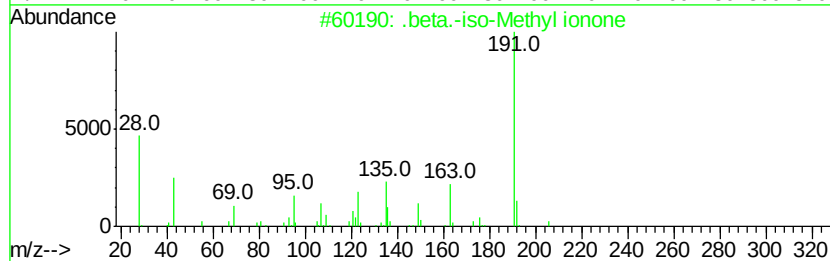
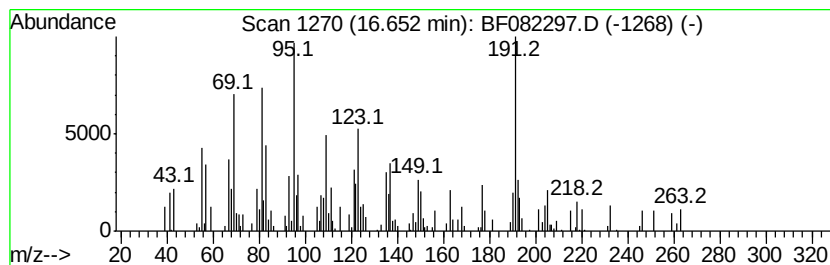
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 unknown16.65 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.65	5.37 ng	132093	Perylene-d12	15.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	46
2		Anthracene, 9-dodecyltetradeca-hy...	360	C26H48	055401-75-7	43
3		Arthole	152	C10H16O	1000281-70-0	38
4		Bicyclo[3.3.1]nonan-9-one, 1,2,4...	225	C12H19NO3	129967-65-3	38
5		Cyclopentane, 1-methyl-1-(2-meth...	138	C10H18	074764-47-9	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082297.D
Acq On : 15 Oct 2015 9:09
Operator : UM/IZ
Sample : G4016-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CS-4

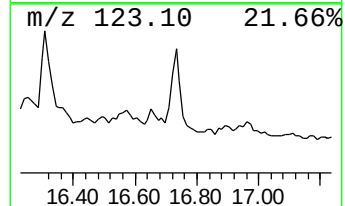
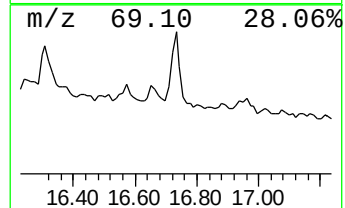
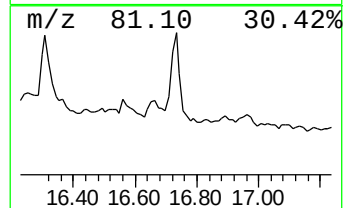
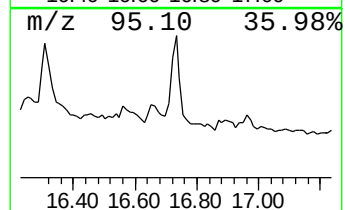
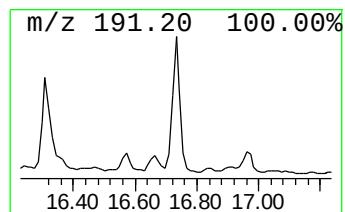
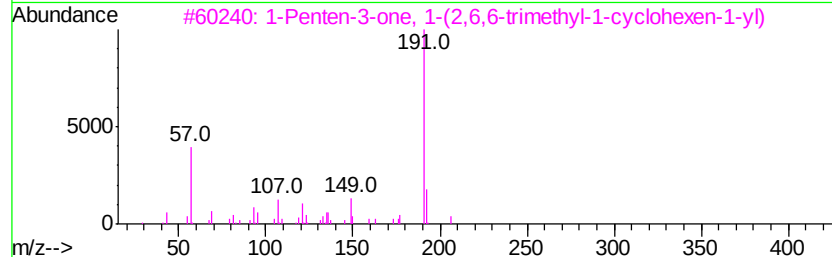
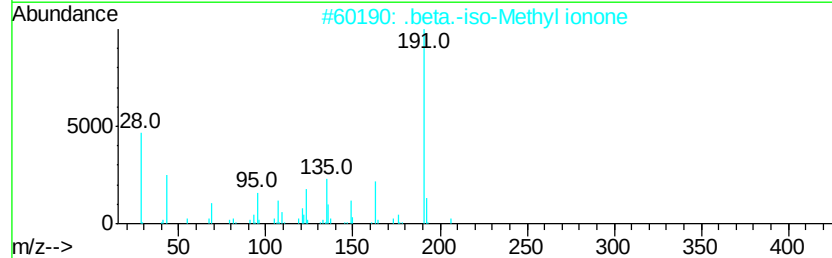
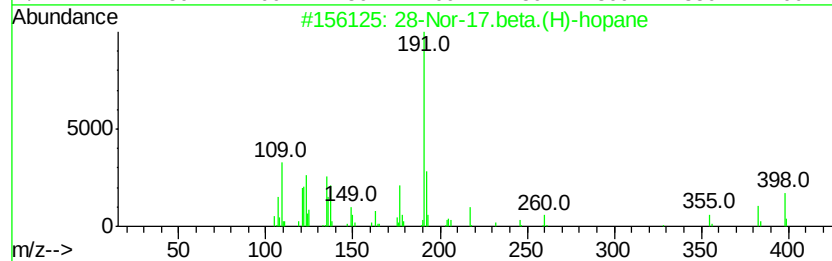
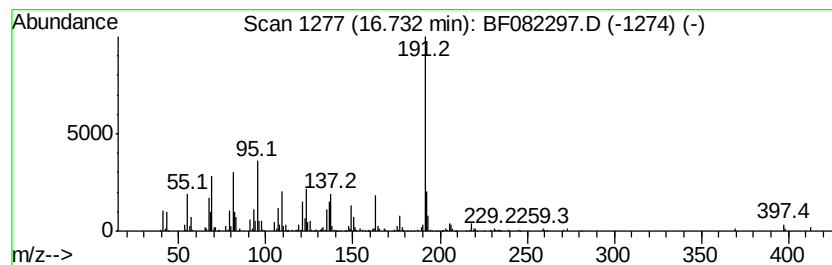
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 1-Penten-3-one, 1-(2,6,6-tr... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.73	20.81 ng	511576	Perylene-d12	15.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	28-Nor-17.beta.(H)-hopane			398	C29H50	036728-72-0	72
2	.beta.-iso-Methyl ionone			206	C14H22O	1000285-40-2	70
3	1-Penten-3-one, 1-(2,6,6-trimeth...			206	C14H22O	000127-43-5	53
4	1-Cyclohexene, 1,3,3-trimethyl-2...			206	C14H22O	1000197-08-4	42
5	2,4,5,5,8a-Pentamethyl-6,7,8,8a-...			206	C14H22O	1000195-40-9	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082297.D
Acq On : 15 Oct 2015 9:09
Operator : UM/IZ
Sample : G4016-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CS-4

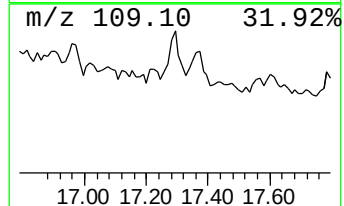
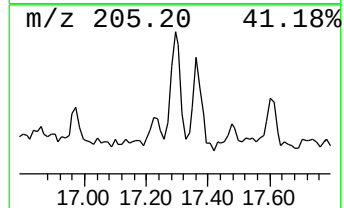
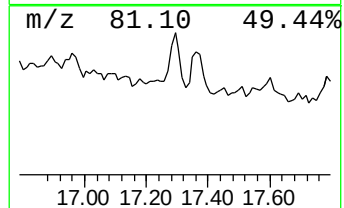
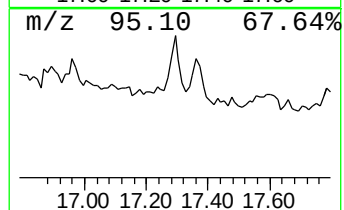
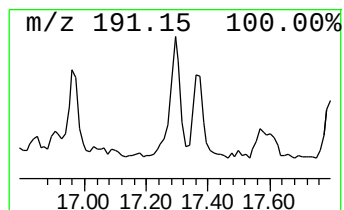
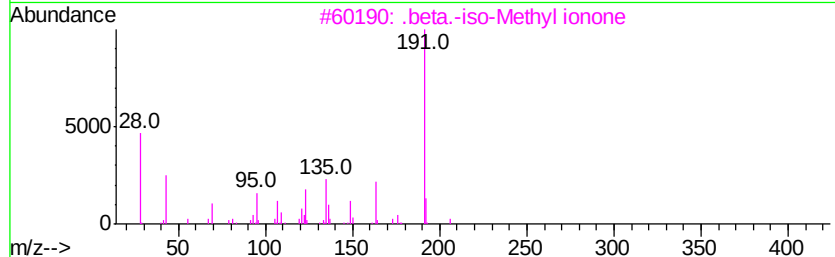
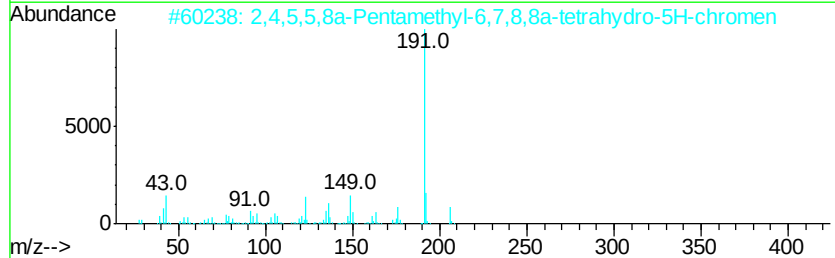
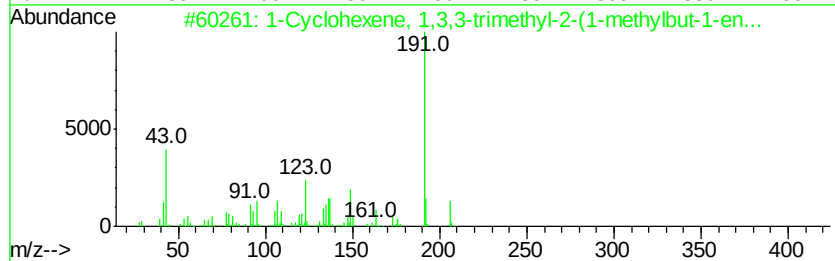
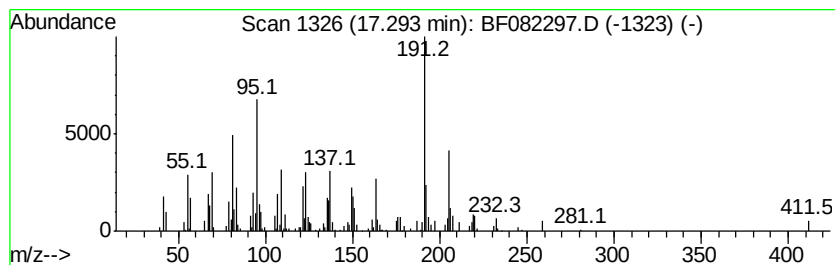
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown17.29 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.29	5.92 ng	145609	Perylene-d12	15.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	49
2		2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	38
3		.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	35
4		5-Hexen-1-one, 1-(1H-imidazol-4-...	192	C11H16N2O	069393-41-5	35
5		Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	32



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082297.D
Acq On : 15 Oct 2015 9:09
Operator : UM/IZ
Sample : G4016-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CS-4

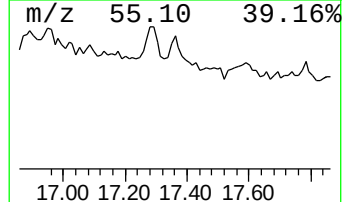
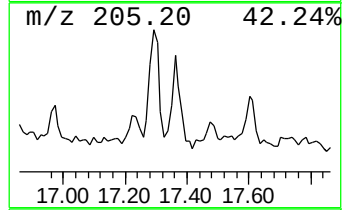
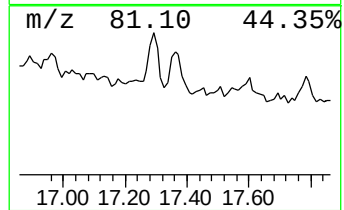
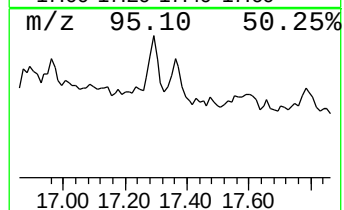
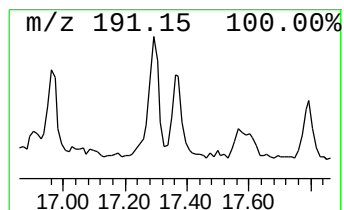
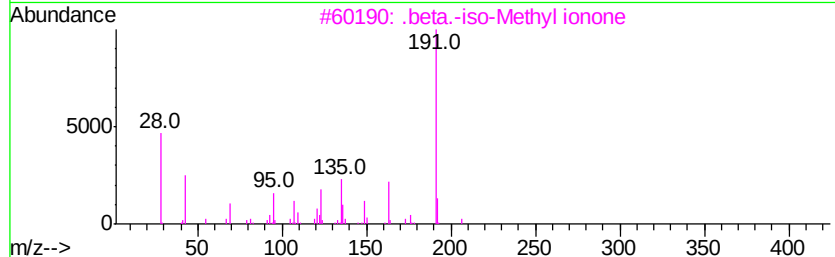
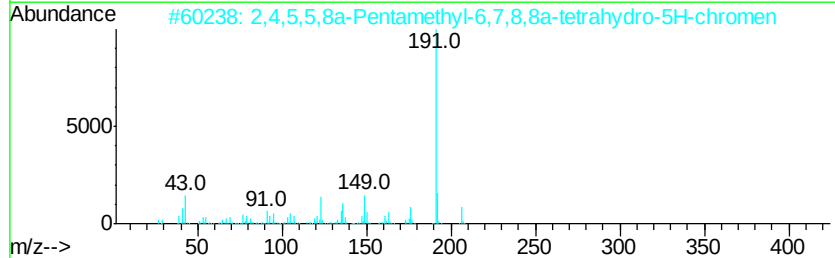
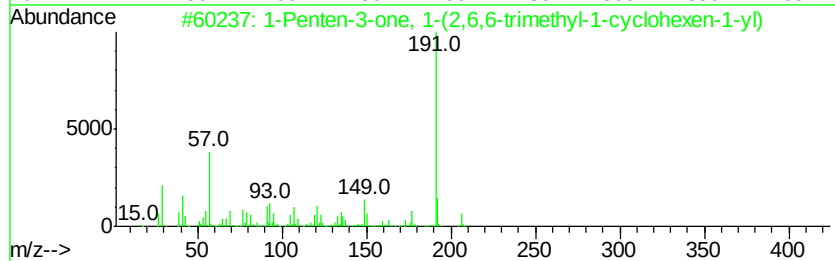
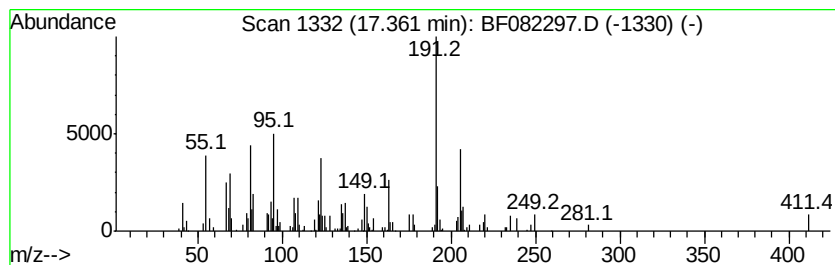
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 unknown17.36 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.36	5.51 ng	135330	Perylene-d12	15.64

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	50
2			2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	47
3			.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	43
4			2-Fluoro-4-(trifluoromethyl)acet...	206	C9H6F4O	122023-29-4	35
5			Anthracene, 9-cyclohexyltetradec...	274	C20H34	055255-70-4	32



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
 Data File : BF082297.D
 Acq On : 15 Oct 2015 9:09
 Operator : UM/IZ
 Sample : G4016-04
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CS-4

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.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
3-Penten-2-one, 4...	4.31	108.8	ng	2324860	1	6.82	427310	20.0
2-Pentanone, 4-hy...	5.04	339.6	ng	7255320	1	6.82	427310	20.0
Hexadecane	10.77	5.2	ng	189962	4	11.36	725165	20.0
5-Octadecene, (E)-	11.58	2.0	ng	73993	4	11.36	725165	20.0
Isopropyl Myristate	12.14	2.3	ng	84492	4	11.36	725165	20.0
9-Octadecenamide, ...	13.45	2.9	ng	89409	5	14.07	622609	20.0
11-Tricosene	13.90	4.2	ng	130214	5	14.07	622609	20.0
Heptacosane	14.61	2.8	ng	88294	5	14.07	622609	20.0
Tetradecanamide	14.94	60.3	ng	1482510	6	15.64	491634	20.0
.beta.-iso-Methyl...	15.58	4.7	ng	115883	6	15.64	491634	20.0
unknown15.76	15.76	5.3	ng	131134	6	15.64	491634	20.0
28-Nor-17.beta.(H...	16.31	23.9	ng	586840	6	15.64	491634	20.0
unknown16.65	16.65	5.4	ng	132093	6	15.64	491634	20.0
1-Penten-3-one, 1...	16.73	20.8	ng	511576	6	15.64	491634	20.0
unknown17.29	17.29	5.9	ng	145609	6	15.64	491634	20.0
unknown17.36	17.36	5.5	ng	135330	6	15.64	491634	20.0