

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
 Data File : BF082304.D
 Acq On : 15 Oct 2015 12:47
 Operator : UM/IZ
 Sample : G4016-09
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CS-10

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-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	4.983	247	249	252	rBV	513679	821704	50.64%	4.279%
2	5.417	285	287	289	rBV	1196515	1206353	74.34%	6.282%
3	5.817	320	322	325	rVB	20199	17895	1.10%	0.093%
4	6.457	376	378	381	rBV	1123109	1106157	68.17%	5.760%
5	6.583	387	389	392	rBV	925803	1292475	79.65%	6.730%
6	6.823	408	410	412	rBV	494883	428500	26.41%	2.231%
7	6.983	421	424	426	rBV	792096	1056369	65.10%	5.501%
8	7.395	457	460	462	rBV	784297	801935	49.42%	4.176%
9	7.486	466	468	474	rVB2	15247	25947	1.60%	0.135%
10	8.115	520	523	525	rBV	615022	578369	35.64%	3.012%
11	9.189	615	617	620	rBV	1657854	1622762	100.00%	8.450%
12	9.589	650	652	654	rBV	293271	240081	14.79%	1.250%
13	9.875	674	677	679	rBV	613534	691369	42.60%	3.600%
14	10.252	708	710	713	rBV2	41618	51281	3.16%	0.267%
15	10.298	713	714	716	rVB	34126	25909	1.60%	0.135%
16	10.515	729	733	737	rBV	11913	23246	1.43%	0.121%
17	10.663	743	746	748	rBV	1066869	1041150	64.16%	5.422%
18	10.778	754	756	760	rVB	90386	117065	7.21%	0.610%
19	10.972	771	773	775	rBV	79508	113740	7.01%	0.592%
20	11.155	786	789	792	rBV3	9503	20390	1.26%	0.106%
21	11.224	793	795	796	rVV	49687	45812	2.82%	0.239%
22	11.246	796	797	800	rVB	19321	22083	1.36%	0.115%
23	11.361	805	807	809	rBV	755919	703656	43.36%	3.664%
24	11.646	830	832	834	rBV	45633	46255	2.85%	0.241%
25	11.818	844	847	849	rBV2	14775	35786	2.21%	0.186%
26	11.898	851	854	857	rBV2	56834	123200	7.59%	0.642%
27	12.058	867	868	870	rBV	62400	54598	3.36%	0.284%
28	12.344	891	893	894	rBV	32533	47676	2.94%	0.248%
29	12.447	900	902	904	rBV	46775	70285	4.33%	0.366%
30	12.961	945	947	950	rBV	1327420	1577636	97.22%	8.215%
31	13.921	1029	1031	1034	rVB	167718	240062	14.79%	1.250%
32	14.092	1044	1046	1048	rBV	416862	509943	31.42%	2.655%
33	14.961	1120	1122	1124	rBV	266404	394135	24.29%	2.052%
34	15.670	1182	1184	1193	rVB	408664	900016	55.46%	4.687%

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Start Thrs: 0.2 Max Peaks: 100
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35	15.795	1193	1195	1198	rBV	194910	296357	18.26%	1.543%
36	16.355	1241	1244	1252	rVB	502232	1308139	80.61%	6.812%
37	16.778	1278	1281	1289	rVB	567797	1212933	74.74%	6.316%
38	17.338	1328	1330	1334	rVB	163185	332443	20.49%	1.731%

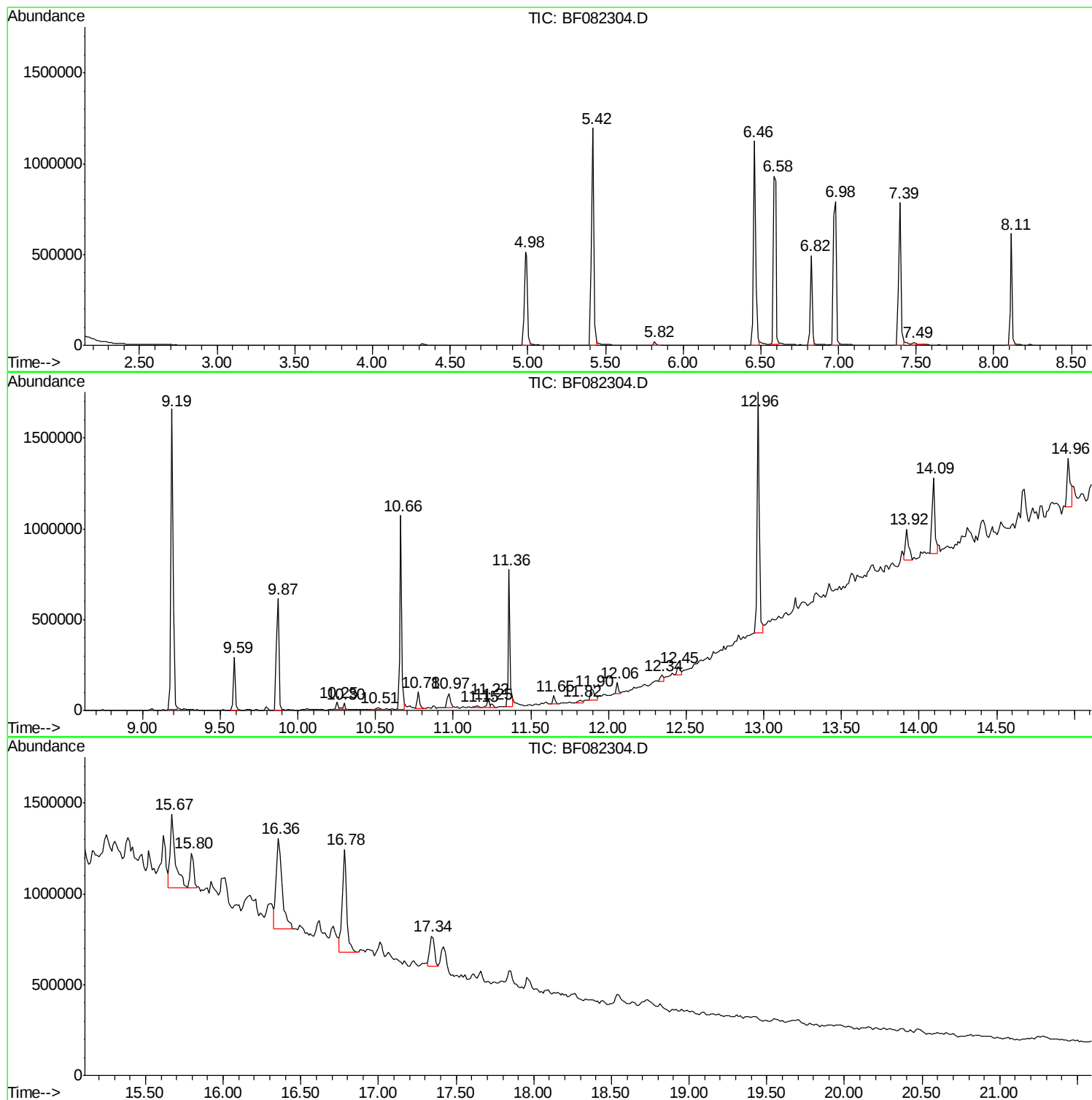
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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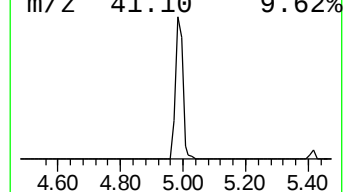
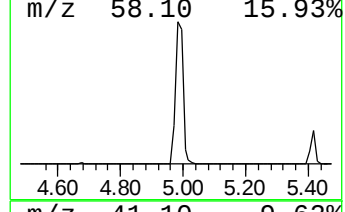
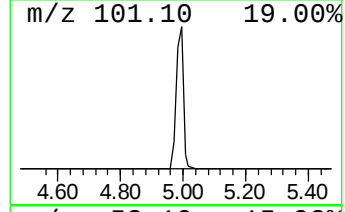
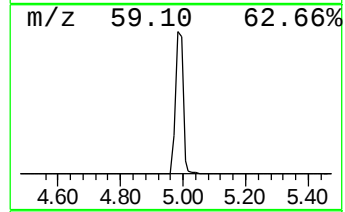
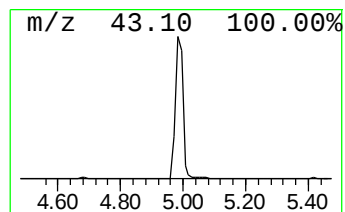
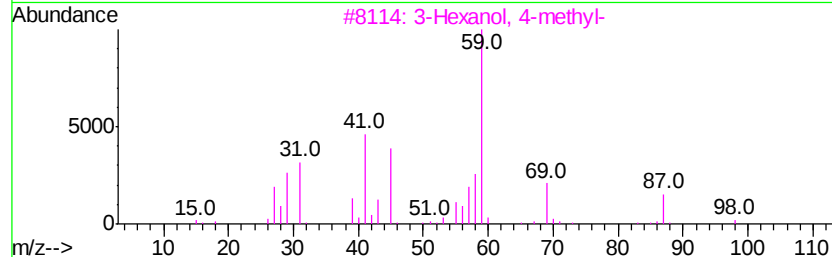
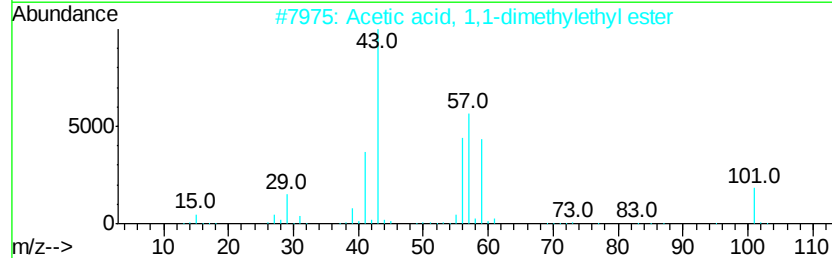
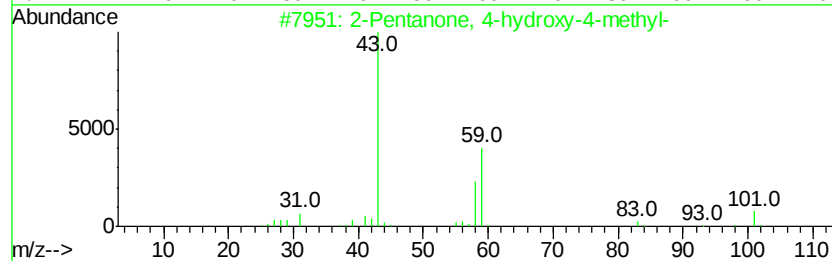
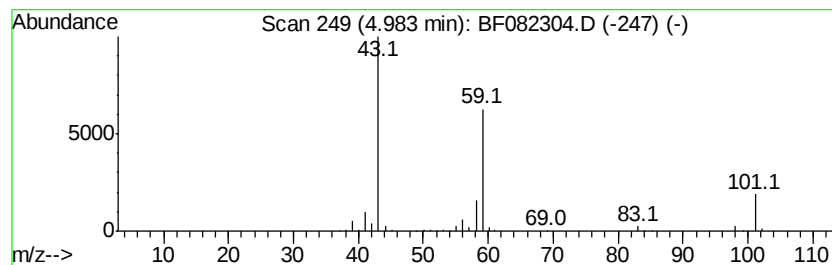
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.98	38.35 ng	821704	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32



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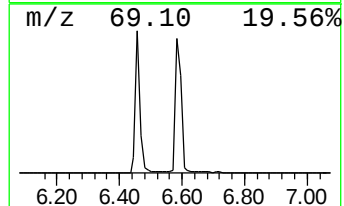
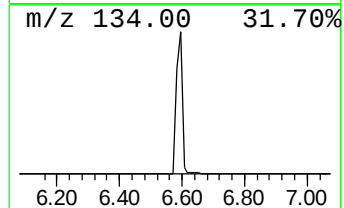
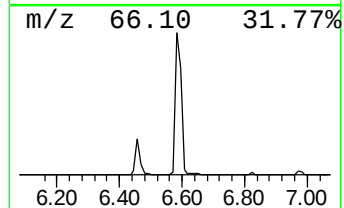
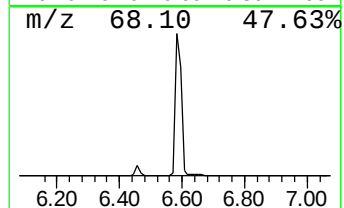
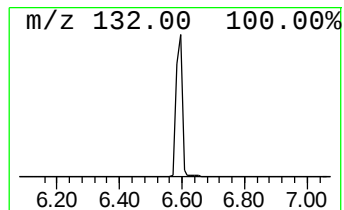
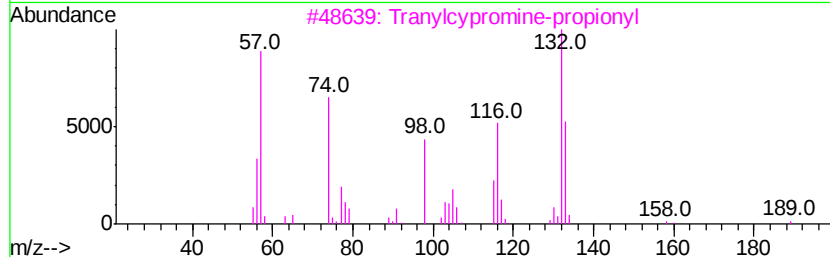
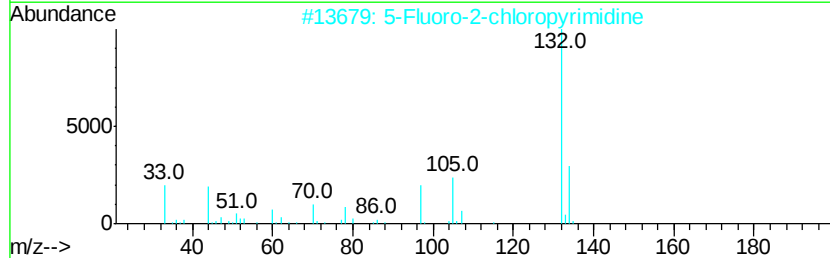
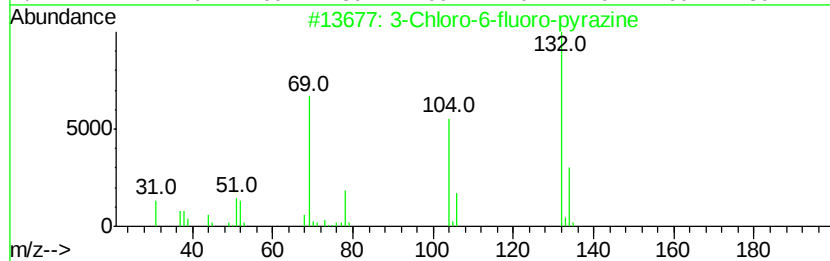
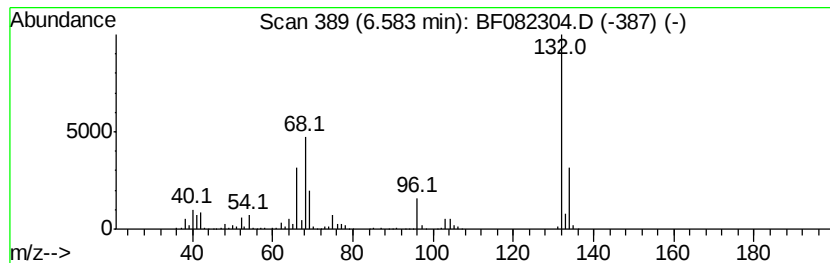
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 unknown6.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	60.33 ng	1292480	1,4-Dichlorobenzene-d4	6.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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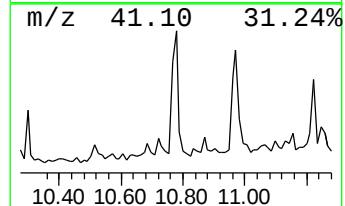
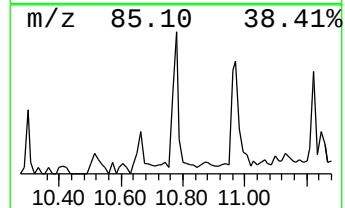
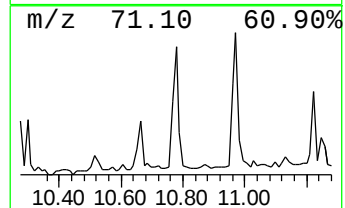
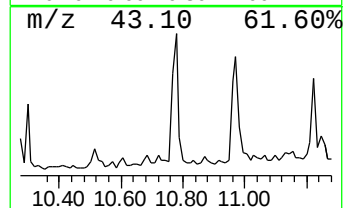
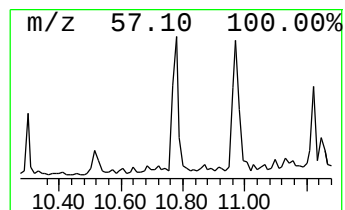
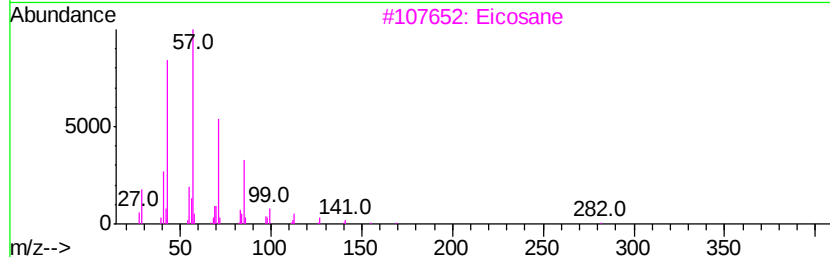
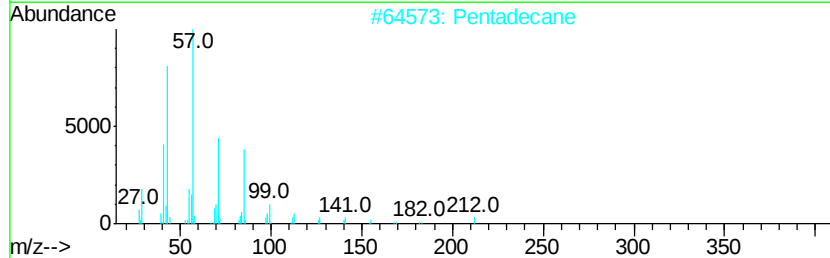
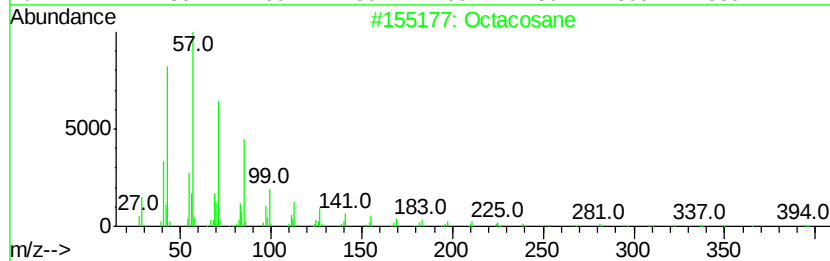
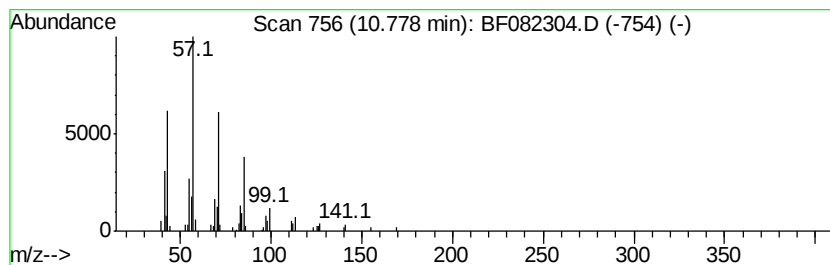
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 4 Octacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.78	3.33 ng	117065	Phenanthrene-d10	11.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octacosane	394	C28H58	000630-02-4	91
2		Pentadecane	212	C15H32	000629-62-9	91
3		Eicosane	282	C20H42	000112-95-8	91
4		Tetradecane	198	C14H30	000629-59-4	91
5		Tetratriacontane	479	C34H70	014167-59-0	90



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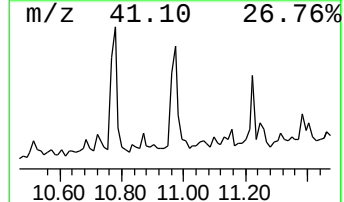
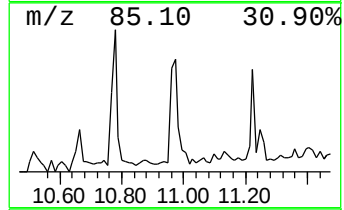
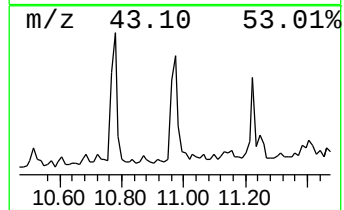
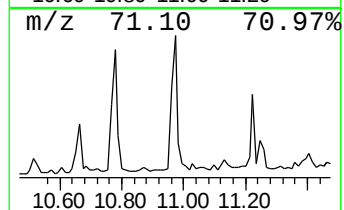
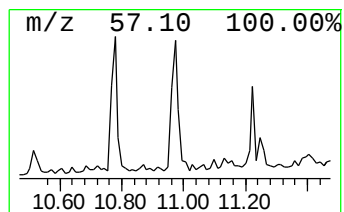
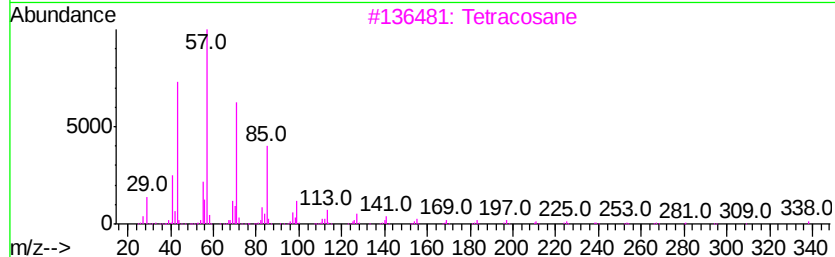
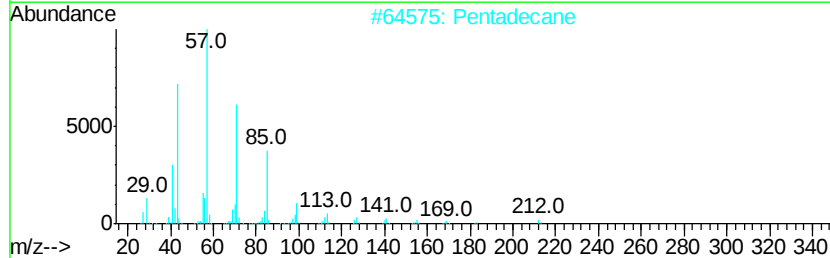
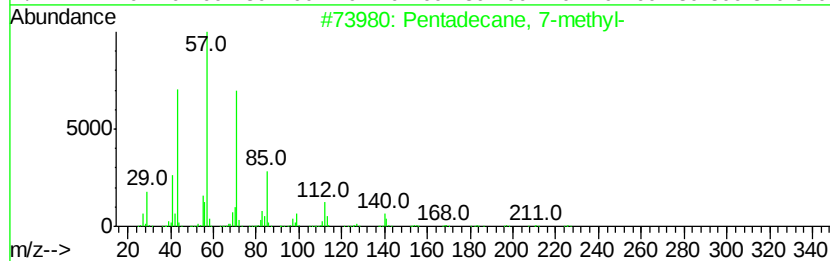
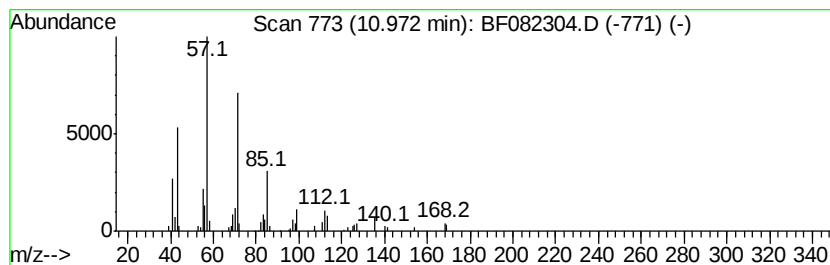
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Pentadecane, 7-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.97	3.23 ng	113740	Phenanthrene-d10	11.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 7-methyl-	226	C16H34	006165-40-8	80	
2	Pentadecane	212	C15H32	000629-62-9	80	
3	Tetracosane	338	C24H50	000646-31-1	72	
4	Tetradecane	198	C14H30	000629-59-4	72	
5	Heptacosane	380	C27H56	000593-49-7	72	



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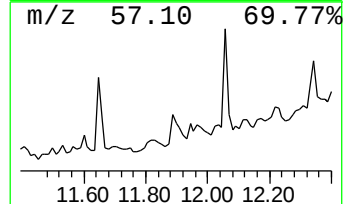
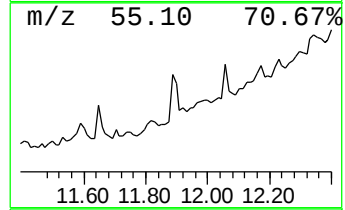
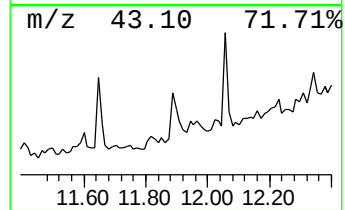
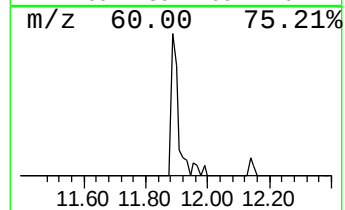
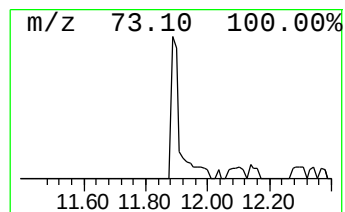
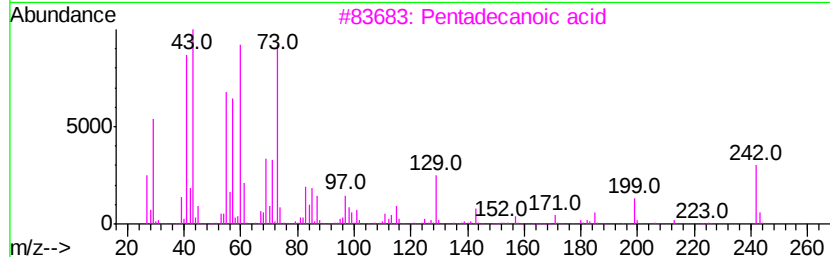
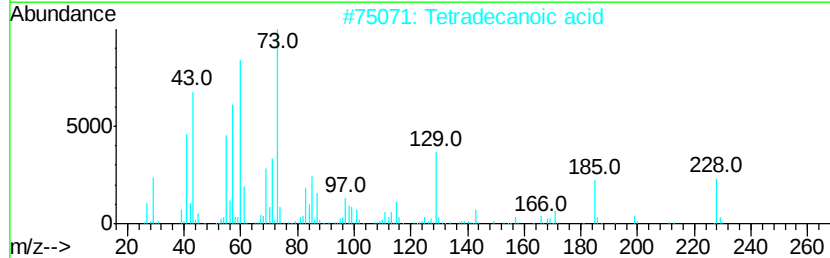
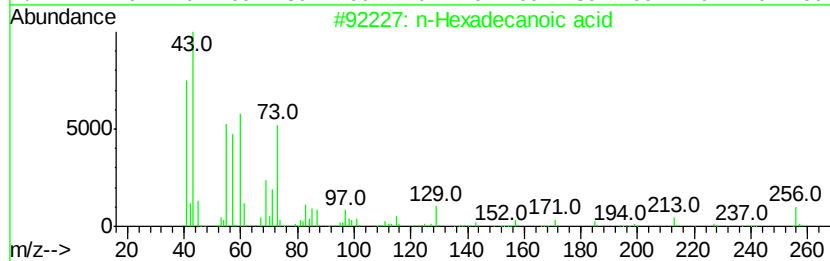
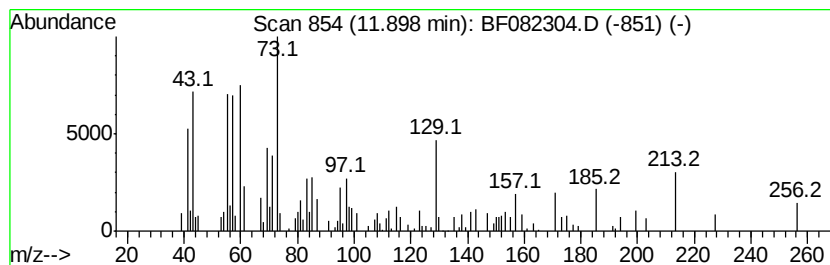
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Peak Number 6 n-Hexadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	3.50 ng	123200	Phenanthrene-d10	11.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	58
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	50
4	Tridecanoic acid	214	C13H26O2	000638-53-9	49
5	Dodecanoic acid	200	C12H24O2	000143-07-7	47



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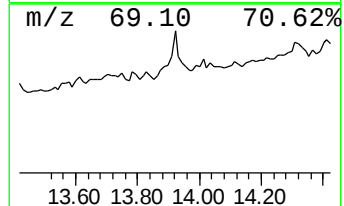
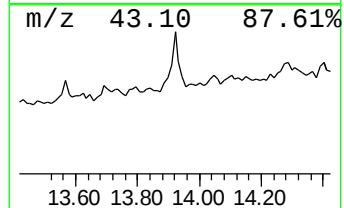
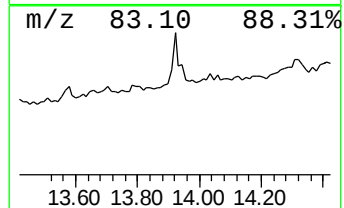
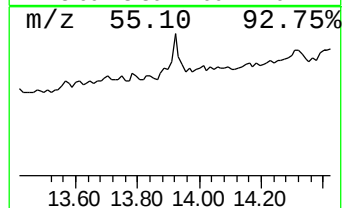
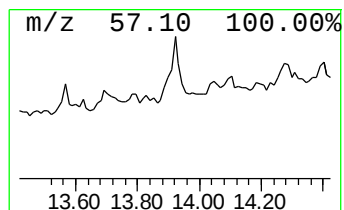
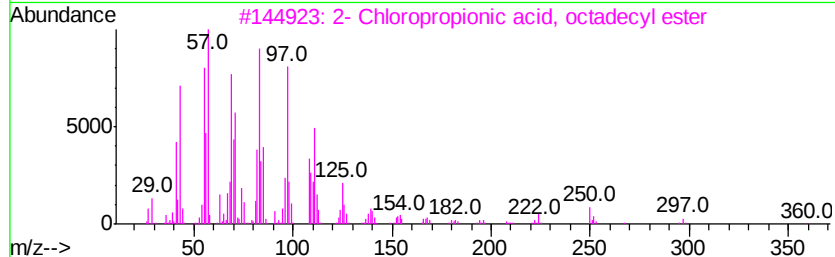
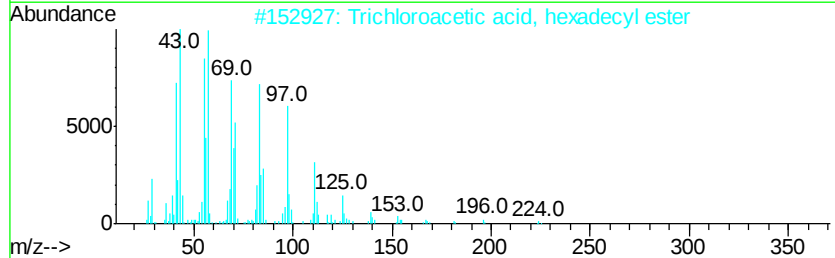
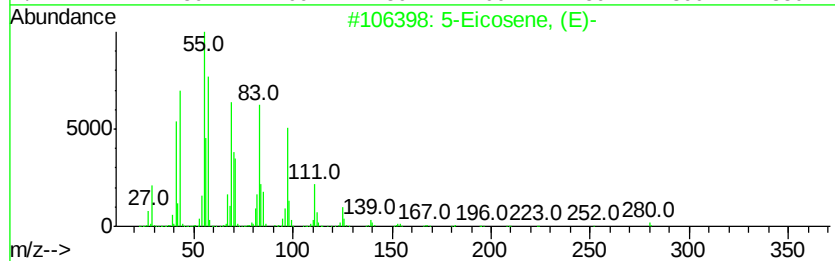
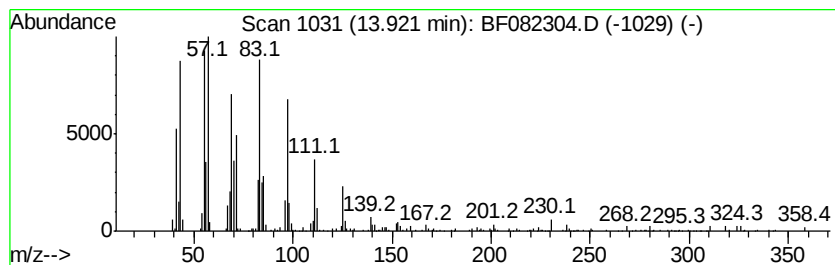
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF101015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 5-Eicosene, (E)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.92	9.42 ng	240062	Chrysene-d12	14.09
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	5-Eicosene, (E)-	280 C20H40	074685-30-6	95
2	Trichloroacetic acid, hexadecyl ...	386 C18H33Cl3O2	074339-54-1	91
3	2- Chloropropionic acid, octadec...	360 C21H41ClO2	088104-31-8	91
4	Cyclotetracosane	336 C24H48	000297-03-0	91
5	Dichloroacetic acid, tetradecyl ...	324 C16H30Cl2O2	083005-02-1	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF101615\
Data File : BF082304.D
Acq On : 15 Oct 2015 12:47
Operator : UM/IZ
Sample : G4016-09
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CS-10

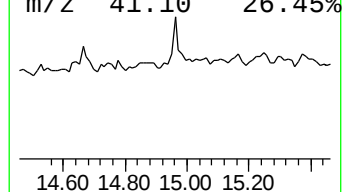
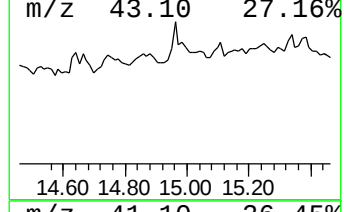
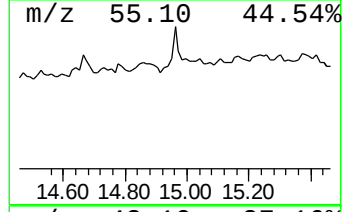
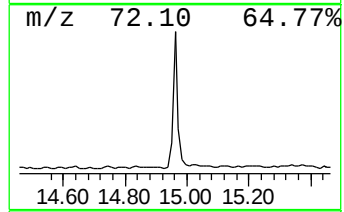
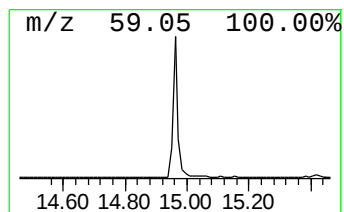
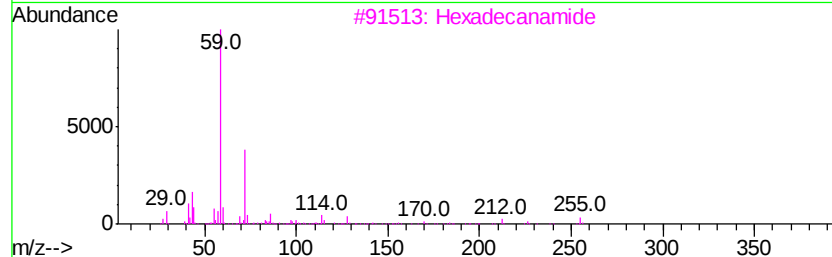
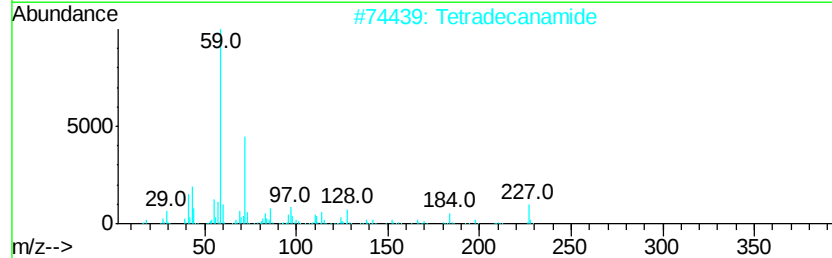
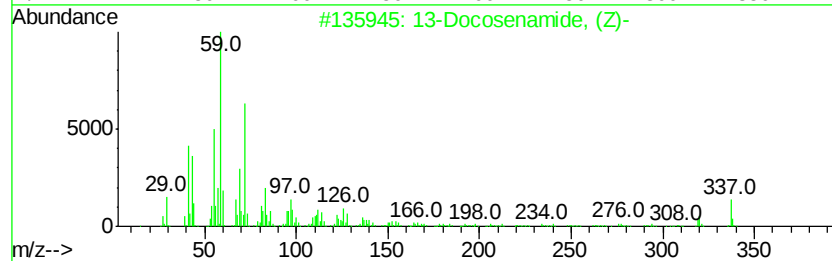
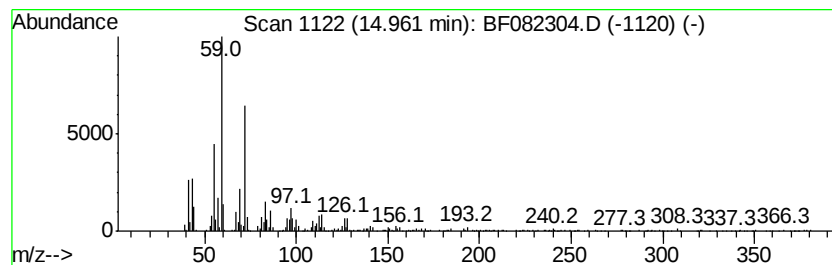
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 13-Docosenamide, (Z)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	8.76 ng	394135	Perylene-d12	15.67

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	87
2		Tetradecanamide	227	C14H29NO	000638-58-4	78
3		Hexadecanamide	255	C16H33NO	000629-54-9	78
4		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	68
5		Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FN03	061338-98-5	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082304.D
Acq On : 15 Oct 2015 12:47
Operator : UM/IZ
Sample : G4016-09
Misc :
ALS Vial : 29 Sample Multiplier: 1

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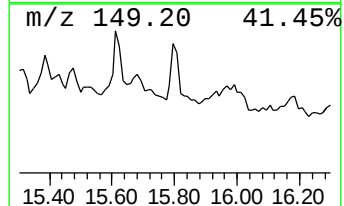
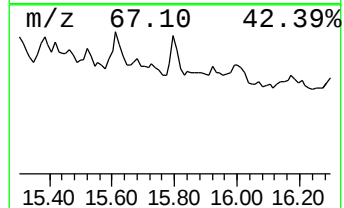
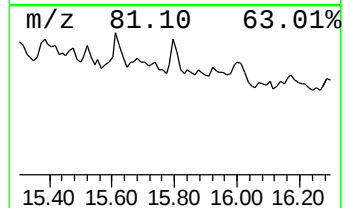
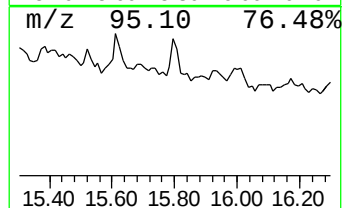
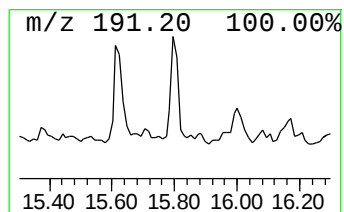
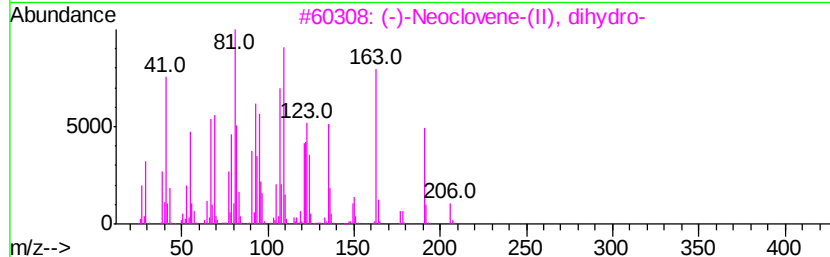
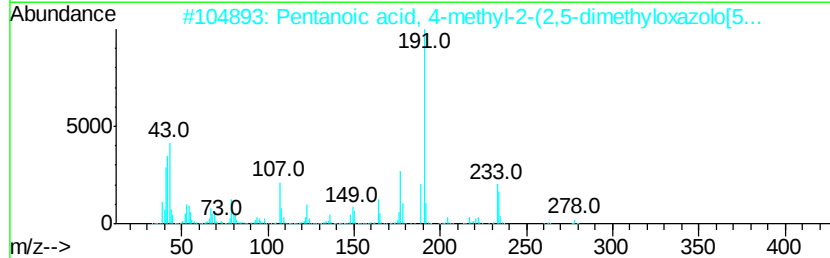
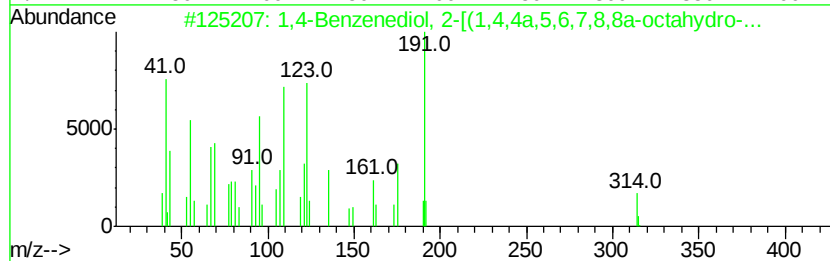
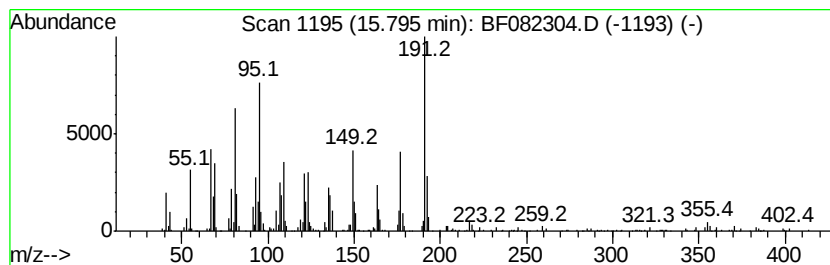
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown15.80 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.80	6.59 ng	296357	Perylene-d12	15.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4-Benzenediol, 2-[(1,4,4a,5,6,...	314	C21H30O2	039707-55-6	32
2		Pentanoic acid, 4-methyl-2-(2,5-...	278	C13H18N4O3	1000294-63-5	16
3		(-)-Neoclovene-(II), dihydro-	206	C15H26	1000152-83-6	14
4		2H-Pyran-2,2-dicarboxylic acid, ...	242	C12H18O5	024588-59-8	14
5		5-(1-Isopropenyl-4,5-dimethylbic...	332	C22H36O2	1000195-69-8	14



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082304.D
Acq On : 15 Oct 2015 12:47
Operator : UM/IZ
Sample : G4016-09
Misc :
ALS Vial : 29 Sample Multiplier: 1

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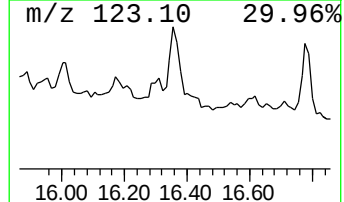
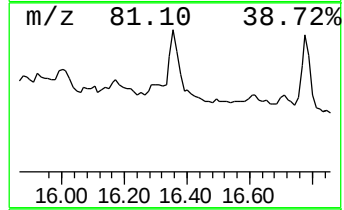
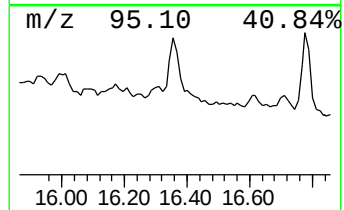
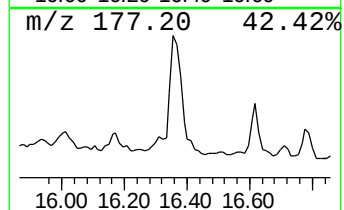
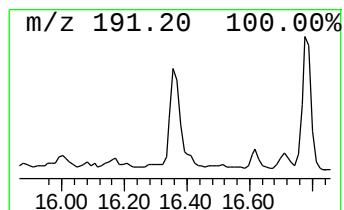
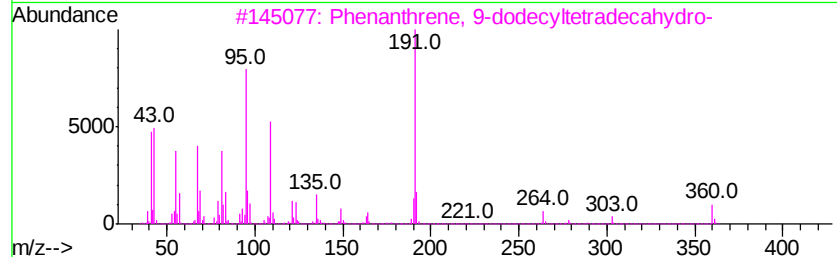
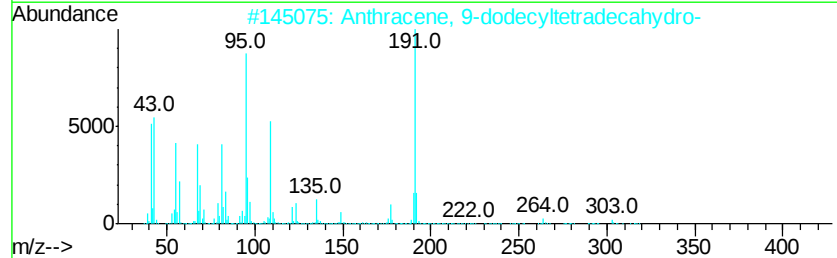
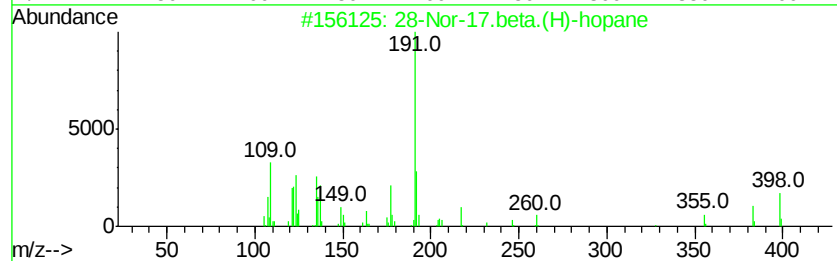
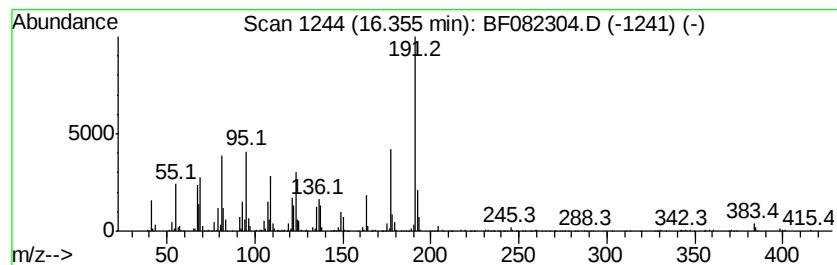
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 28-Nor-17.beta.(H)-hopane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.36	29.07 ng	1308140	Perylene-d12	15.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	62
2		Anthracene, 9-dodecyltetradeca...	360	C26H48	055401-75-7	40
3		Phenanthrene, 9-dodecyltetradeca...	360	C26H48	055334-01-5	40
4		.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	37
5		9H-Fluorene-9-carbonitrile	191	C14H9N	001529-40-4	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
Data File : BF082304.D
Acq On : 15 Oct 2015 12:47
Operator : UM/IZ
Sample : G4016-09
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CS-10

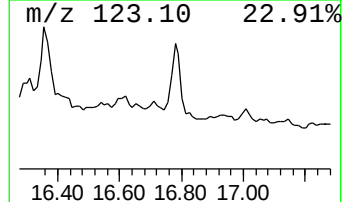
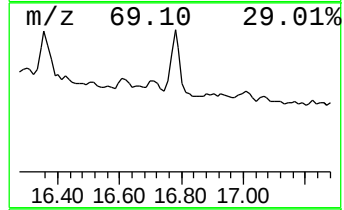
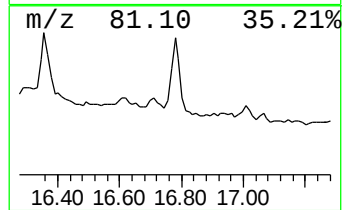
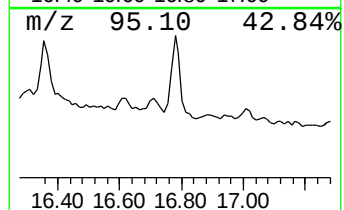
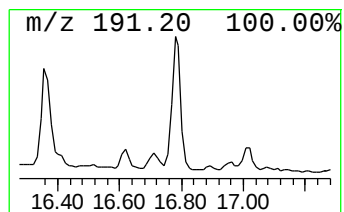
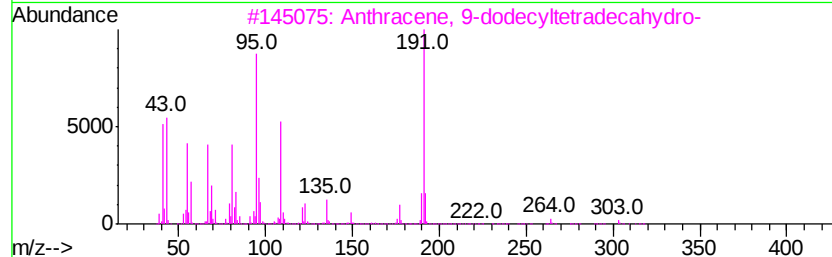
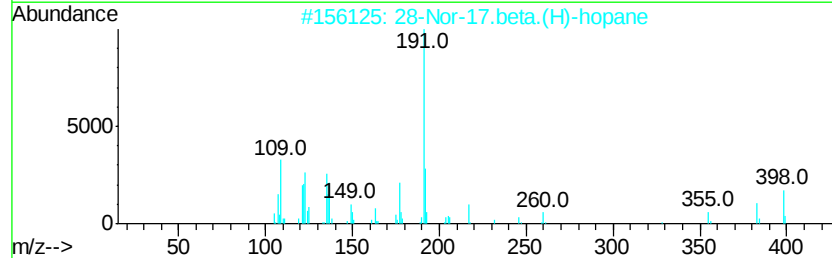
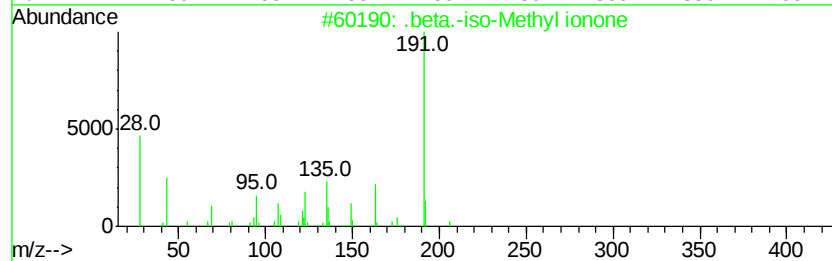
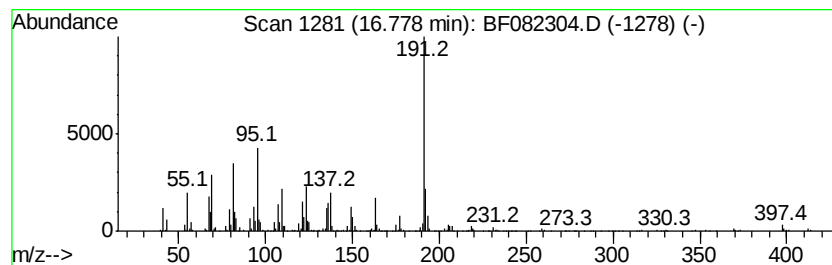
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 .beta.-iso-Methyl ionone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.78	26.95 ng	1212930	Perylene-d12	15.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	60
2		28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	50
3		Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	50
4		1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	40
5		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
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Acq On : 15 Oct 2015 12:47
Operator : UM/IZ
Sample : G4016-09
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CS-10

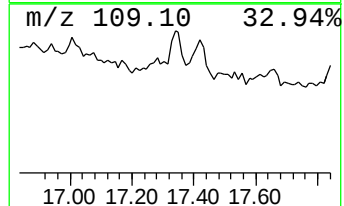
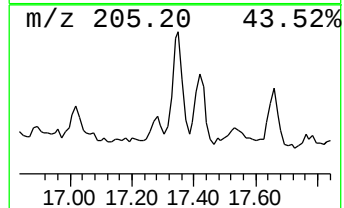
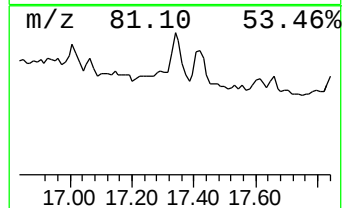
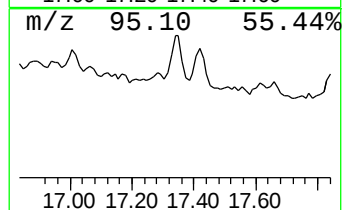
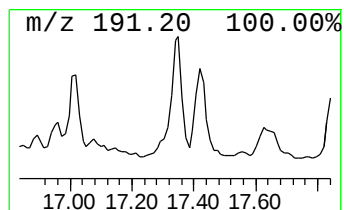
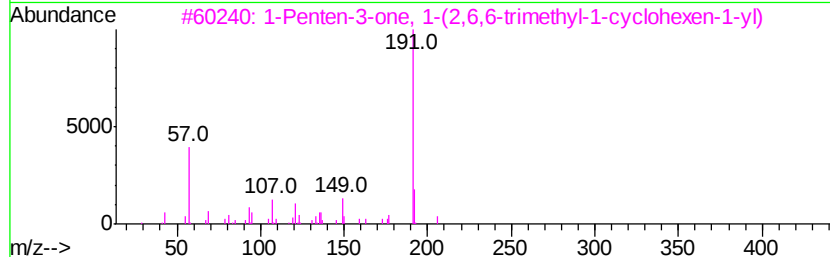
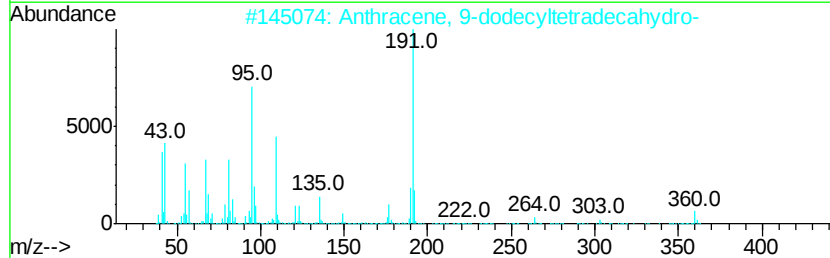
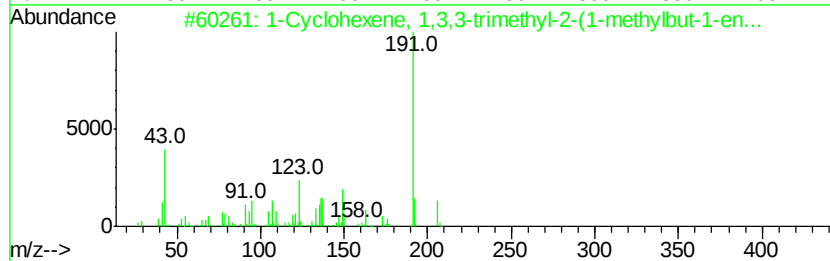
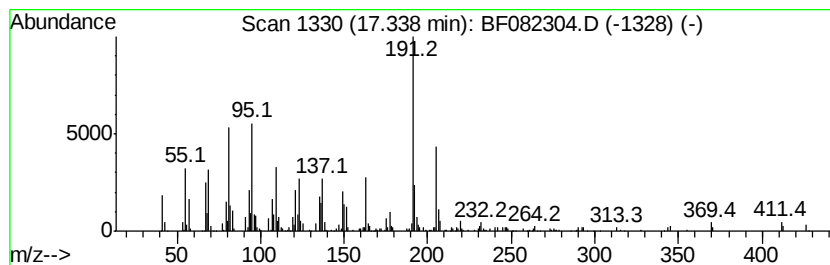
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown17.34 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.34	7.39 ng	332443	Perylene-d12	15.67

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	47
2		Anthracene, 9-dodecyltetradecahy...	360	C26H48	055401-75-7	38
3		1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	35
4		2-(2-((2-Chloro-6-fluorobenzyl)s...	334	C17H16ClFN2S	1000298-18-2	27
5		2'-Fluoro-6'-(trifluoromethyl)-a...	206	C9H6F4O	174013-29-7	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF101615\
 Data File : BF082304.D
 Acq On : 15 Oct 2015 12:47
 Operator : UM/IZ
 Sample : G4016-09
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
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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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown6.58	6.58	60.3	ng	1292480	1	6.82	428500	20.0
Octacosane	10.78	3.3	ng	117065	4	11.36	703656	20.0
Pentadecane, 7-me...	10.97	3.2	ng	113740	4	11.36	703656	20.0
n-Hexadecanoic acid	11.90	3.5	ng	123200	4	11.36	703656	20.0
5-Eicosene, (E)-	13.92	9.4	ng	240062	5	14.09	509943	20.0
13-Docosenamide, ...	14.96	8.8	ng	394135	6	15.67	900016	20.0
unknown15.80	15.80	6.6	ng	296357	6	15.67	900016	20.0
28-Nor-17.beta.(H...	16.36	29.1	ng	1308140	6	15.67	900016	20.0
.beta.-iso-Methyl...	16.78	26.9	ng	1212930	6	15.67	900016	20.0
unknown17.34	17.34	7.4	ng	332443	6	15.67	900016	20.0