

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
 Data File : BF083623.D
 Acq On : 9 Dec 2015 5:38
 Operator : UM/IZ
 Sample : G4647-03
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BRIDGE14-2(0-2)

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-BF120815.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	3.539	160	162	182	rBV	432694	989538	24.91%	3.028%
2	3.984	198	201	220	rBV	2271347	3923937	98.76%	12.007%
3	5.276	311	314	322	rBV	2161731	3925359	98.80%	12.011%
4	5.402	322	325	338	rVB	2387117	3973234	100.00%	12.157%
5	5.699	348	351	362	rVB	649045	865133	21.77%	2.647%
6	5.916	367	370	378	rBV	2142808	2836100	71.38%	8.678%
7	6.522	421	423	438	rBV	1477677	2437538	61.35%	7.458%
8	7.596	514	517	528	rBV	808490	1047336	26.36%	3.205%
9	9.151	651	653	657	rVB	29529	40269	1.01%	0.123%
10	9.333	666	669	679	rBV	2075937	3193832	80.38%	9.773%
11	9.551	684	688	692	rBV	28920	39800	1.00%	0.122%
12	10.031	727	730	735	rBV	323945	517883	13.03%	1.585%
13	10.385	758	761	767	rBV2	699229	1072668	27.00%	3.282%
14	10.785	790	796	799	rBV2	18984	47156	1.19%	0.144%
15	11.151	824	828	832	rBV4	16144	41763	1.05%	0.128%
16	11.219	832	834	837	rVV	120616	178521	4.49%	0.546%
17	11.585	862	866	868	rBV	36784	81615	2.05%	0.250%
18	11.676	871	874	881	rBV	1015636	1618371	40.73%	4.952%
19	11.996	899	902	906	rBV	135496	216314	5.44%	0.662%
20	12.728	964	966	968	rBV	61364	70441	1.77%	0.216%
21	12.774	968	970	977	rVB	490270	785266	19.76%	2.403%
22	13.620	1038	1044	1046	rBV2	22832	45903	1.16%	0.140%
23	13.825	1059	1062	1067	rBV2	128000	230238	5.79%	0.704%
24	14.202	1092	1095	1099	rVB	25439	54027	1.36%	0.165%
25	14.545	1123	1125	1129	rVB	51763	82074	2.07%	0.251%
26	15.105	1172	1174	1177	rBV2	102886	173044	4.36%	0.529%
27	15.403	1197	1200	1203	rBV	1445451	1944885	48.95%	5.951%
28	15.643	1219	1221	1223	rBV2	32358	57153	1.44%	0.175%
29	16.088	1258	1260	1263	rBV	61847	92838	2.34%	0.284%
30	16.923	1330	1333	1337	rBV2	193849	411226	10.35%	1.258%
31	16.991	1337	1339	1343	rBV	651838	890700	22.42%	2.725%
32	18.637	1482	1483	1486	rVB	349670	431394	10.86%	1.320%
33	19.529	1559	1561	1565	rVB	230655	365837	9.21%	1.119%

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Misc :
ALS Vial : 31 Sample Multiplier: 1

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ClientSampleId :
BRIDGE14-2(0-2)

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-BF120815.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

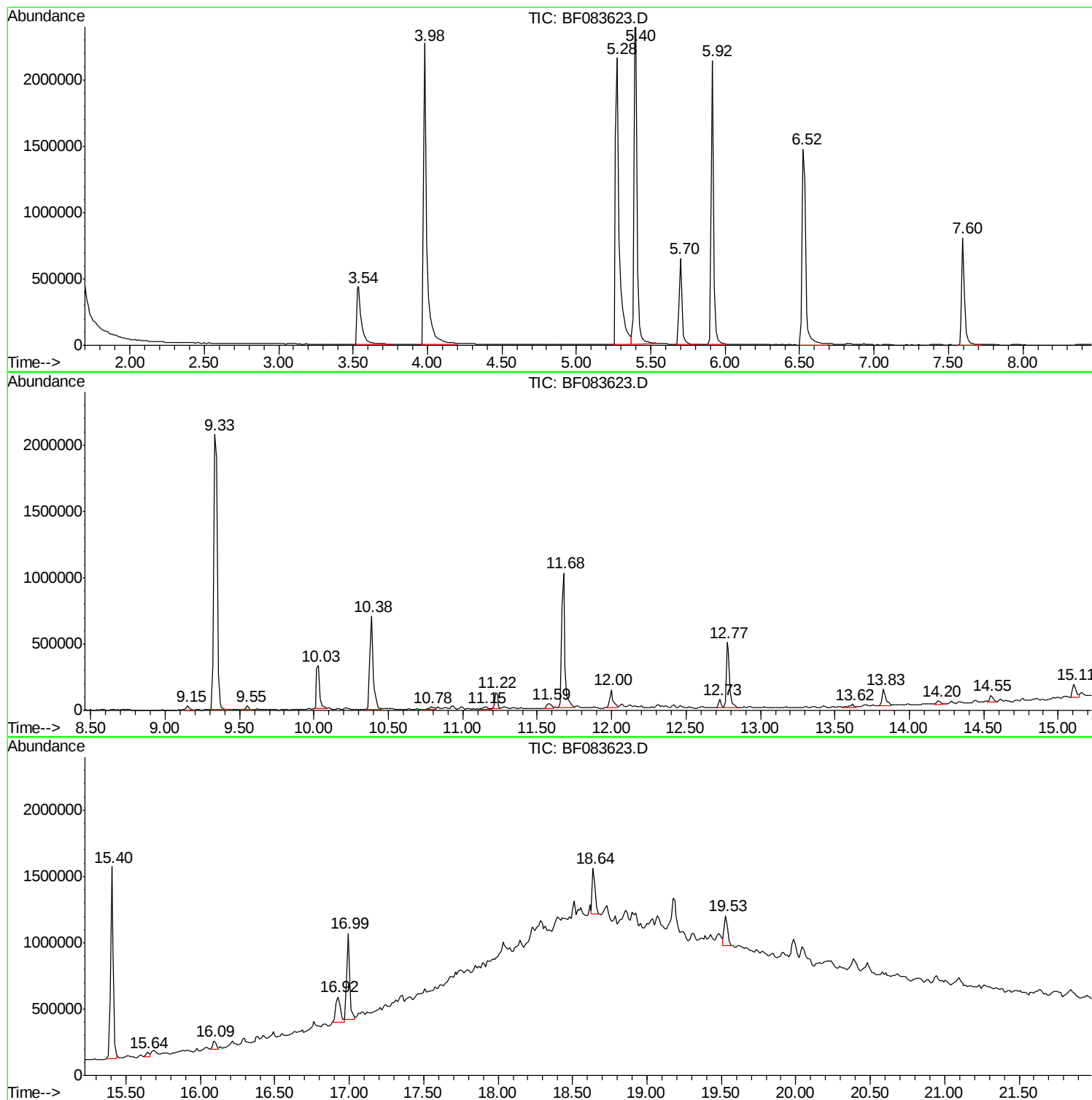
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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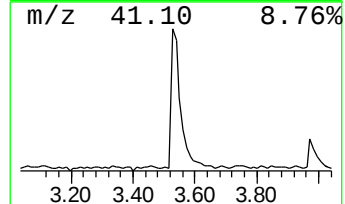
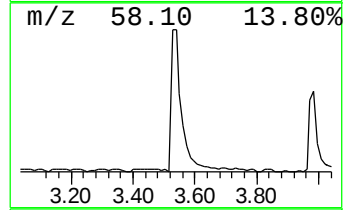
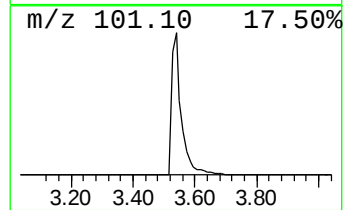
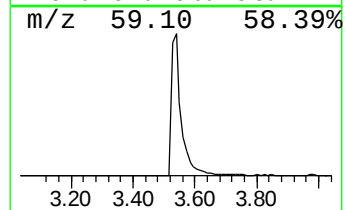
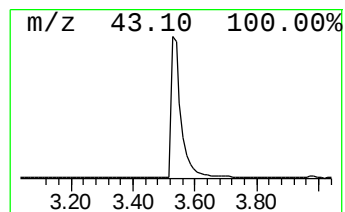
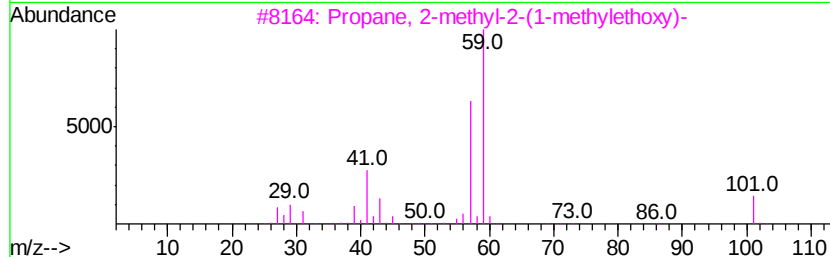
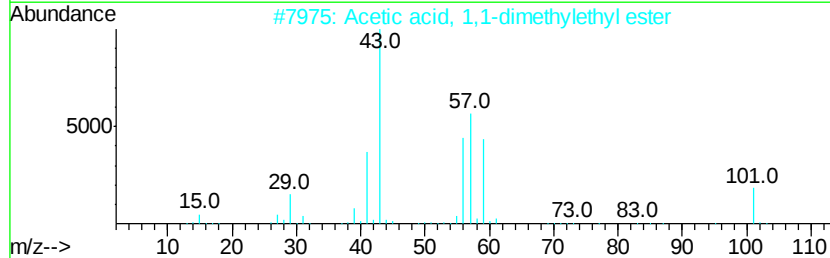
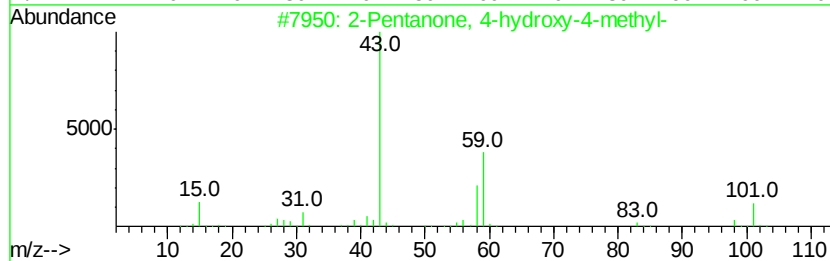
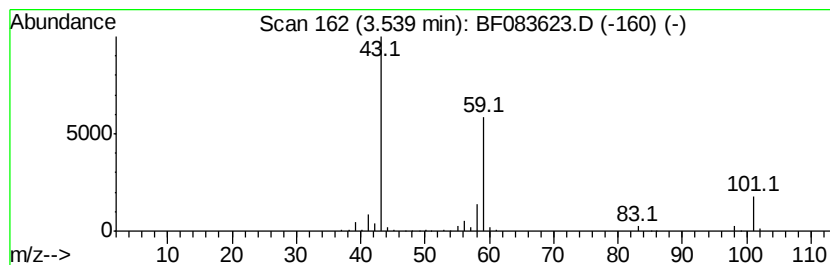
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.54	22.88 ng	989538	1,4-Dichlorobenzene-d4	5.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	33
4		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



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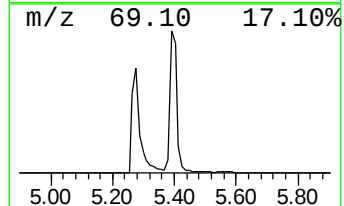
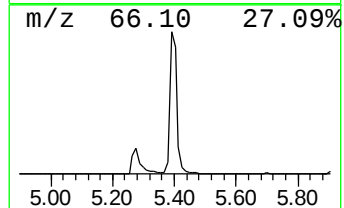
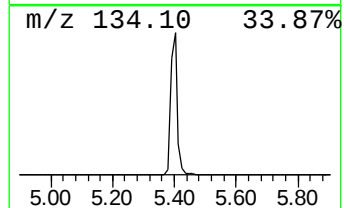
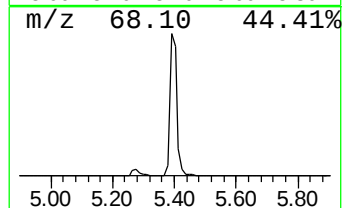
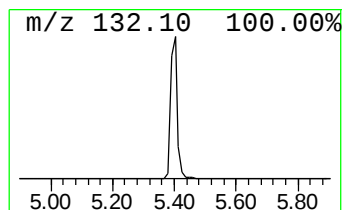
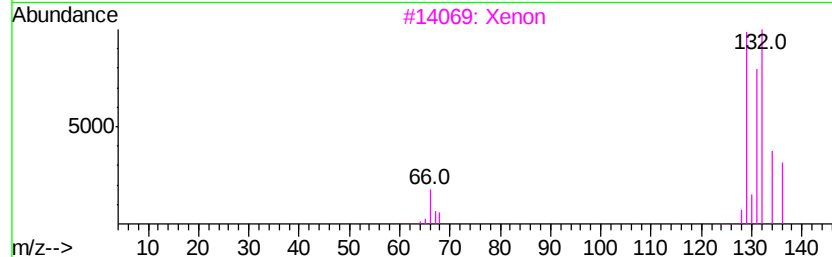
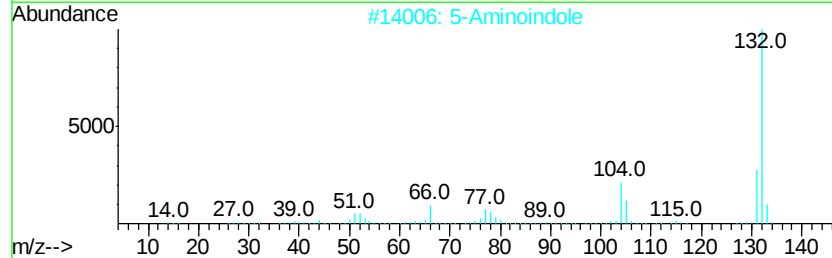
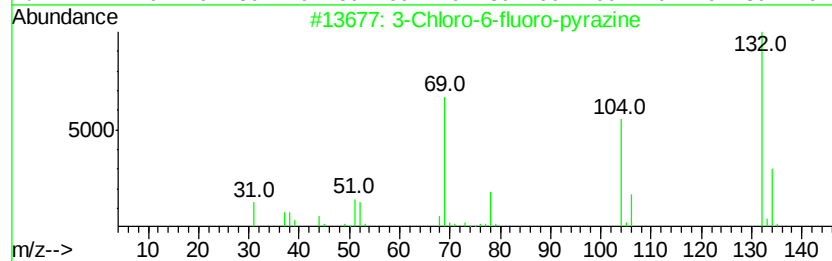
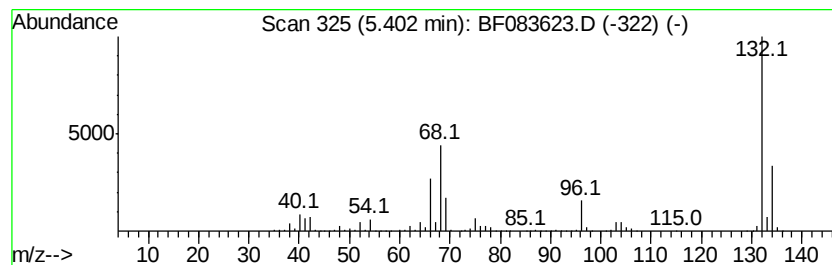
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown5.40 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.40	91.85 ng	3973230	1,4-Dichlorobenzene-d4	5.70

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3	Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5	Aminoindole	132	C8H8N2	005192-03-0	9
3		Xenon	132	Xe	007440-63-3	9
4		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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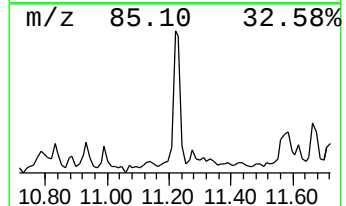
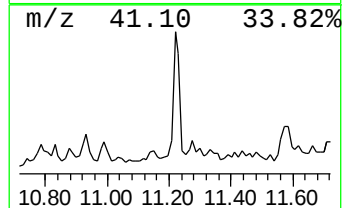
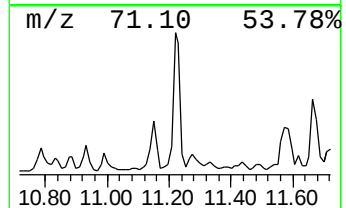
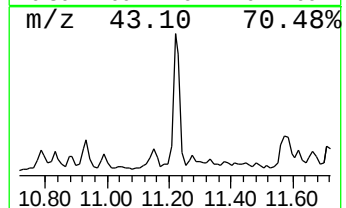
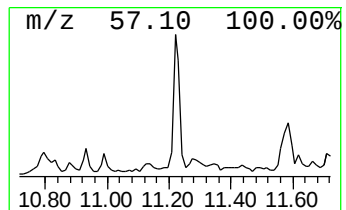
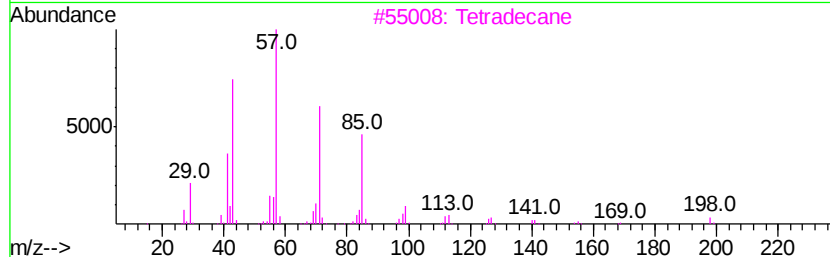
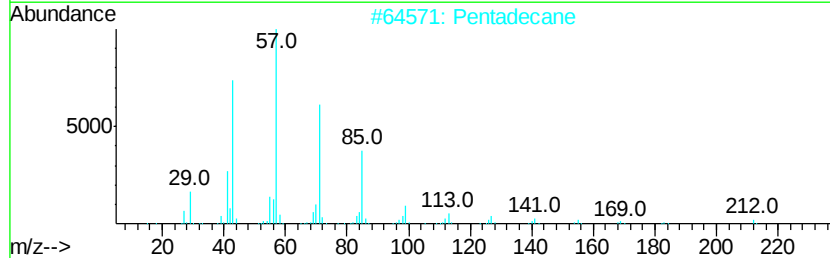
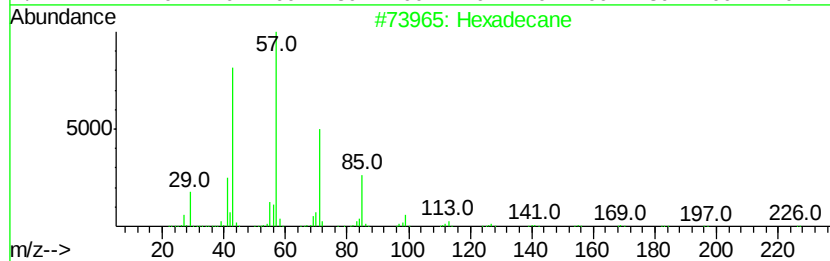
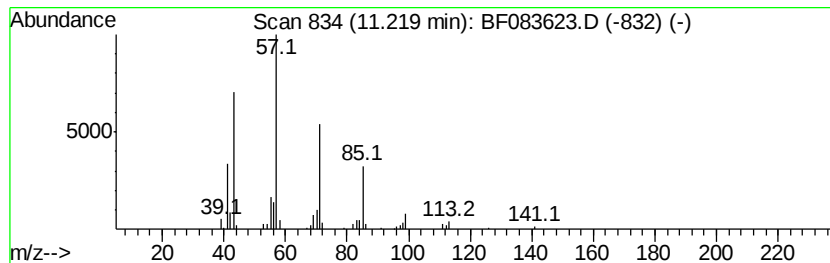
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Hexadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.22	3.33 ng	178521	Acenaphthene-d10	10.38

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	91
2	Pentadecane		212	C15H32	000629-62-9	90
3	Tetradecane		198	C14H30	000629-59-4	90
4	Decane, 2,3,5-trimethyl-		184	C13H28	062238-11-3	86
5	Heptadecane		240	C17H36	000629-78-7	83



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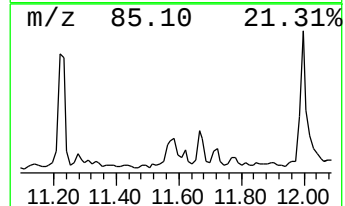
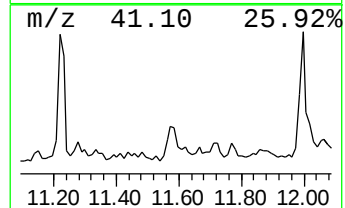
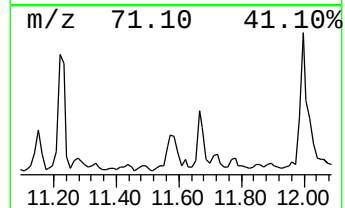
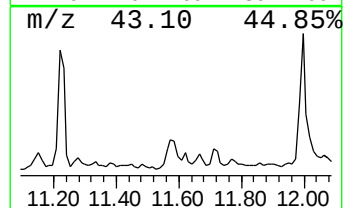
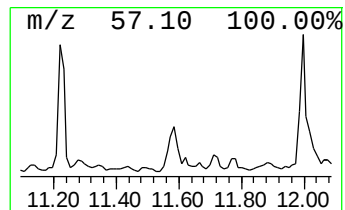
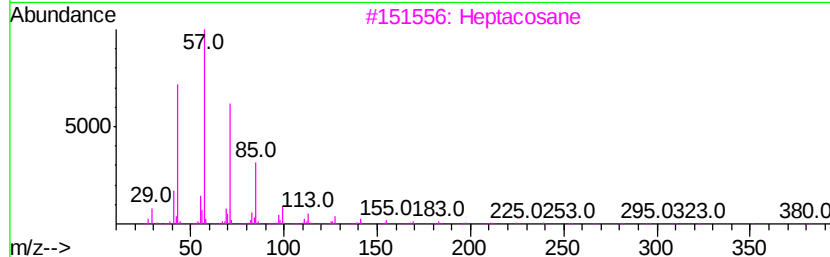
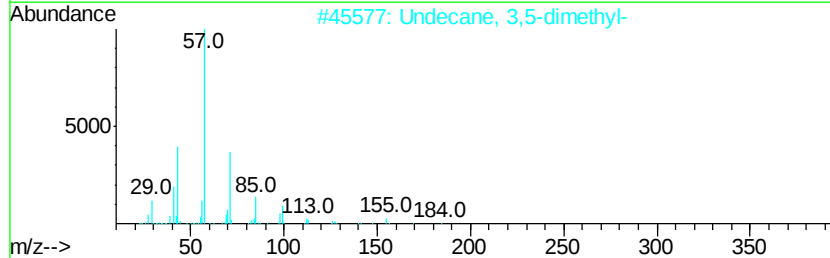
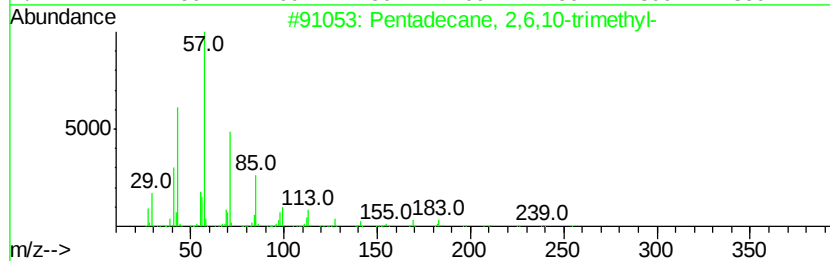
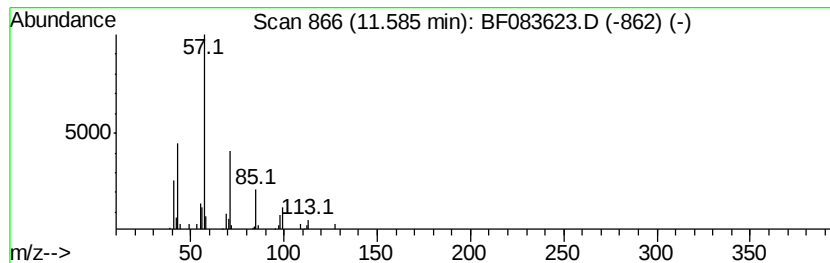
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Pentadecane, 2,6,10-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.59	2.08 ng	81615	Phenanthrene-d10	12.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	78
2		Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	64
3		Heptacosane	380	C27H56	000593-49-7	64
4		Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	59
5		Decane, 2,6,8-trimethyl-	184	C13H28	062108-26-3	59



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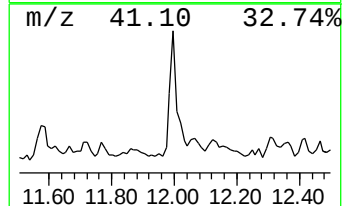
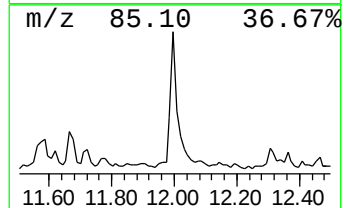
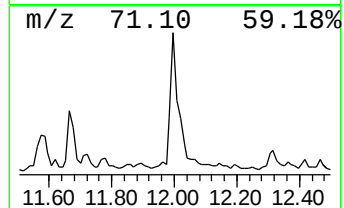
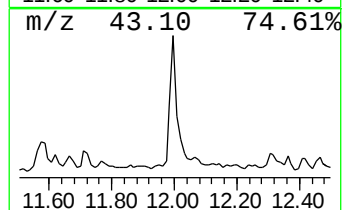
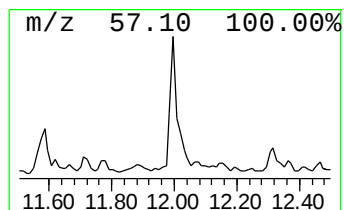
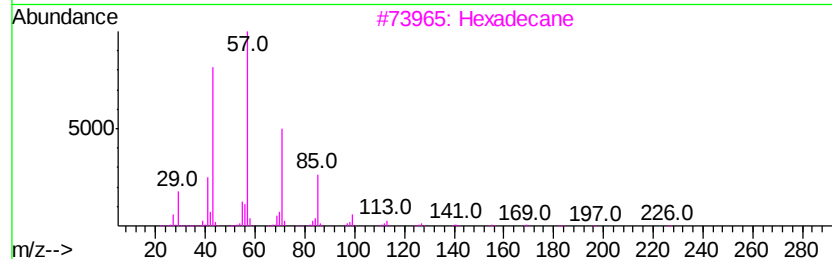
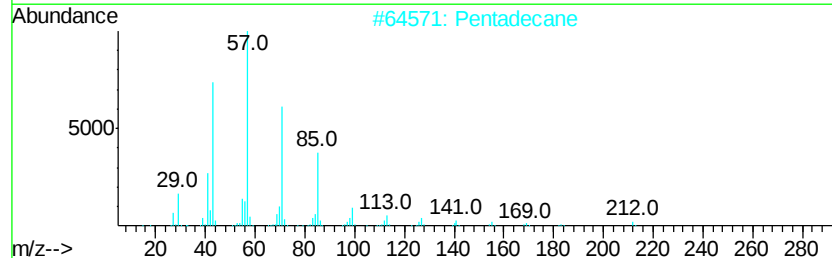
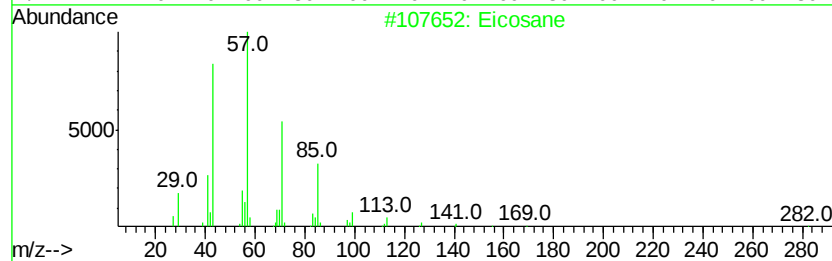
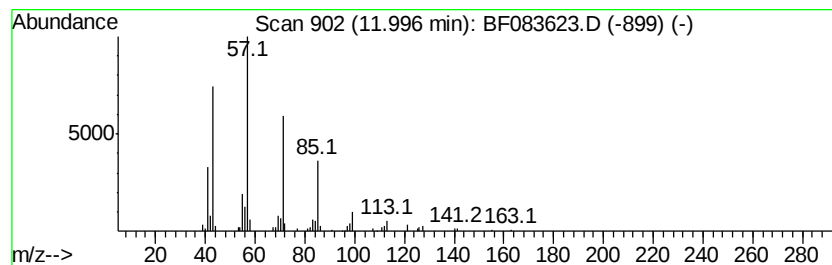
Instrument :
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Eicosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.00	5.51 ng	216314	Phenanthrene-d10	12.77	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	91
2	Pentadecane	212	C15H32	000629-62-9	90
3	Hexadecane	226	C16H34	000544-76-3	90
4	Octadecane	254	C18H38	000593-45-3	86
5	Tetradecane	198	C14H30	000629-59-4	72



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BRIDGE14-2(0-2)

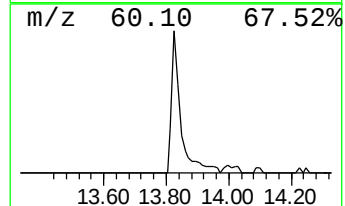
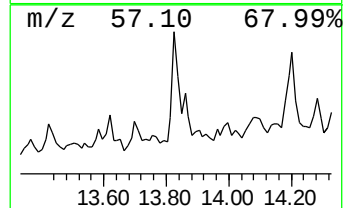
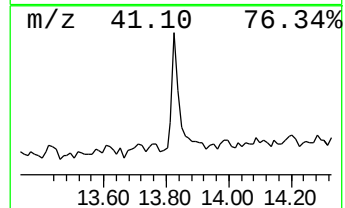
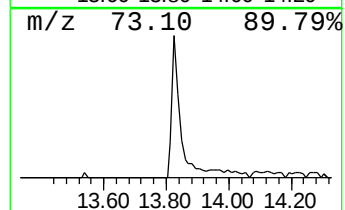
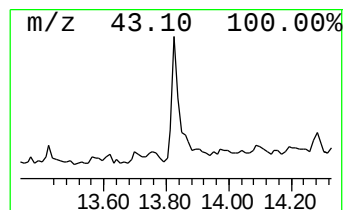
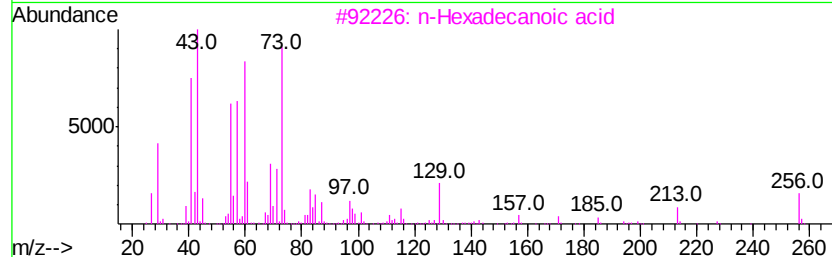
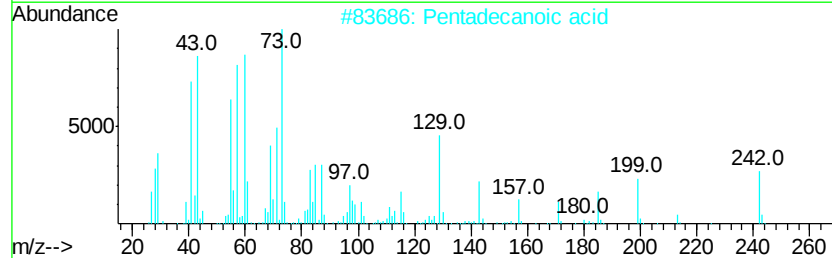
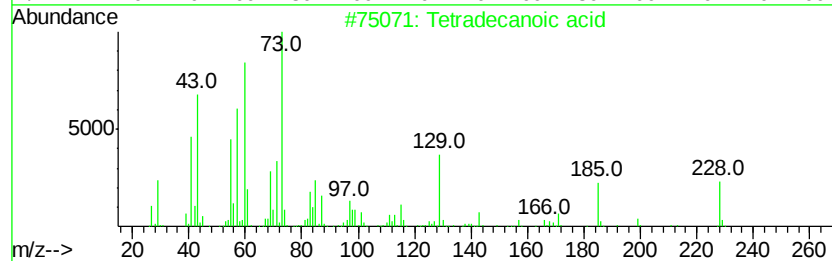
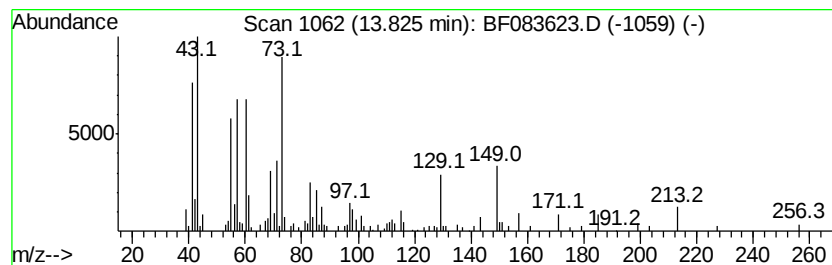
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetradecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	5.86 ng	230238	Phenanthrene-d10	12.77

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecanoic acid	228	C14H28O2	000544-63-8	62
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	59
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	58
4		Dodecanoic acid	200	C12H24O2	000143-07-7	53
5		Tridecanoic acid	214	C13H26O2	000638-53-9	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
Data File : BF083623.D
Acq On : 9 Dec 2015 5:38
Operator : UM/IZ
Sample : G4647-03
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BRIDGE14-2(0-2)

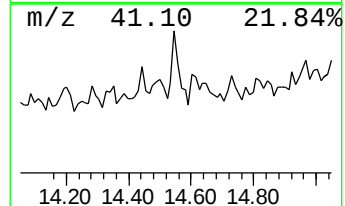
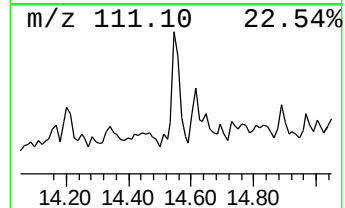
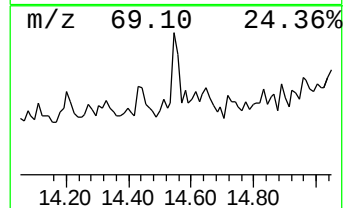
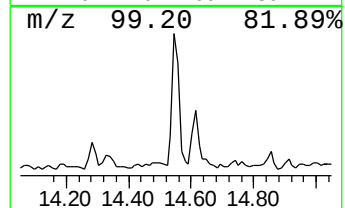
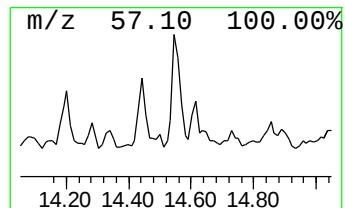
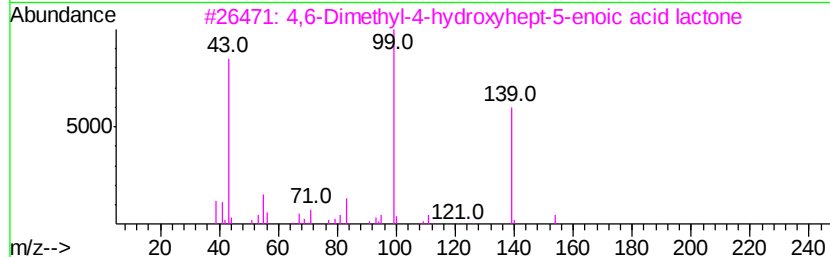
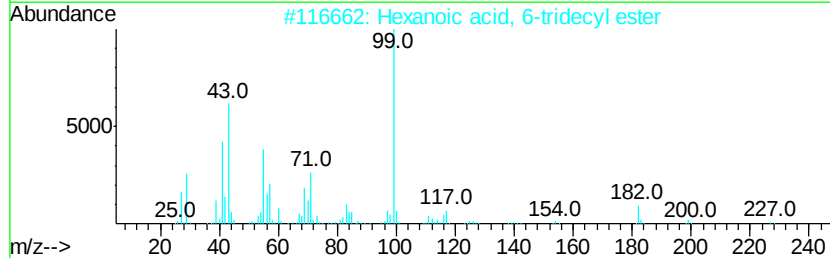
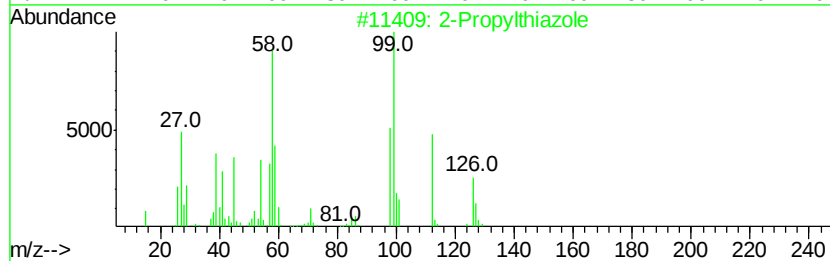
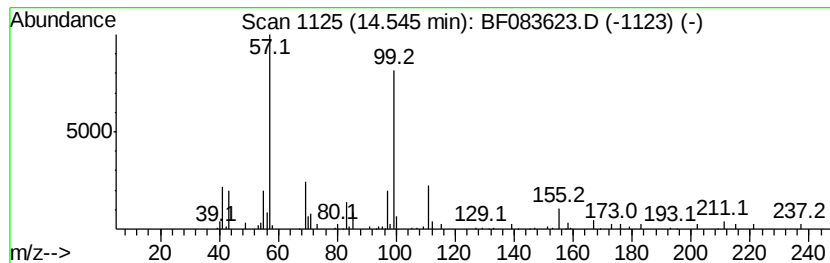
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown14.55 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.55	2.09 ng	82074	Phenanthrene-d10	12.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Propylthiazole	127	C6H9NS	017626-75-4	38
2		Hexanoic acid, 6-tridecyl ester	298	C19H38O2	1000279-29-5	32
3		4,6-Dimethyl-4-hydroxyhept-5-eno...	154	C9H14O2	1000122-47-0	25
4		Cyclohexanol, 1-ethyl-2,2-dimeth...	168	C11H20O	054345-64-1	25
5		Spiro(1,3-dioxolane)-2,3'-(pregn...	476	C27H40O7	1000194-70-1	25



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
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Operator : UM/IZ
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Instrument :
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ClientSampleId :
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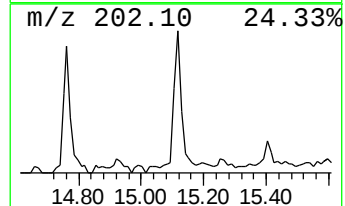
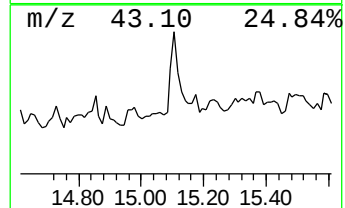
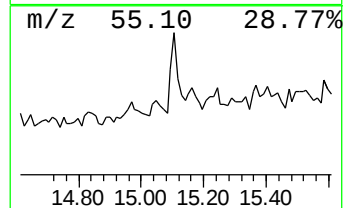
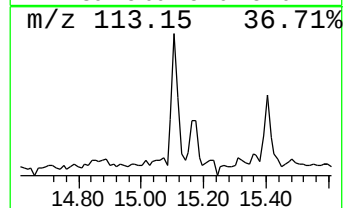
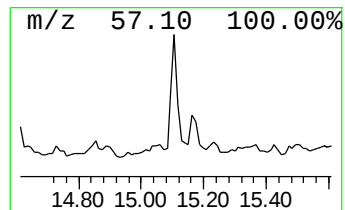
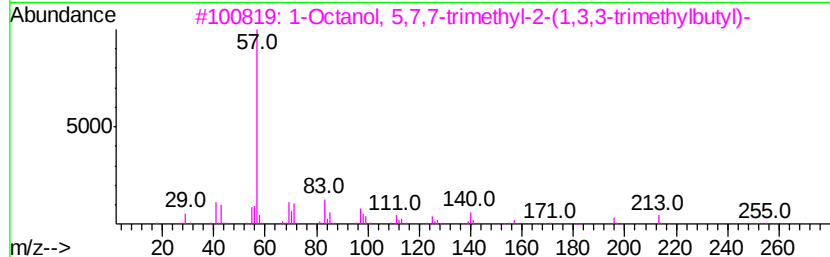
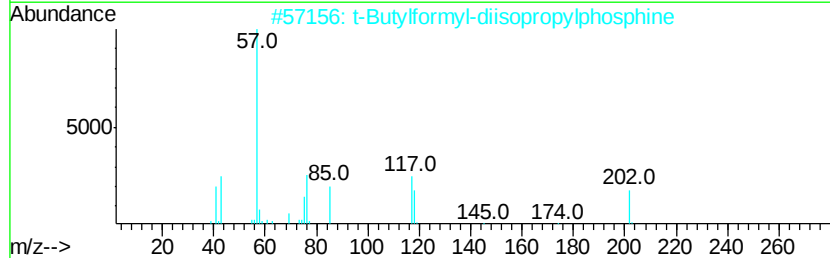
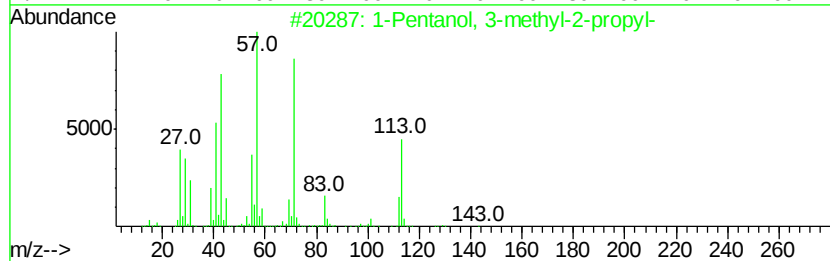
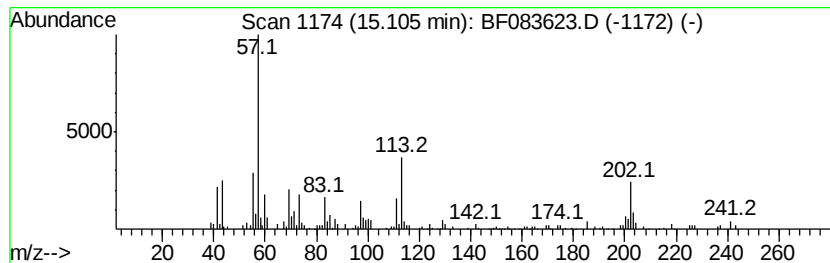
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 unknown15.11 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.11	3.89 ng	173044	Chrysene-d12	16.99

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	27
2			t-Butylformyl-diisopropylphosphine	202	C11H23OP	051049-58-2	22
3			1-Octanol, 5,7,7-trimethyl-2-(1,...	270	C18H38O	036400-98-3	14
4			2-Isothiocyano-2,4,4,6,6-pentame...	227	C13H25NS	1000293-28-0	14
5			Hentriacontane	437	C31H64	000630-04-6	12



Data Path : Z:\HPCHEM1\BNA F\DATA\BF120815\
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Operator : UM/IZ
Sample : G4647-03
Misc :
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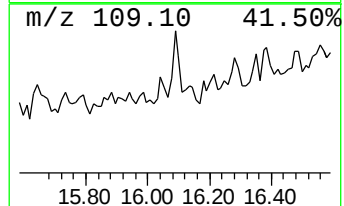
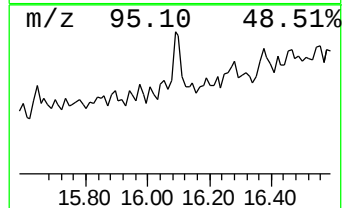
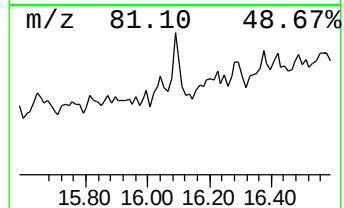
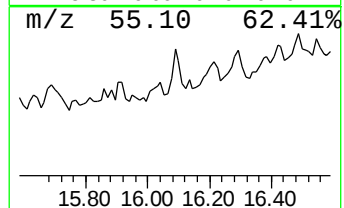
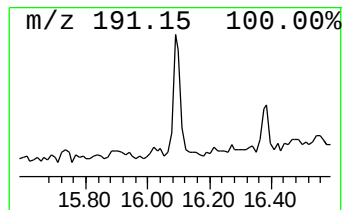
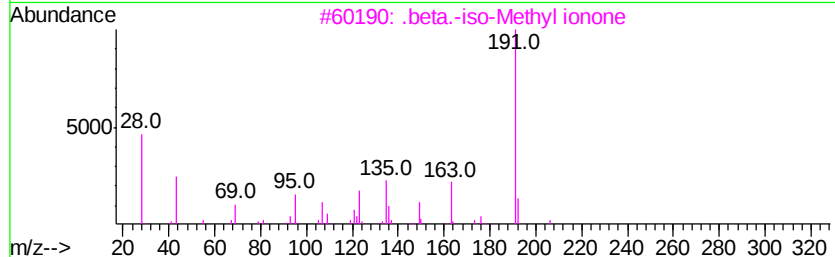
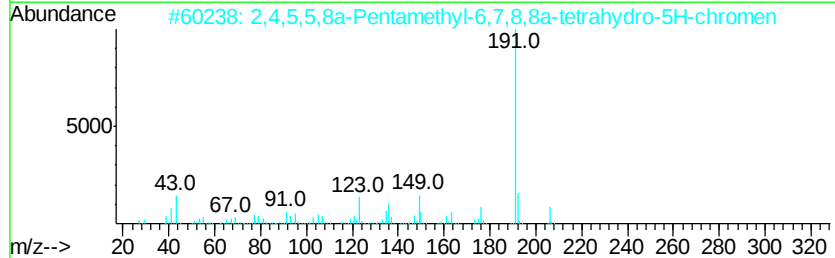
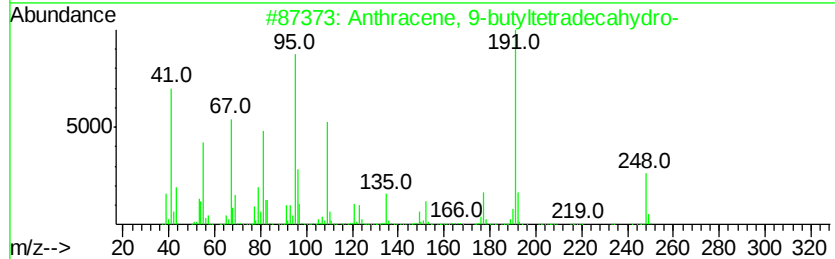
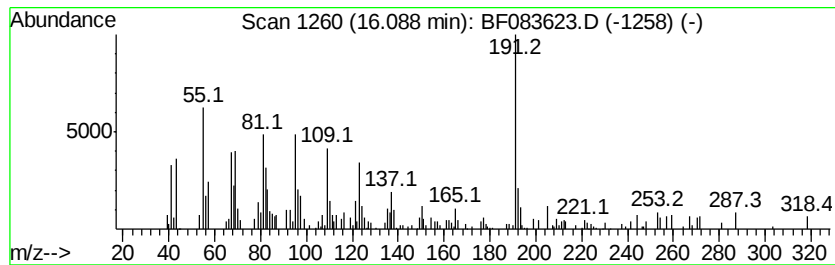
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown16.09 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.09	2.08 ng	92838	Chrysene-d12	16.99	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 9-butyltetradecahydro-	248	C18H32	055133-89-6	40
2	2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	35
3	.beta.-iso-Methyl ionone	206	C14H22O	1000285-40-2	32
4	Tridecanedial	212	C13H24O2	063521-76-6	25
5	Anthracene, 9-(3-butenyl)-	232	C18H16	097451-55-3	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
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Acq On : 9 Dec 2015 5:38
Operator : UM/IZ
Sample : G4647-03
Misc :
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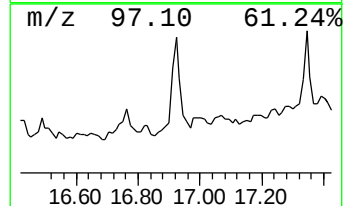
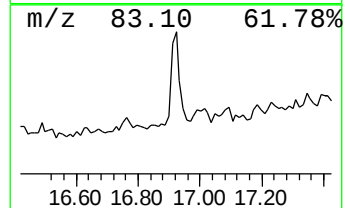
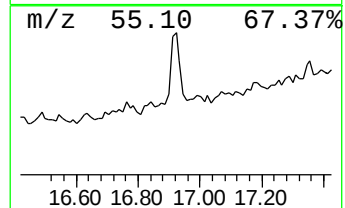
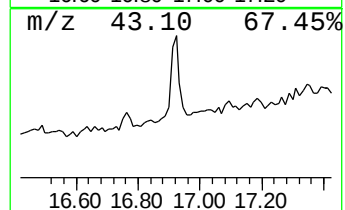
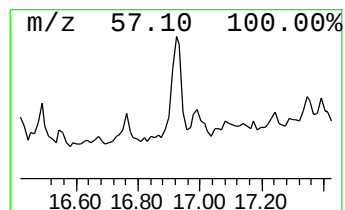
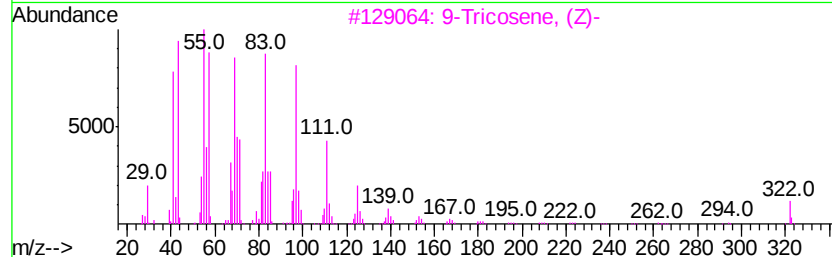
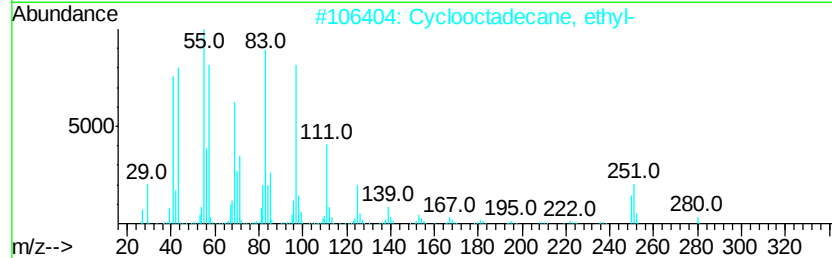
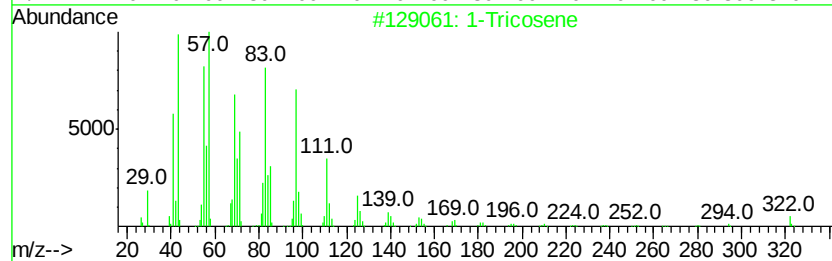
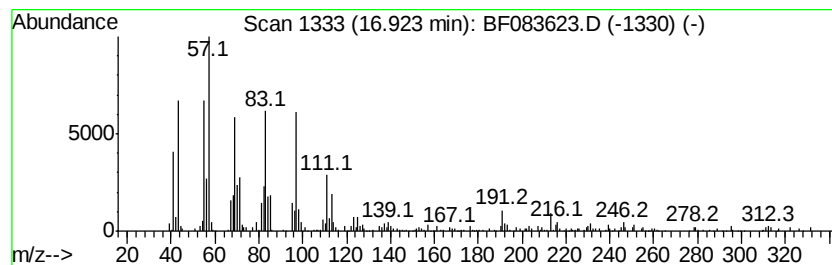
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Tricosene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.92	9.23 ng	411226	Chrysene-d12	16.99	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tricosene	322	C23H46	018835-32-0	80
2	Cyclooctadecane, ethyl-	280	C20H40	1000151-22-5	53
3	9-Tricosene, (Z)-	322	C23H46	027519-02-4	53
4	Decane, 5,6-bis(2,2-dimethylprop...	278	C20H38	073002-85-4	53
5	Acetic acid, chloro-, hexadecyl ...	318	C18H35ClO2	052132-58-8	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
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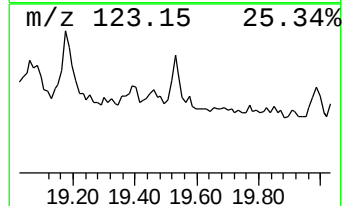
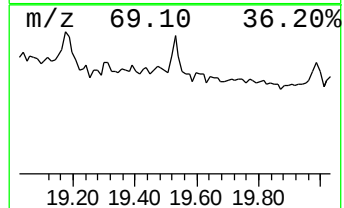
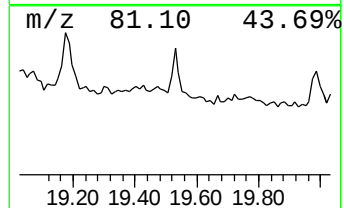
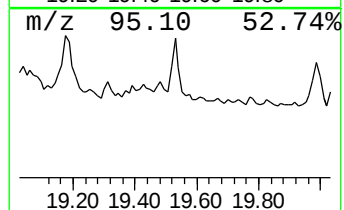
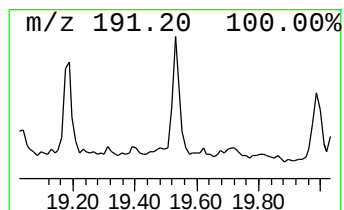
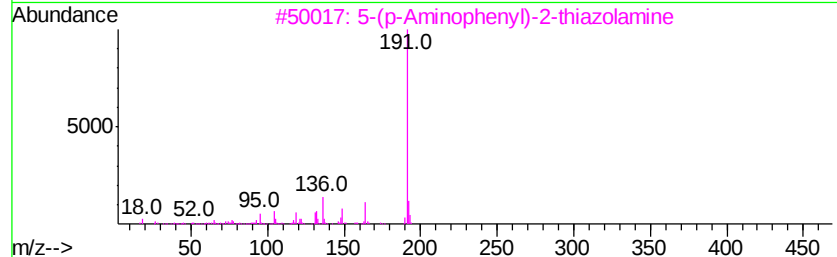
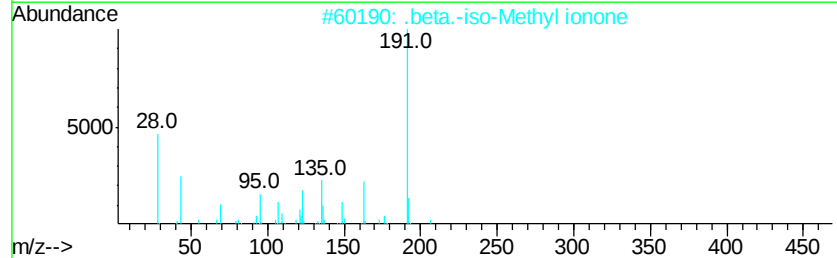
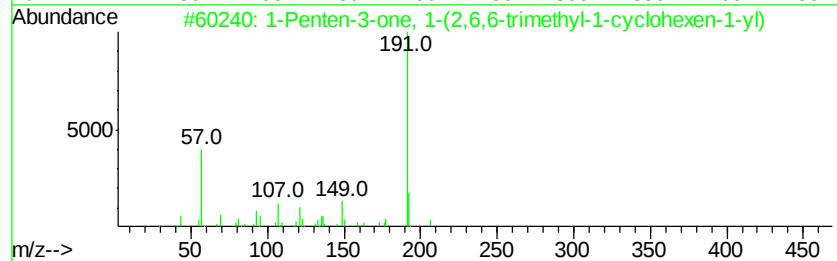
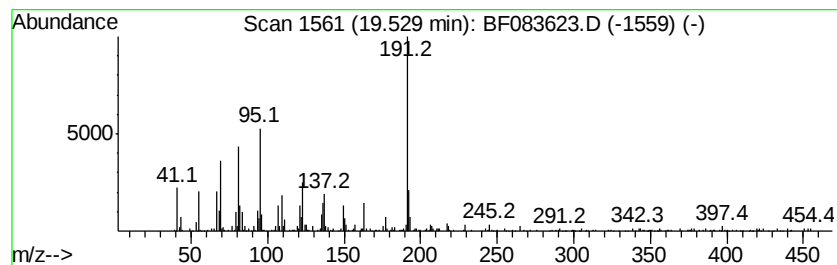
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Penten-3-one, 1-(2,6,6-tr... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.53	16.96 ng	365837	Perylene-d12	18.64
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Penten-3-one, 1-(2,6,6-trimeth...	206 C14H22O	000127-43-5	55
2	.beta.-iso-Methyl ionone	206 C14H22O	1000285-40-2	53
3	5-(p-Aminophenyl)-2-thiazolamine	191 C9H9N3S	090349-87-4	46
4	Anthracene, 9-dodecyltetradecahy...	360 C26H48	055401-75-7	45
5	1-Cyclohexene, 1,3,3-trimethyl-2...	206 C14H22O	1000197-08-4	45



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF120815\
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Acq On : 9 Dec 2015 5:38
Operator : UM/IZ
Sample : G4647-03
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF120815.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	3.54	22.9	ng	989538	1	5.70	865133	20.0
unknown5.40	5.40	91.8	ng	3973230	1	5.70	865133	20.0
Hexadecane	11.22	3.3	ng	178521	3	10.38	1072670	20.0
Pentadecane, 2,6,...	11.59	2.1	ng	81615	4	12.77	785266	20.0
Eicosane	12.00	5.5	ng	216314	4	12.77	785266	20.0
Tetradecanoic acid	13.83	5.9	ng	230238	4	12.77	785266	20.0
unknown14.55	14.55	2.1	ng	82074	4	12.77	785266	20.0
unknown15.11	15.11	3.9	ng	173044	5	16.99	890700	20.0
unknown16.09	16.09	2.1	ng	92838	5	16.99	890700	20.0
1-Tricosene	16.92	9.2	ng	411226	5	16.99	890700	20.0
1-Penten-3-one, 1...	19.53	17.0	ng	365837	6	18.64	431394	20.0