

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110415\
 Data File : BF082692.D
 Acq On : 4 Nov 2015 9:11
 Operator : UM/IZ
 Sample : G4241-11
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-12(2-4)

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-BF103015.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.070	258	261	271	rVB	1530973	2253838	38.13%	4.695%
2	5.493	295	298	300	rVB	5165988	5401991	91.40%	11.253%
3	6.522	384	388	390	rVB	4092302	5910492	100.00%	12.312%
4	6.647	397	399	402	rBV	4098497	5337363	90.30%	11.118%
5	6.887	417	420	421	rBV	974993	1064508	18.01%	2.218%
6	7.047	431	434	436	rVB	2897930	3729700	63.10%	7.769%
7	7.447	466	469	471	rBV	3593435	3656491	61.86%	7.617%
8	7.539	475	477	479	rBV	54017	61888	1.05%	0.129%
9	7.893	507	508	511	rBV2	41061	63680	1.08%	0.133%
10	8.168	530	532	536	rBV	1290379	1338289	22.64%	2.788%
11	8.419	553	554	557	rVB	65767	62556	1.06%	0.130%
12	8.785	583	586	592	rBV2	40735	72553	1.23%	0.151%
13	8.888	592	595	596	rBV	66270	59459	1.01%	0.124%
14	9.128	614	616	620	rVB	106241	122995	2.08%	0.256%
15	9.253	624	627	629	rBV	5129250	4627929	78.30%	9.641%
16	9.642	659	661	663	rBV	1135623	875213	14.81%	1.823%
17	9.871	680	681	684	rBV2	73639	63532	1.07%	0.132%
18	9.928	684	686	689	rVB	1648679	1397149	23.64%	2.910%
19	10.248	710	714	720	rVB3	33125	85403	1.44%	0.178%
20	10.339	720	722	724	rVV	116007	142626	2.41%	0.297%
21	10.373	724	725	726	rVB	105969	72642	1.23%	0.151%
22	10.591	740	744	746	rBV	42653	62554	1.06%	0.130%
23	10.716	752	755	757	rBV	3258327	2951967	49.94%	6.149%
24	10.842	765	766	771	rVB	140170	201018	3.40%	0.419%
25	11.071	782	786	790	rVB6	22735	65316	1.11%	0.136%
26	11.288	804	805	807	rBV	78289	93721	1.59%	0.195%
27	11.414	814	816	818	rBV	1337440	1198649	20.28%	2.497%
28	11.642	834	836	841	rBV	106154	180970	3.06%	0.377%
29	11.722	841	843	845	rVB	96582	81802	1.38%	0.170%
30	11.951	861	863	865	rBV2	337904	334637	5.66%	0.697%
31	12.134	875	879	881	rVB2	56268	99533	1.68%	0.207%
32	12.465	906	908	910	rBV2	42597	61405	1.04%	0.128%
33	12.625	919	922	924	rBV2	40240	70509	1.19%	0.147%
34	12.671	924	926	930	rBV4	58033	98298	1.66%	0.205%

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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

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35	12.739	930	932	934	rBV2	65931	110012	1.86%	0.229%
36	12.831	938	940	943	rBV2	194313	224479	3.80%	0.468%
37	12.888	943	945	948	rBV2	43371	81714	1.38%	0.170%
38	13.002	953	955	957	rBV	3080933	2739191	46.34%	5.706%
39	13.540	1000	1002	1004	rBV2	158899	180123	3.05%	0.375%
40	13.917	1032	1035	1037	rBV	332949	527325	8.92%	1.098%
41	14.054	1045	1047	1050	rBV	985037	1047127	17.72%	2.181%
42	14.557	1089	1091	1096	rBV2	109826	200186	3.39%	0.417%
43	15.528	1173	1176	1178	rBV	763965	993859	16.82%	2.070%

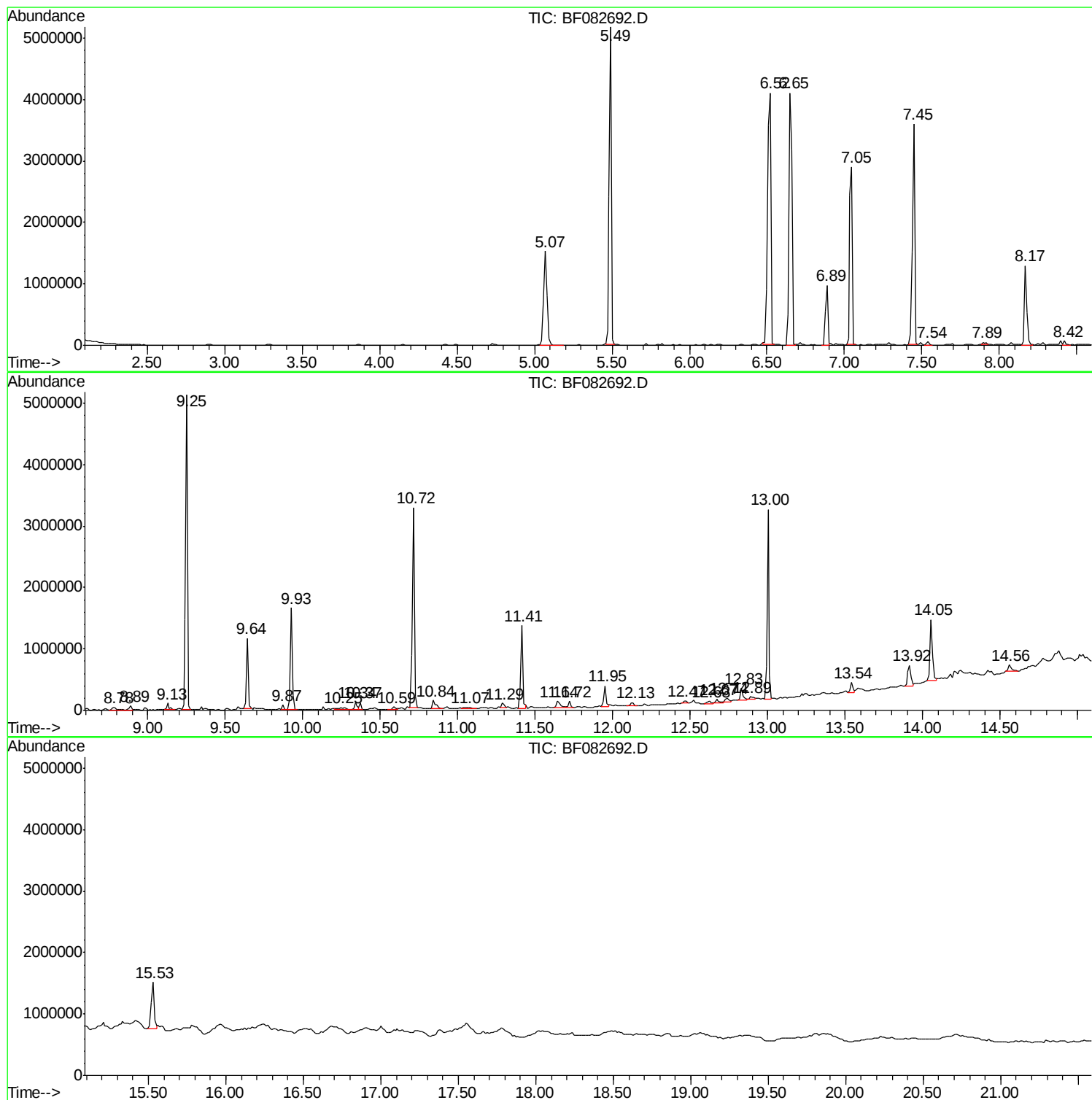
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF103015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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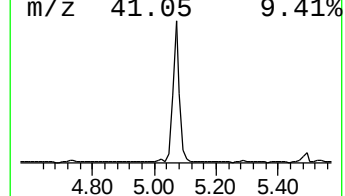
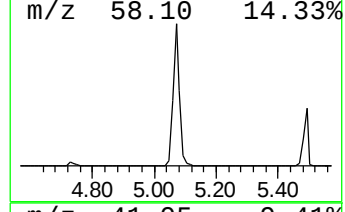
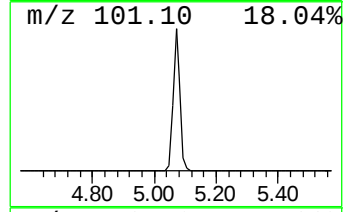
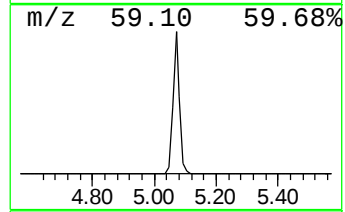
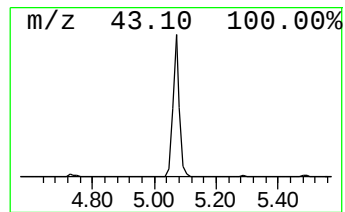
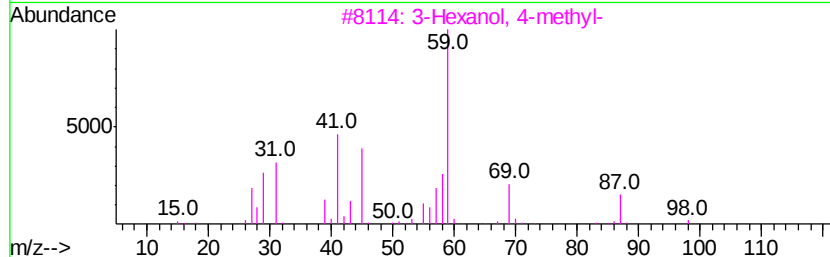
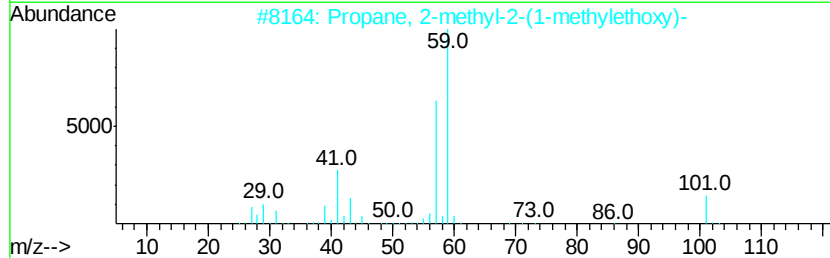
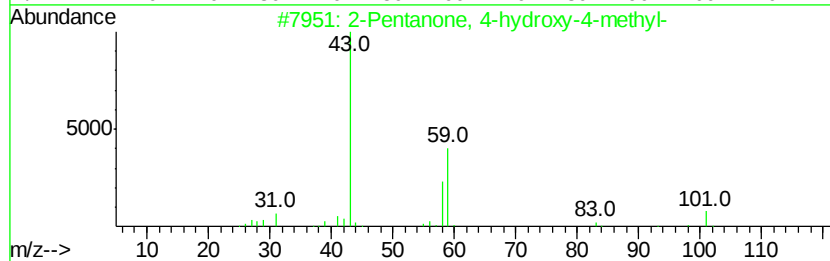
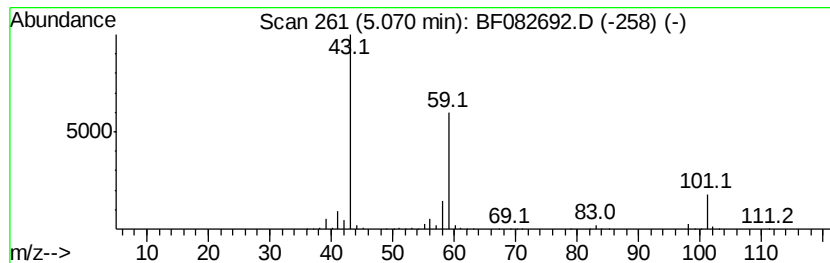
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF103015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
5.07	42.35 ng	2253840	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	33
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	32
5		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28



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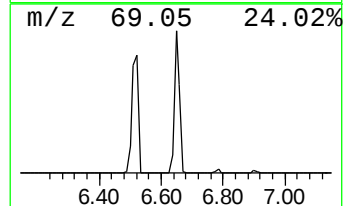
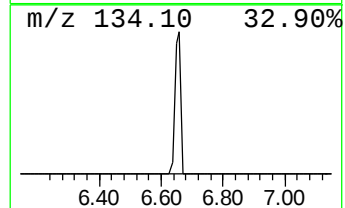
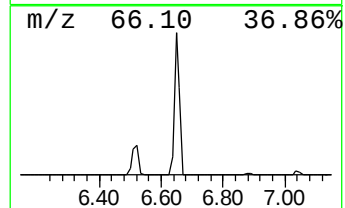
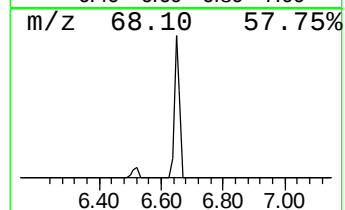
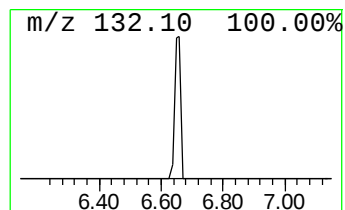
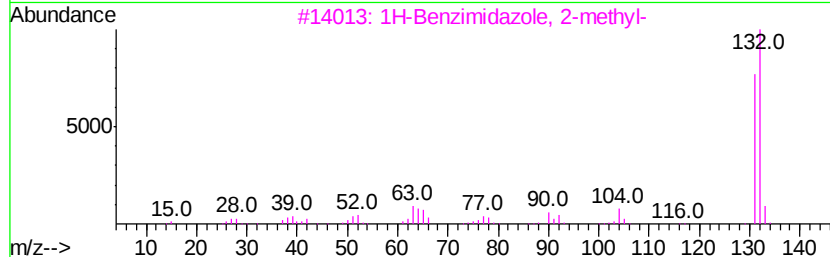
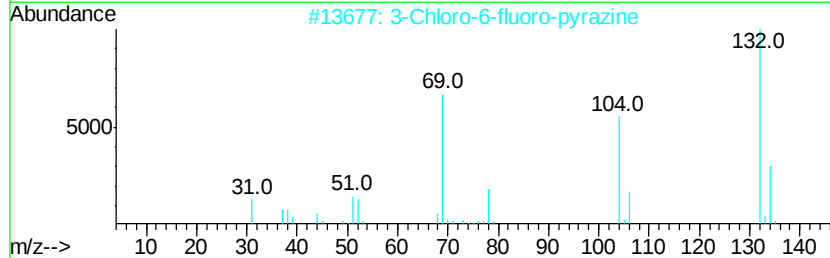
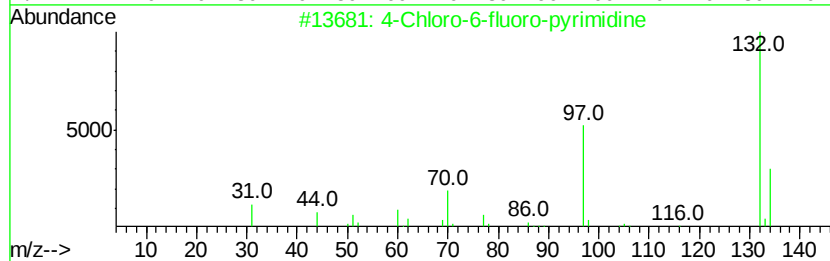
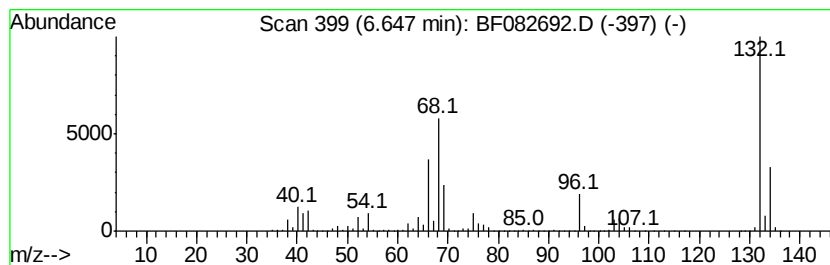
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Peak Number 2 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	100.28 ng	5337360	1,4-Dichlorobenzene-d4	6.89

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
5		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	10



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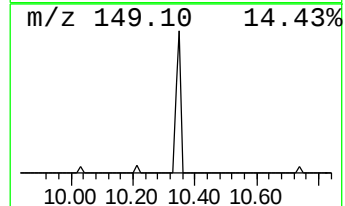
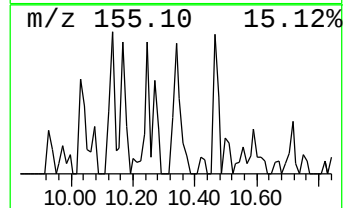
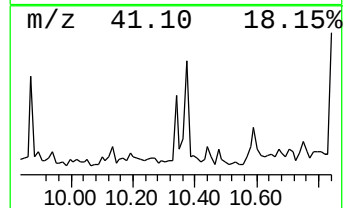
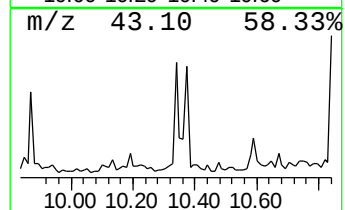
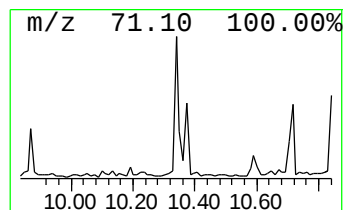
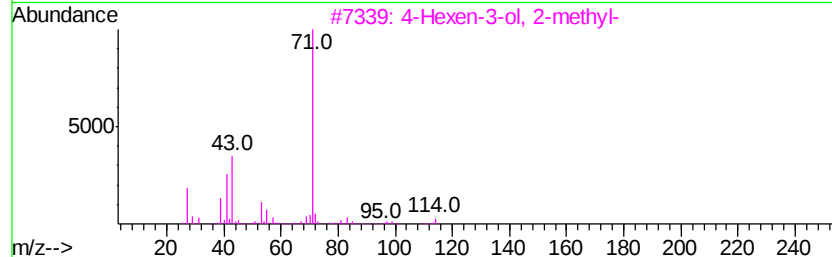
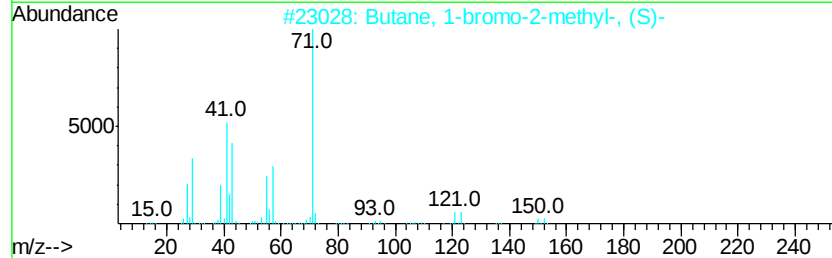
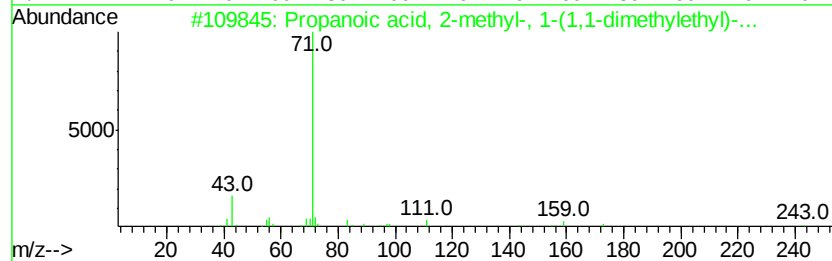
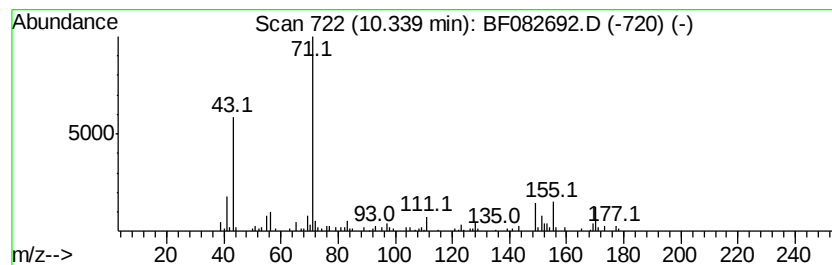
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 unknown10.34 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.34	2.04 ng	142626	Acenaphthene-d10	9.93
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Propanoic acid, 2-methyl-, 1-(1,...	286 C16H30O4	074381-40-1 47
2		Butane, 1-bromo-2-methyl-, (S)-	150 C5H11Br	000534-00-9 47
3		4-Hexen-3-ol, 2-methyl-	114 C7H14O	004798-60-1 43
4		Cyclohexanol, 2,2-dichloro-1-met...	182 C7H12Cl2O	1000115-65-2 43
5		3-Buten-2-ol, 1-bromo-2-methyl-	164 C5H9BrO	036219-40-6 40



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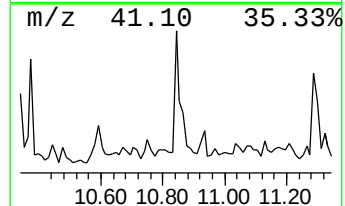
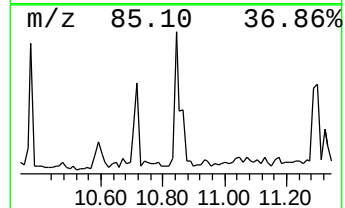
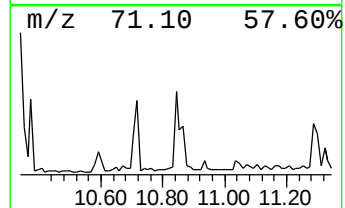
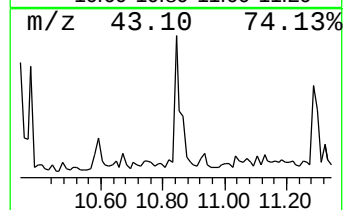
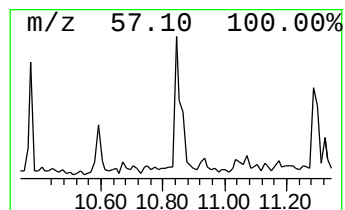
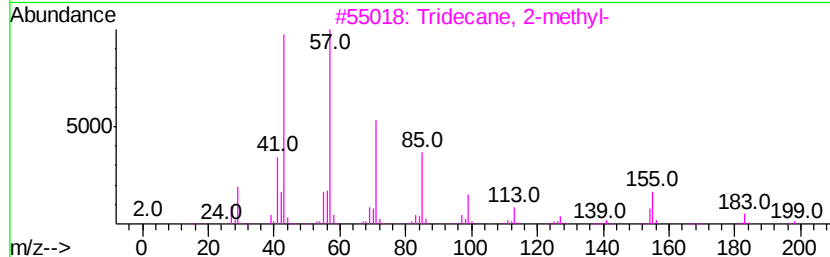
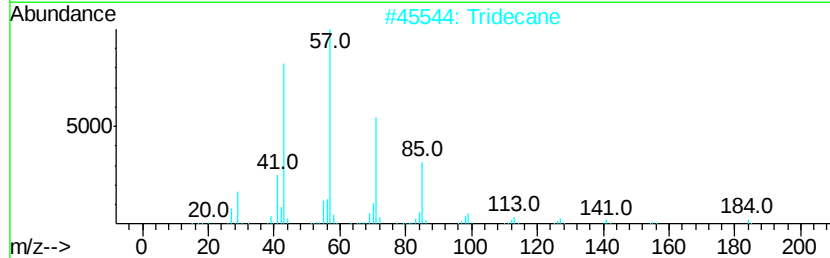
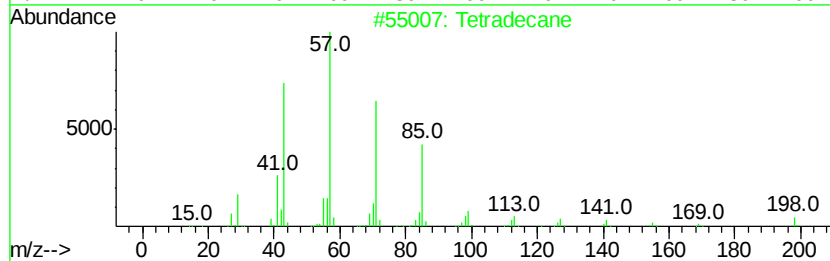
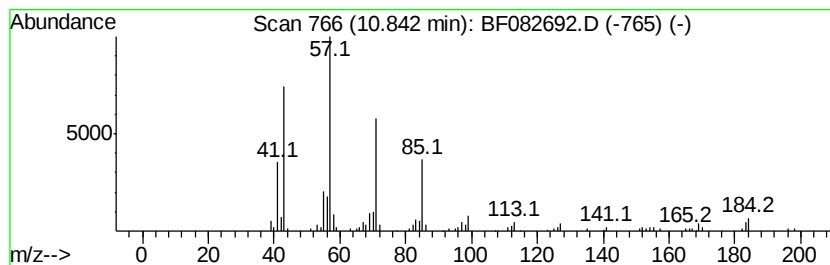
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Tetradecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.84	3.35 ng	201018	Phenanthrene-d10	11.41

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane	198	C14H30	000629-59-4	94
2		Tridecane	184	C13H28	000629-50-5	93
3		Tridecane, 2-methyl-	198	C14H30	001560-96-9	83
4		Dodecane	170	C12H26	000112-40-3	81
5		Eicosane	282	C20H42	000112-95-8	81



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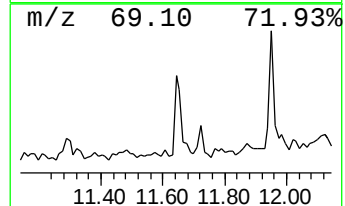
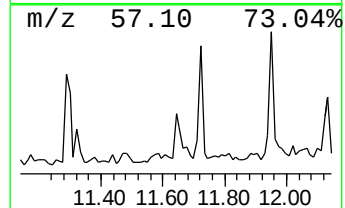
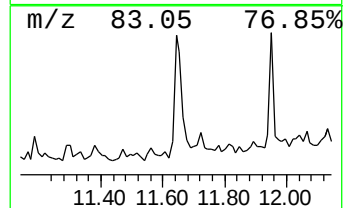
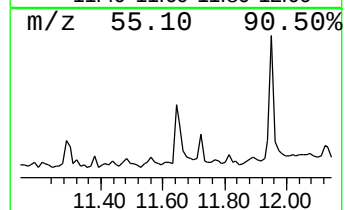
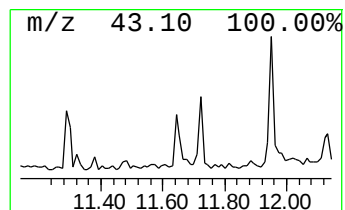
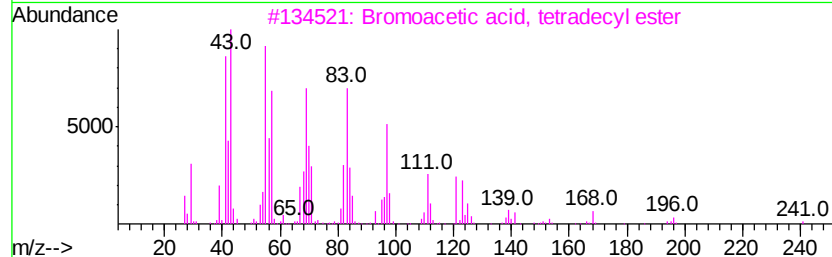
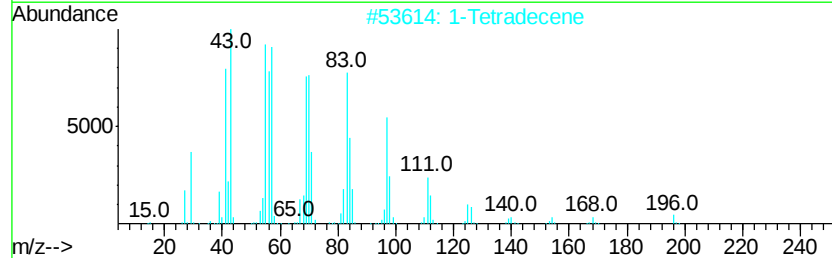
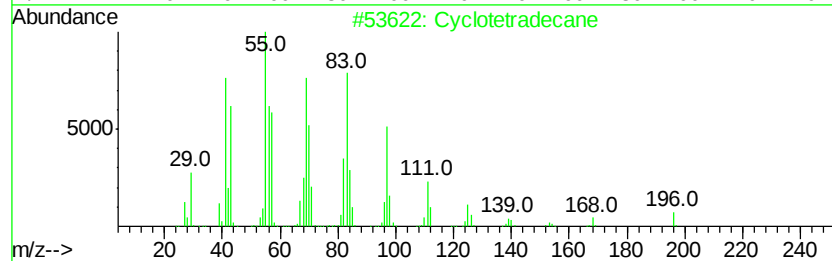
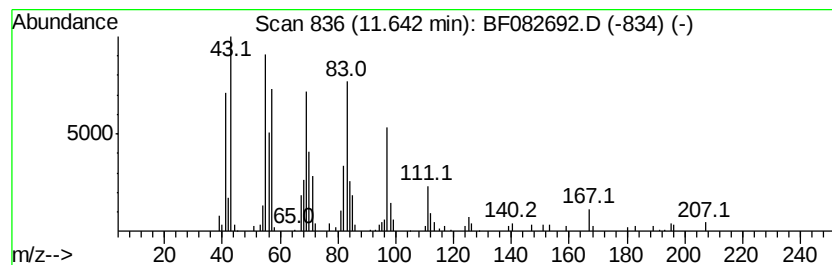
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF103015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Cyclotetradecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.64	3.02 ng	180970	Phenanthrene-d10	11.41

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclotetradecane	196	C14H28	000295-17-0	94
2		1-Tetradecene	196	C14H28	001120-36-1	92
3		Bromoacetic acid, tetradecyl ester	334	C16H31BrO2	018992-01-3	91
4		Hexadecen-1-ol, trans-9-	240	C16H32O	064437-47-4	91
5		1-Heptadecanol	256	C17H36O	001454-85-9	90



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110415\
Data File : BF082692.D
Acq On : 4 Nov 2015 9:11
Operator : UM/IZ
Sample : G4241-11
Misc :
ALS Vial : 36 Sample Multiplier: 1

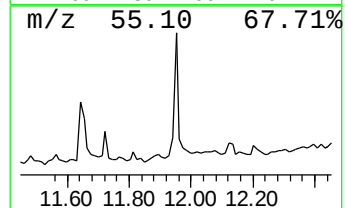
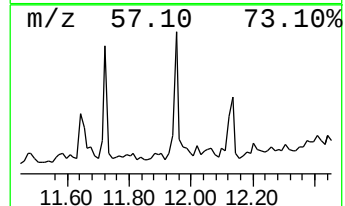
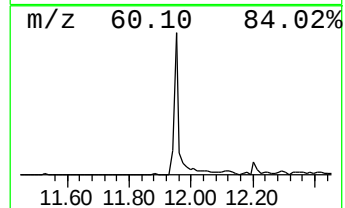
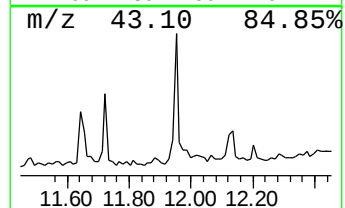
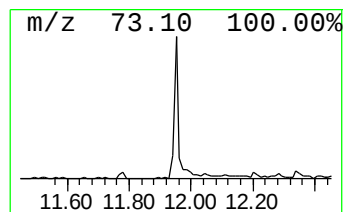
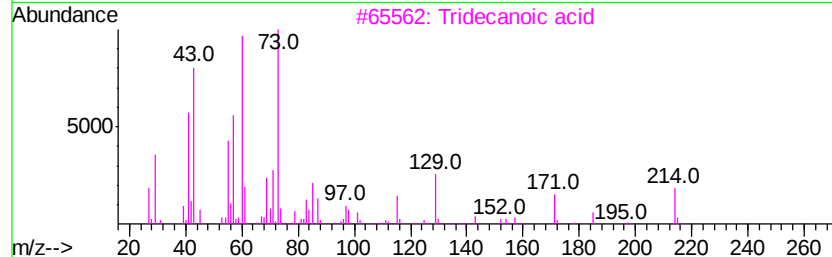
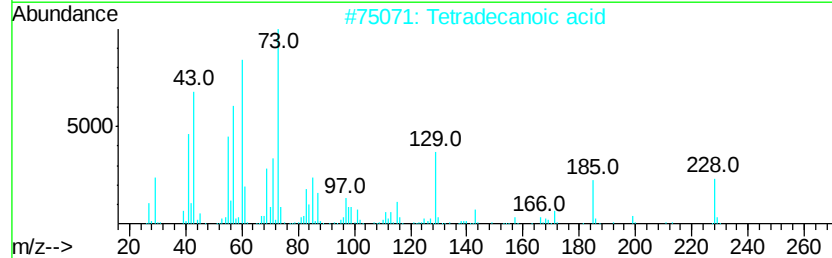
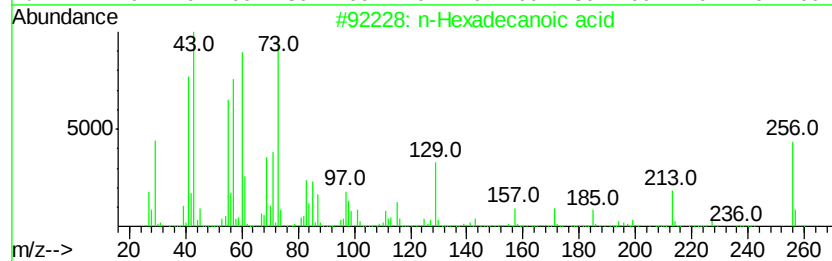
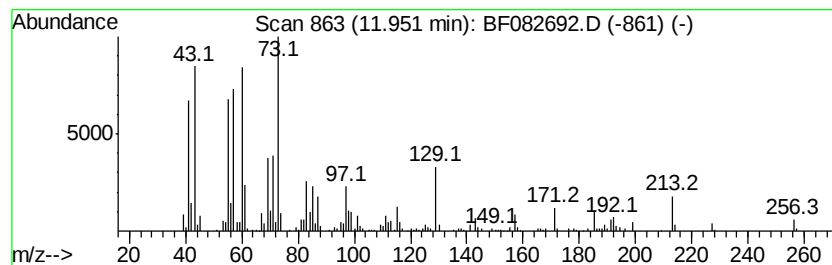
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ClientSampleId :
SB-12(2-4)

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF103015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.95	5.58 ng	334637	Phenanthrene-d10	11.41	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	93
3	Tridecanoic acid	214	C13H26O2	000638-53-9	76
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	50
5	Amyl Nitrite	117	C5H11NO2	000110-46-3	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110415\
Data File : BF082692.D
Acq On : 4 Nov 2015 9:11
Operator : UM/IZ
Sample : G4241-11
Misc :
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Instrument :
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ClientSampleId :
SB-12(2-4)

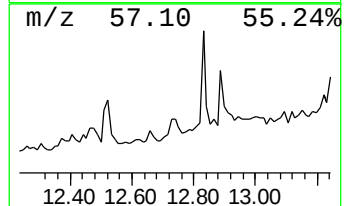
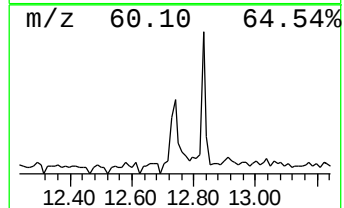
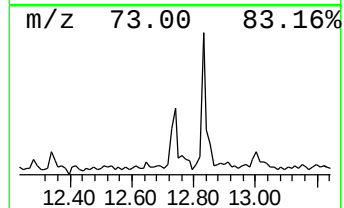
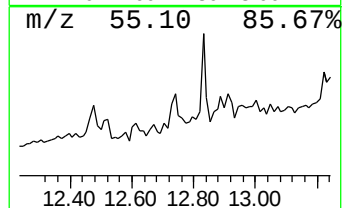
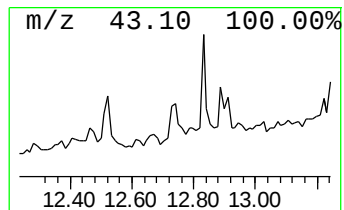
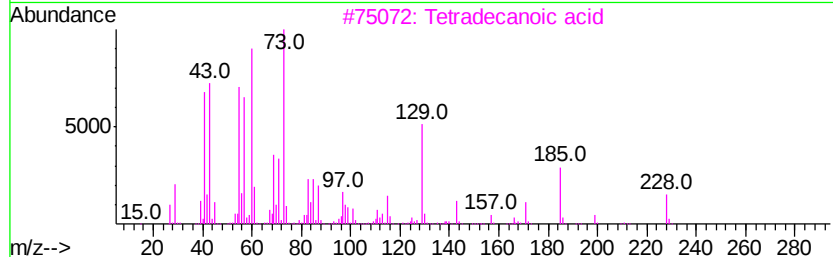
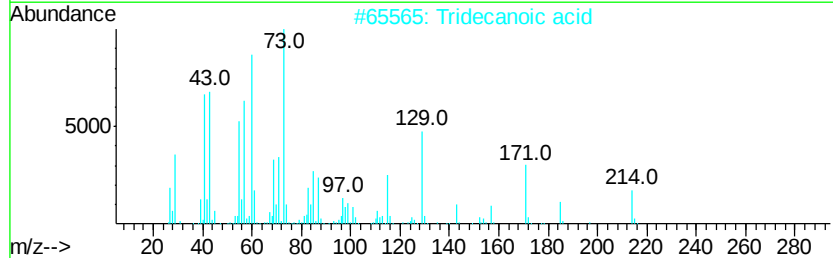
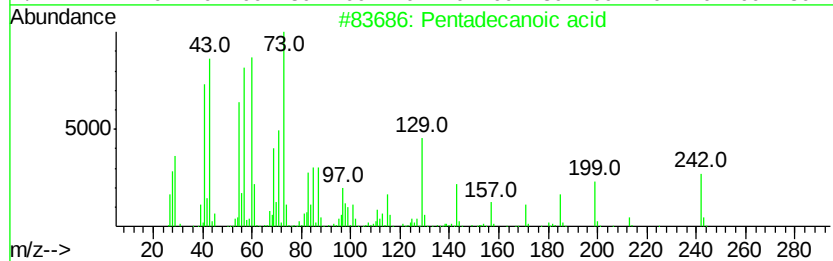
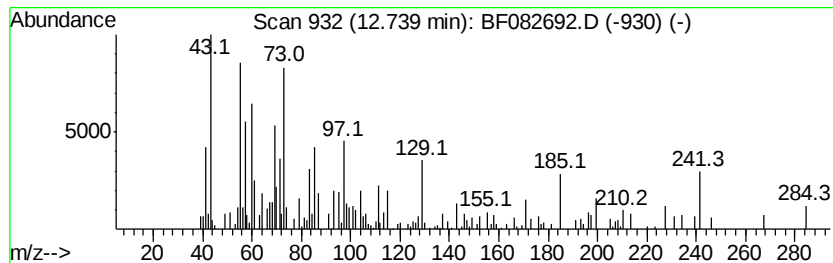
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 unknown12.74 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	2.10 ng	110012	Chrysene-d12	14.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecanoic acid	242	C15H30O2	001002-84-2	38
2			Tridecanoic acid	214	C13H26O2	000638-53-9	30
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	22
4			5-Acetoxy-pentadecane	270	C17H34O2	1000245-62-3	22
5			Cyclopropanecarboxylic acid, 4-h...	310	C20H38O2	1000282-78-4	22



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110415\
Data File : BF082692.D
Acq On : 4 Nov 2015 9:11
Operator : UM/IZ
Sample : G4241-11
Misc :
ALS Vial : 36 Sample Multiplier: 1

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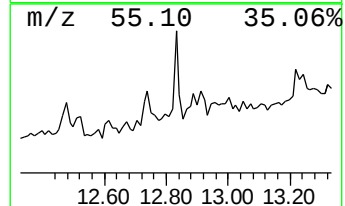
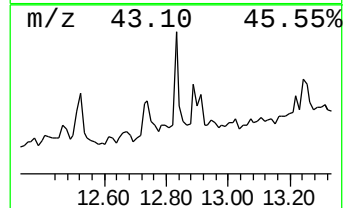
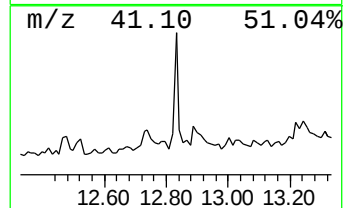
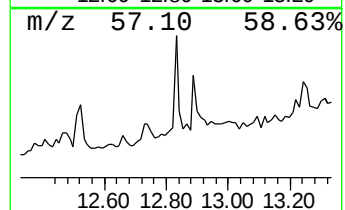
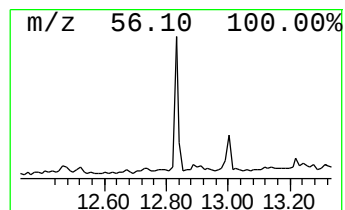
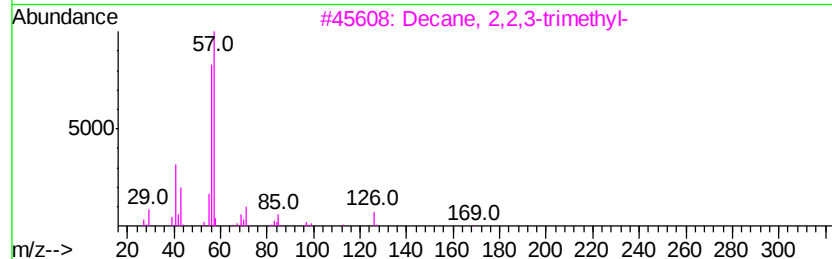
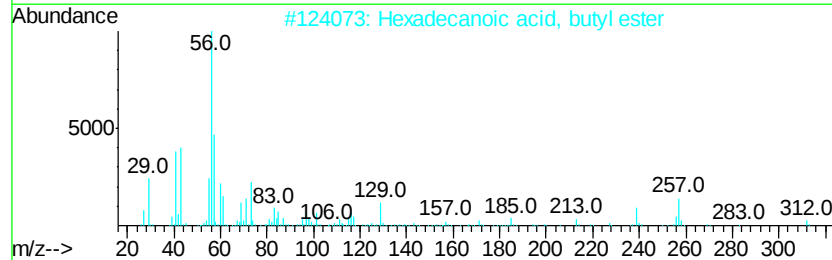
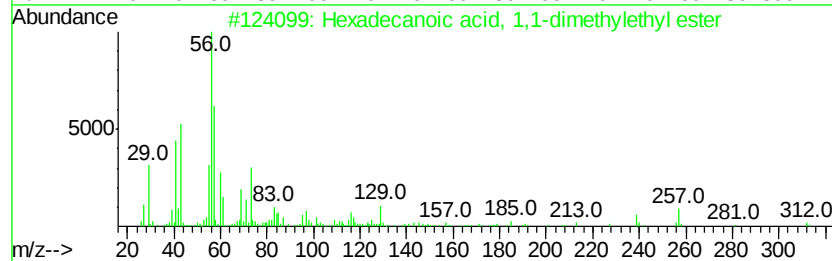
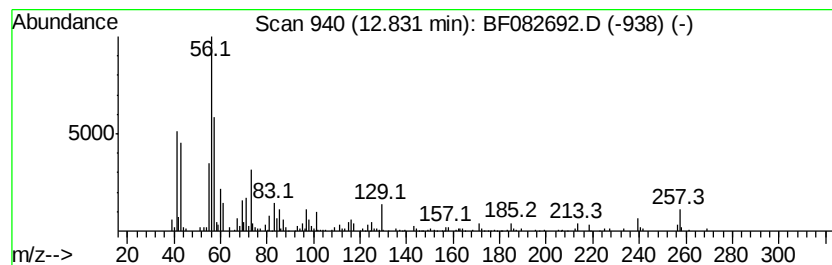
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Hexadecanoic acid, 1,1-dime... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.83	4.29 ng	224479	Chrysene-d12	14.05

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	90
2		Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	86
3		Decane, 2,2,3-trimethyl-	184	C13H28	062338-09-4	38
4		1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	38
5		2-Methyl-1-octadecene	266	C19H38	061868-20-0	38



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110415\
Data File : BF082692.D
Acq On : 4 Nov 2015 9:11
Operator : UM/IZ
Sample : G4241-11
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SB-12(2-4)

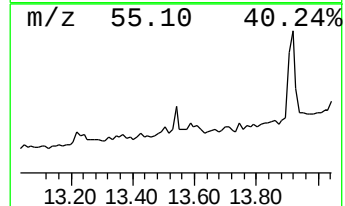
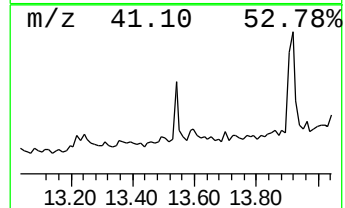
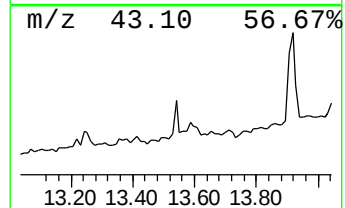
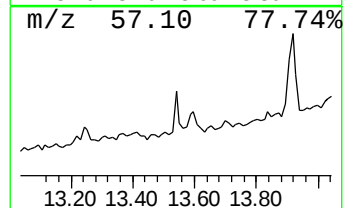
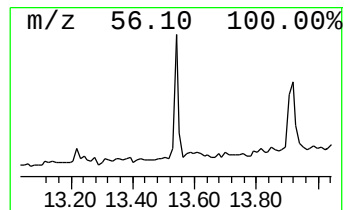
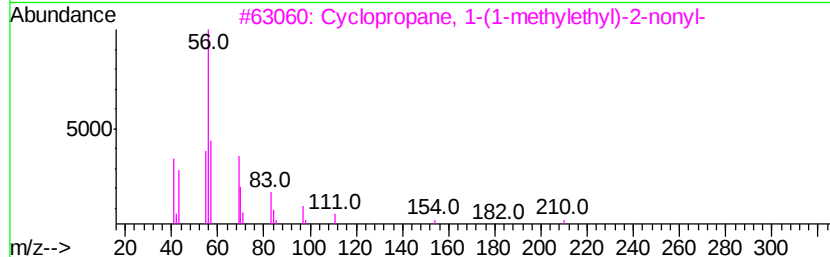
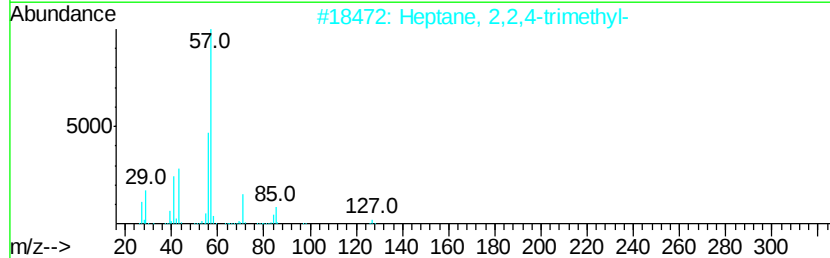
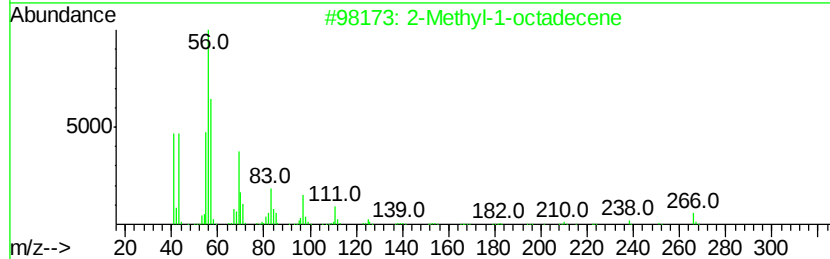
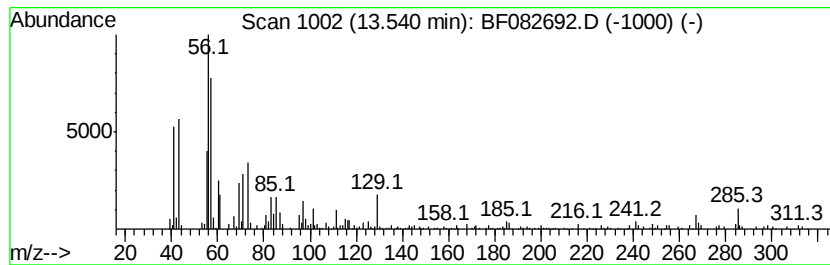
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown13.54 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	3.44 ng	180123	Chrysene-d12	14.05

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methyl-1-octadecene	266	C19H38	061868-20-0	46
2			Heptane, 2,2,4-trimethyl-	142	C10H22	014720-74-2	43
3			Cyclopropane, 1-(1-methylethyl)-...	210	C15H30	041977-39-3	43
4			1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	40
5			Cyclopentane, (1,1-dimethylethyl)-	126	C9H18	003875-52-3	35



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110415\
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Sample : G4241-11
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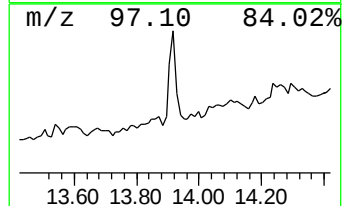
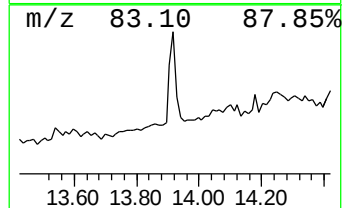
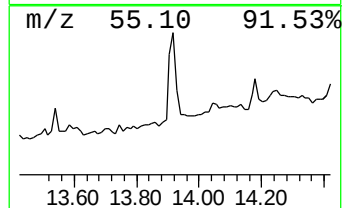
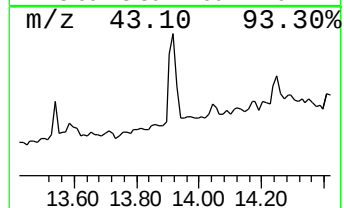
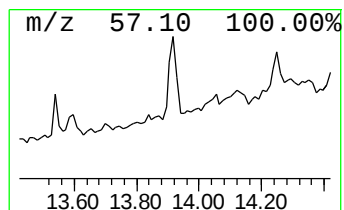
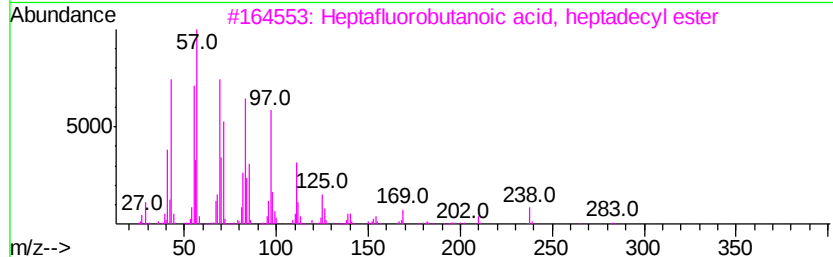
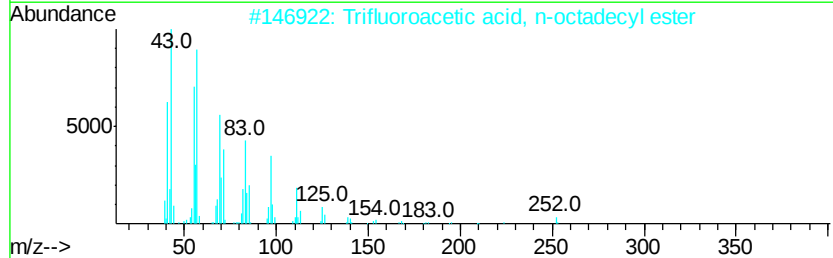
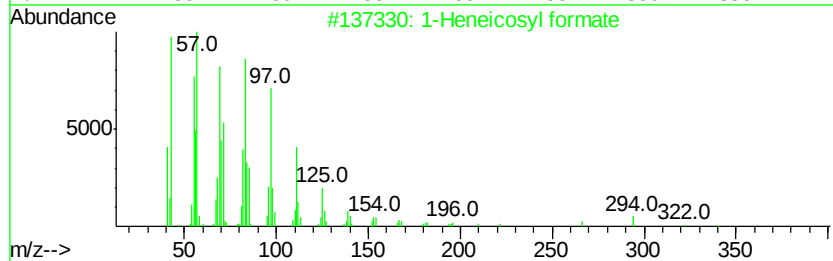
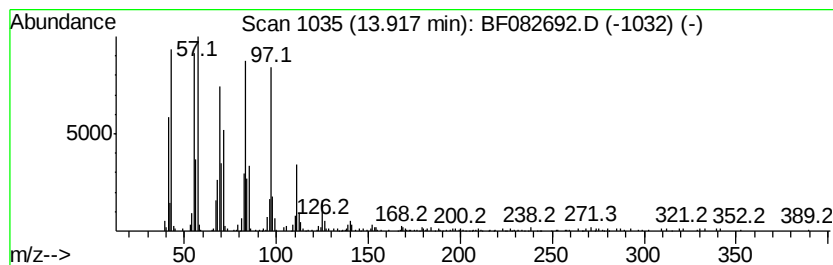
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1-Heneicosyl formate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.92	10.07 ng	527325	Chrysene-d12	14.05		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosyl formate	340	C22H44O2	077899-03-7	93
2		Trifluoroacetic acid, n-octadecyl...	366	C20H37F3O2	079392-43-1	91
3		Heptafluorobutanoic acid, heptadecyl...	452	C21H35F7O2	1000282-97-3	91
4		Cyclopentadecane	210	C15H30	000295-48-7	89
5		2-Chloropropionic acid, octadecyl...	360	C21H41ClO2	088104-31-8	87



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110415\
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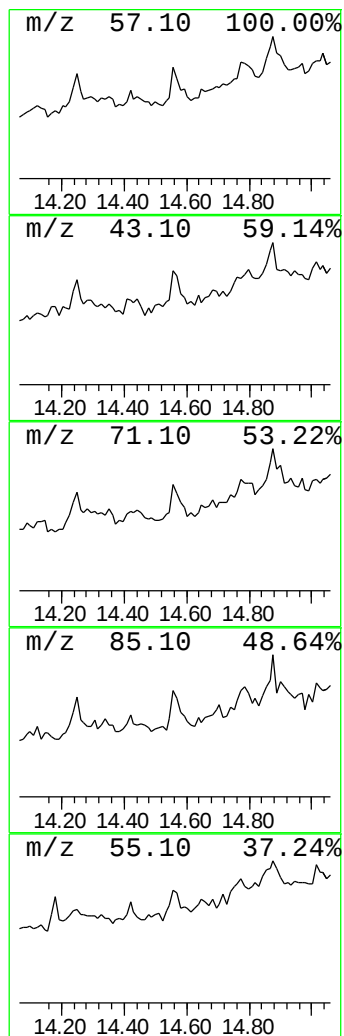
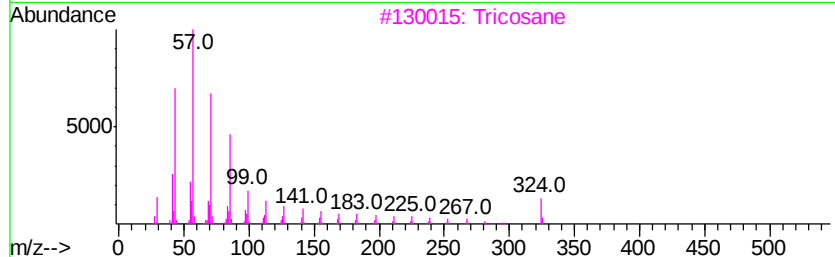
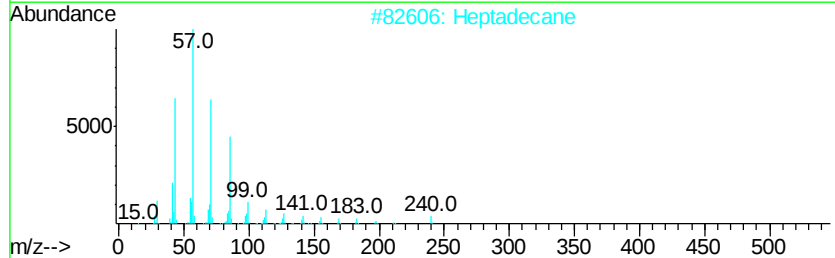
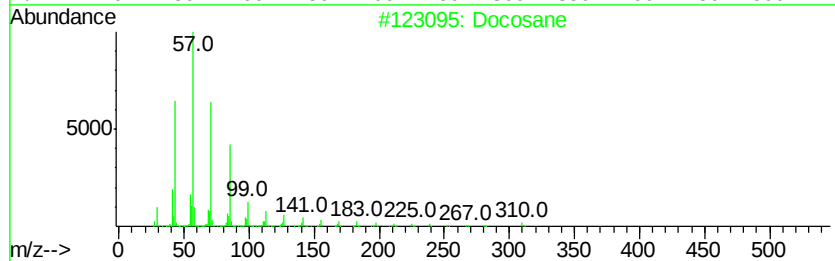
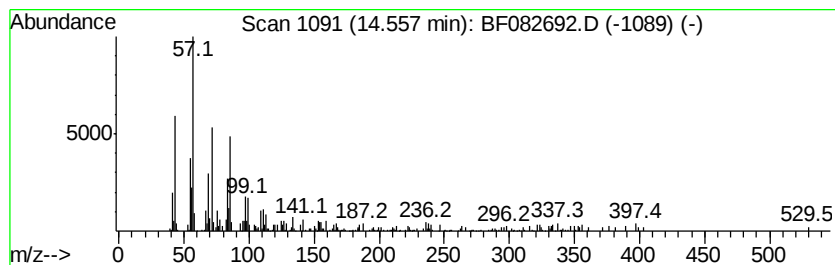
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Docosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.56	3.82 ng	200186	Chrysene-d12	14.05
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Docosane		310 C22H46	000629-97-0 92
2	Heptadecane		240 C17H36	000629-78-7 83
3	Tricosane		324 C23H48	000638-67-5 64
4	Pentadecane		212 C15H32	000629-62-9 58
5	Dodecane, 2-methyl-		184 C13H28	001560-97-0 58



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110415\
Data File : BF082692.D
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ALS Vial : 36 Sample Multiplier: 1

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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF103015.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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unknown6.65	6.65	100.3	ng	5337360	1	6.89	1064510	20.0
unknown10.34	10.34	2.0	ng	142626	3	9.93	1397150	20.0
Tetradecane	10.84	3.4	ng	201018	4	11.41	1198650	20.0
Cyclotetradecane	11.64	3.0	ng	180970	4	11.41	1198650	20.0
n-Hexadecanoic acid	11.95	5.6	ng	334637	4	11.41	1198650	20.0
unknown12.74	12.74	2.1	ng	110012	5	14.05	1047130	20.0
Hexadecanoic acid...	12.83	4.3	ng	224479	5	14.05	1047130	20.0
unknown13.54	13.54	3.4	ng	180123	5	14.05	1047130	20.0
1-Heneicosyl formate	13.92	10.1	ng	527325	5	14.05	1047130	20.0
Docosane	14.56	3.8	ng	200186	5	14.05	1047130	20.0