

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112515\
 Data File : BF083293.D
 Acq On : 26 Nov 2015 00:45
 Operator : UM/IZ
 Sample : G4569-06
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IM-TOPSOIL-001

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-BF112115.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	2.170	15	16	31	rVB3	33145	102174	1.93%	0.230%
2	4.718	236	239	254	rVB	1211769	1968343	37.15%	4.433%
3	5.187	278	280	293	rBV	4356100	4746021	89.56%	10.689%
4	6.261	371	374	376	rBV	4843455	5299004	100.00%	11.934%
5	6.364	381	383	385	rBV	5197593	4979275	93.97%	11.214%
6	6.582	400	402	405	rBV	828544	998181	18.84%	2.248%
7	6.742	414	416	419	rBV	3710585	3578428	67.53%	8.059%
8	7.164	450	453	455	rBV	2408593	3189742	60.20%	7.184%
9	7.873	513	515	518	rBV	1367903	1244418	23.48%	2.803%
10	8.959	607	610	612	rBV	5279253	4880140	92.10%	10.991%
11	9.359	643	645	647	rBV	765633	669155	12.63%	1.507%
12	9.576	662	664	666	rBV	79711	83835	1.58%	0.189%
13	9.622	666	668	671	rVB	1439279	1283573	24.22%	2.891%
14	10.079	701	708	709	rBV	153360	168404	3.18%	0.379%
15	10.296	724	727	730	rBV	53481	90677	1.71%	0.204%
16	10.410	735	737	739	rBV	2439908	2350187	44.35%	5.293%
17	10.548	747	749	753	rBV	178976	215615	4.07%	0.486%
18	10.993	783	788	789	rBV	91873	96477	1.82%	0.217%
19	11.096	794	797	799	rVV	1253871	1120682	21.15%	2.524%
20	11.153	801	802	808	rVB2	52002	79316	1.50%	0.179%
21	11.656	844	846	849	rBV2	156888	207573	3.92%	0.467%
22	12.434	912	914	919	rBV	59306	76047	1.44%	0.171%
23	12.685	933	936	938	rBV	2996589	3148271	59.41%	7.090%
24	13.588	1013	1015	1019	rBV	191347	284430	5.37%	0.641%
25	13.714	1024	1026	1029	rBV	914795	1109108	20.93%	2.498%
26	14.228	1068	1071	1074	rBV	86694	151368	2.86%	0.341%
27	14.445	1086	1090	1092	rBV4	23510	56271	1.06%	0.127%
28	14.799	1119	1121	1128	rBV2	134896	234271	4.42%	0.528%
29	15.074	1142	1145	1147	rBV	682676	954725	18.02%	2.150%
30	15.291	1162	1164	1167	rVB	67338	82747	1.56%	0.186%
31	15.485	1178	1181	1183	rBV	140393	201002	3.79%	0.453%
32	15.531	1183	1185	1189	rVB2	58295	103665	1.96%	0.233%
33	16.137	1235	1238	1241	rVB	72758	125566	2.37%	0.283%
34	16.388	1257	1260	1262	rBV	37239	71102	1.34%	0.160%

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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-BF112115.M
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35	16.765	1290	1293	1298	rBV6	49971	126947	2.40%	0.286%
36	16.937	1306	1308	1314	rVB4	40961	97112	1.83%	0.219%
37	17.611	1363	1367	1371	rVB5	47667	147183	2.78%	0.331%
38	18.503	1442	1445	1449	rBV5	27630	81459	1.54%	0.183%

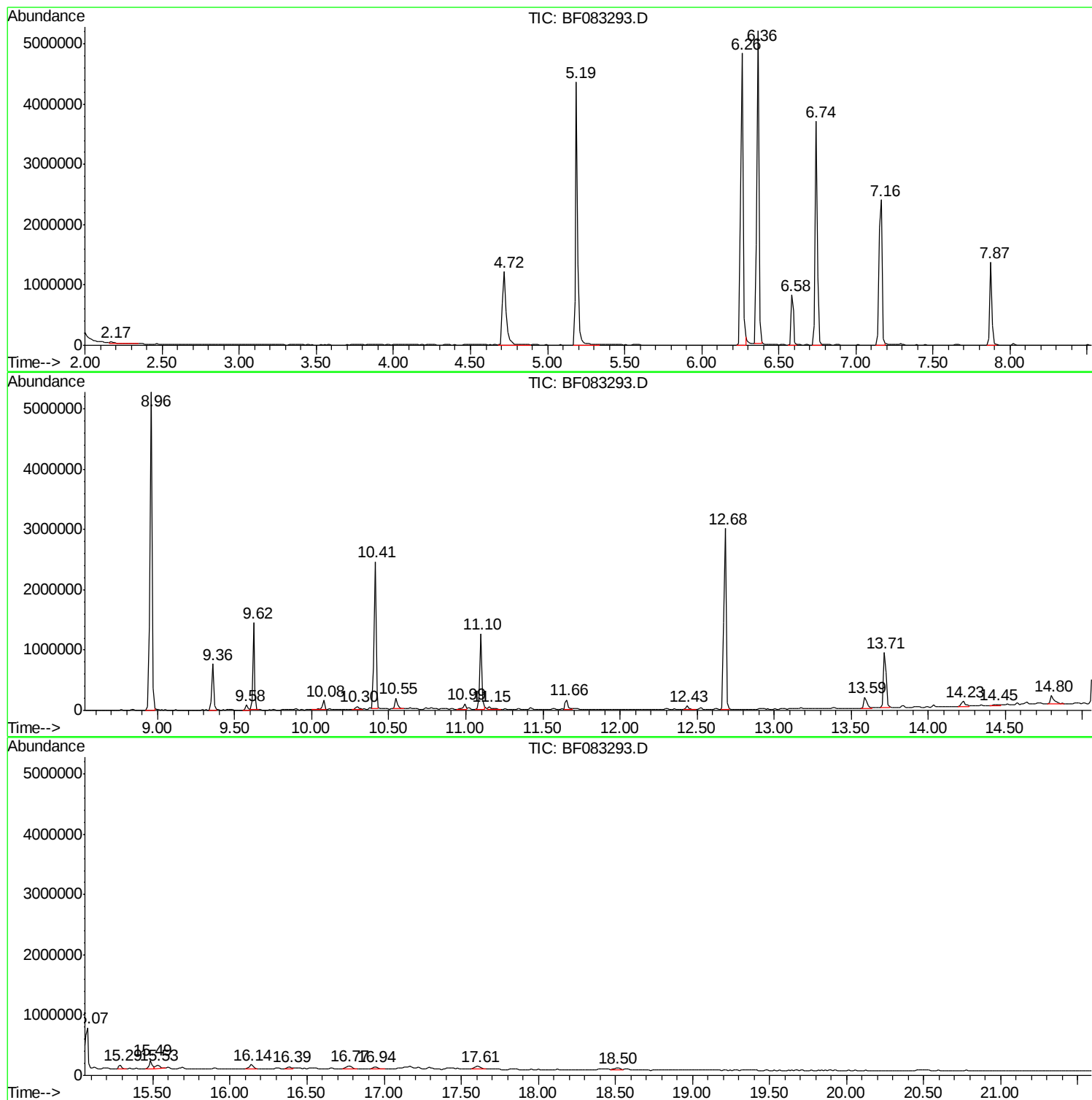
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Quant Method : Z:\HPCHEM1\BNA_F\Methods\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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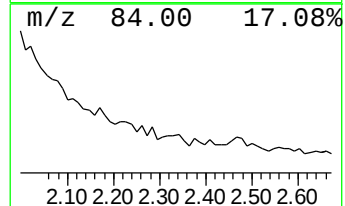
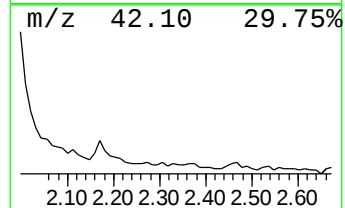
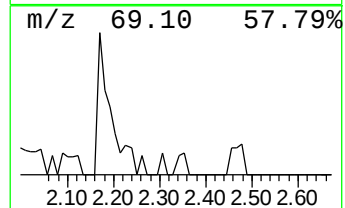
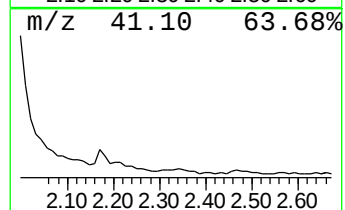
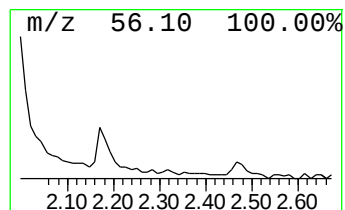
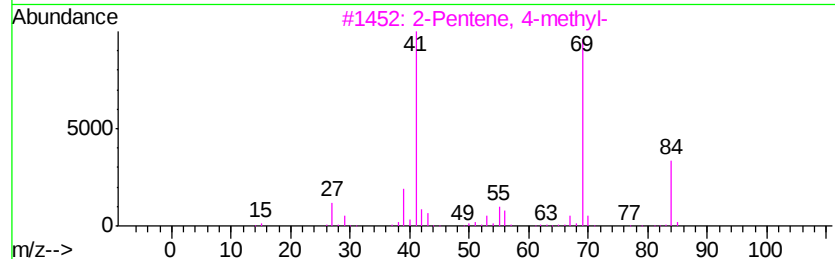
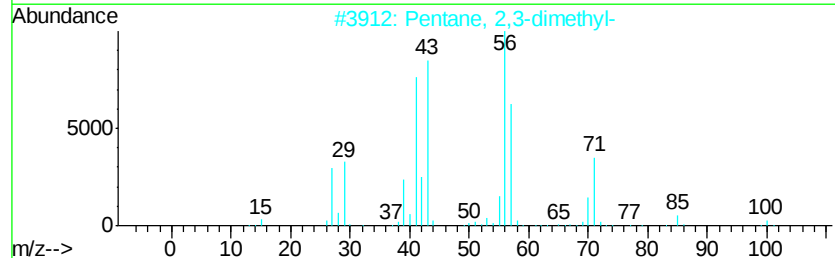
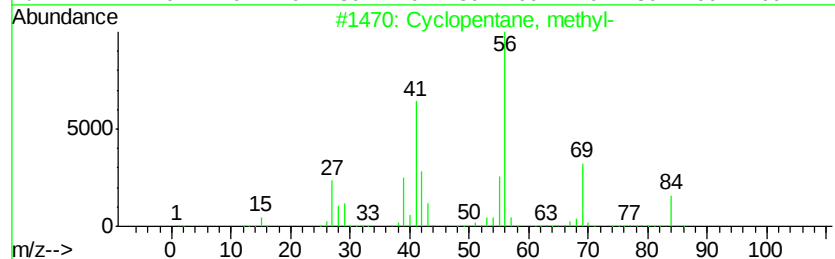
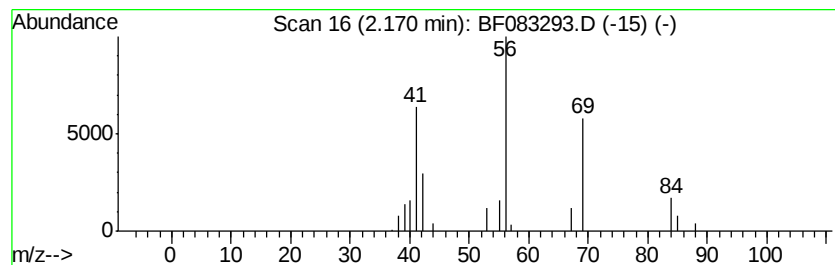
Quant Method : Z:\HPCHEM1\BNA_F\Methods\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown2.17 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.17	2.05 ng	102174	1,4-Dichlorobenzene-d4	6.58

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	38
2			Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	36
3			2-Pentene, 4-methyl-	84	C6H12	004461-48-7	9
4			2-Pentene, 4-methyl-, (Z)-	84	C6H12	000691-38-3	9
5			Cyclopropane, 1,2,3-trimethyl-	84	C6H12	042984-19-0	9



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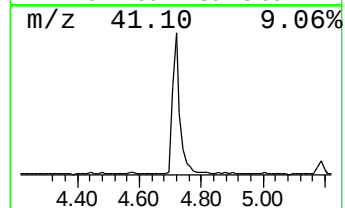
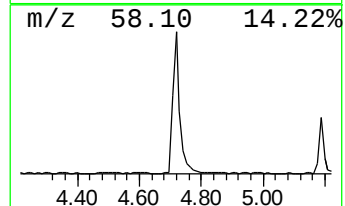
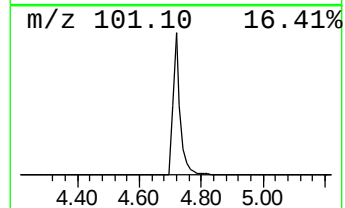
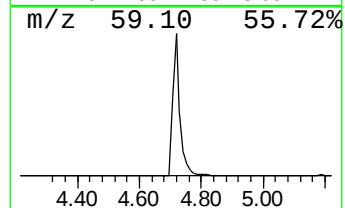
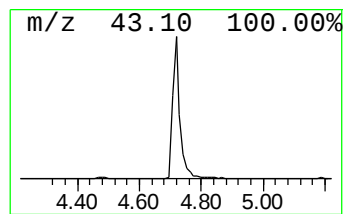
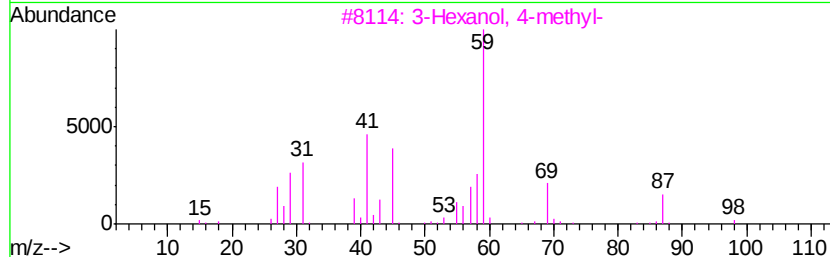
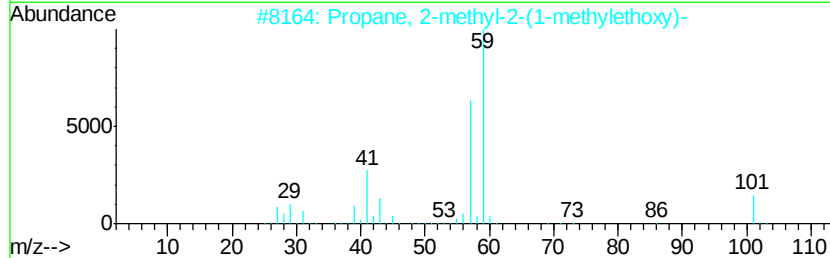
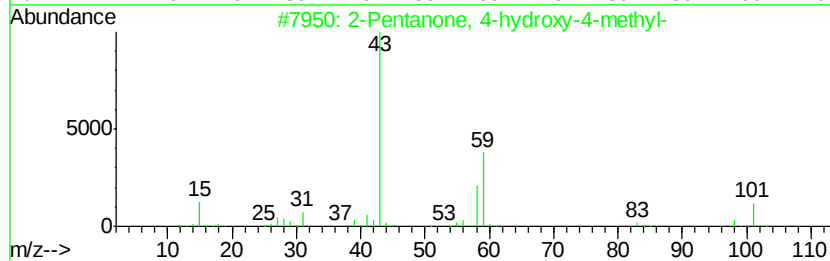
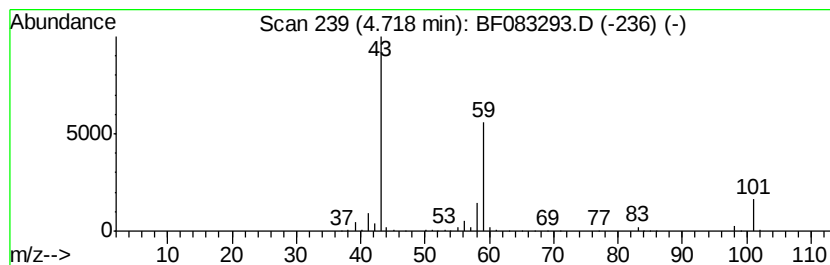
Quant Method : Z:\HPCHEM1\BNA_F\Methods\8270-BF112115.M
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.72	39.44 ng	1968340	1,4-Dichlorobenzene-d4	6.58

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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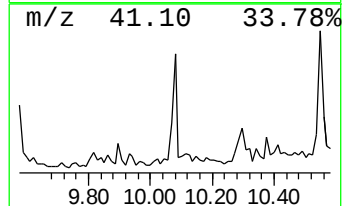
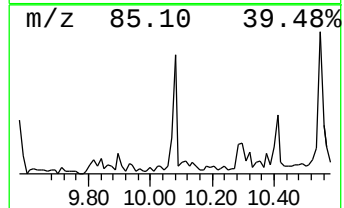
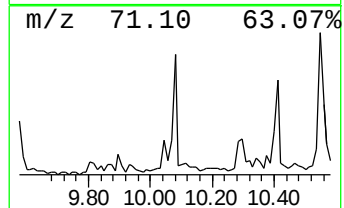
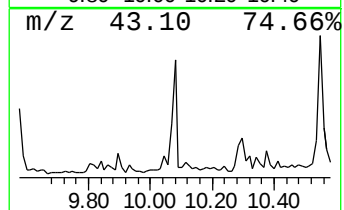
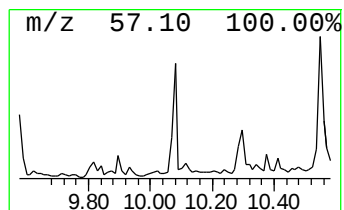
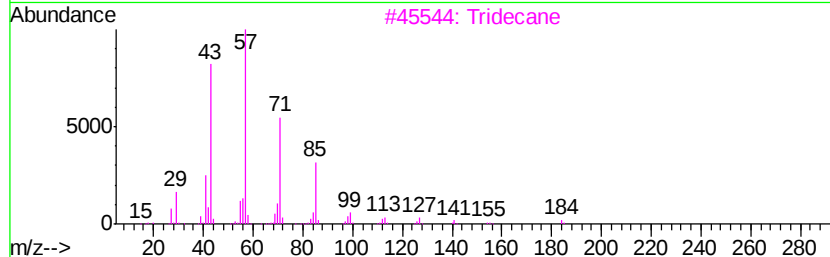
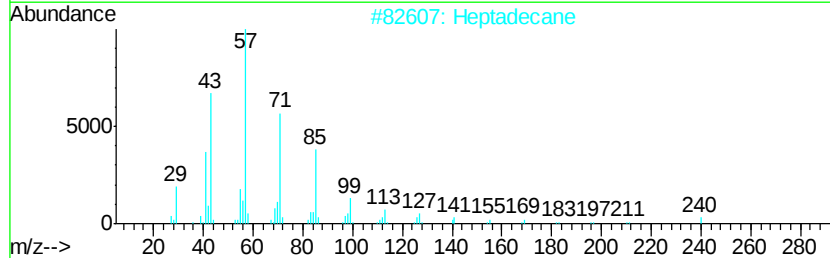
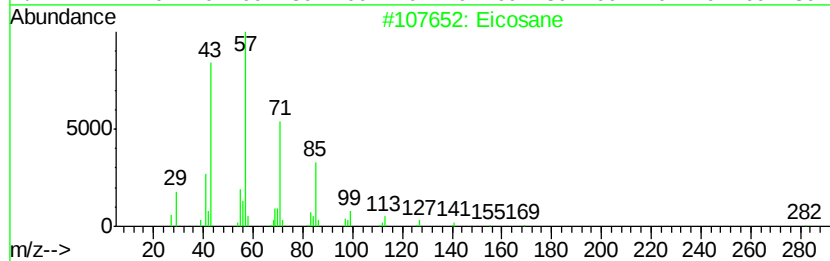
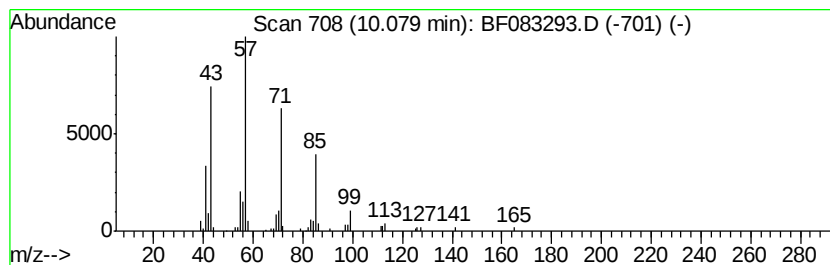
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Eicosane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.08	2.62 ng	168404	Acenaphthene-d10	9.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	91
2		Heptadecane	240	C17H36	000629-78-7	90
3		Tridecane	184	C13H28	000629-50-5	86
4		Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	83
5		Pentadecane	212	C15H32	000629-62-9	78



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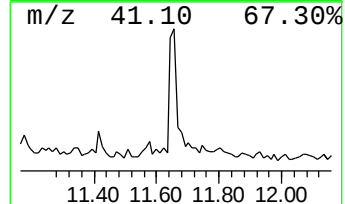
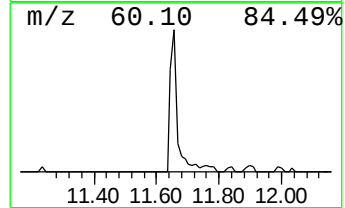
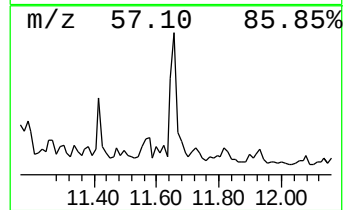
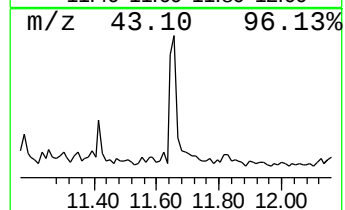
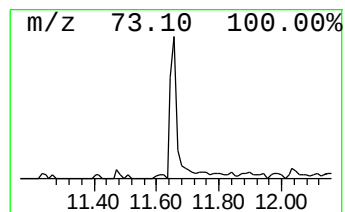
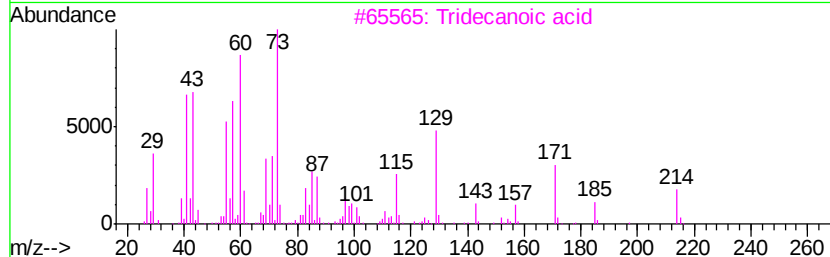
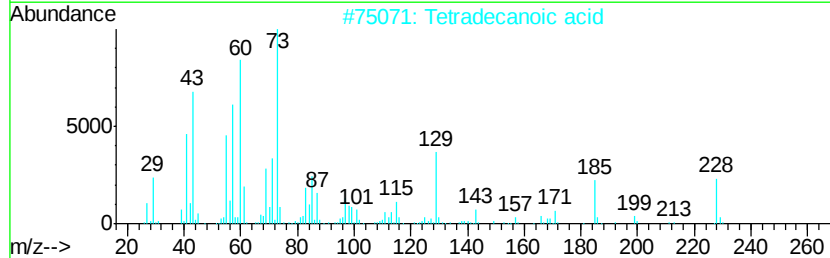
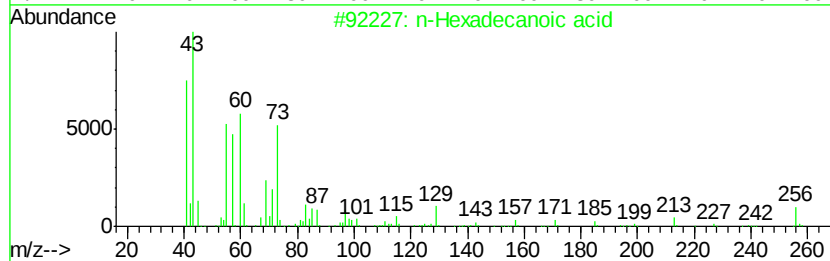
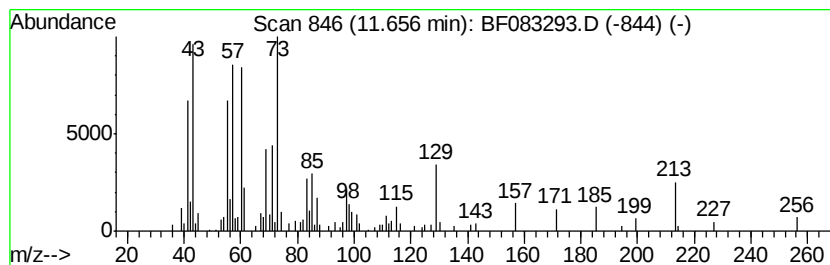
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	3.70 ng	207573	Phenanthrene-d10	11.10

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	80
3		Tridecanoic acid	214	C13H26O2	000638-53-9	72
4		Hexadecane	226	C16H34	000544-76-3	11
5		Methoxyacetic acid, 2-tetradecyl...	286	C17H34O3	1000282-04-8	11



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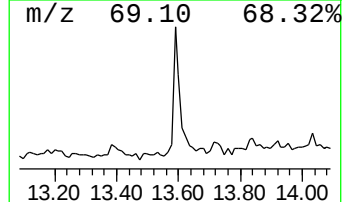
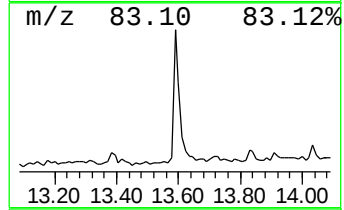
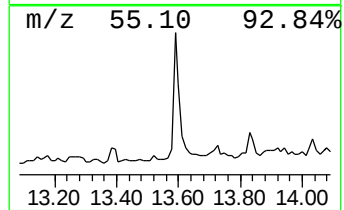
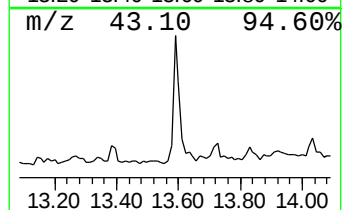
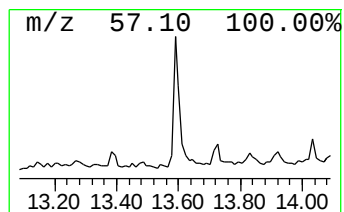
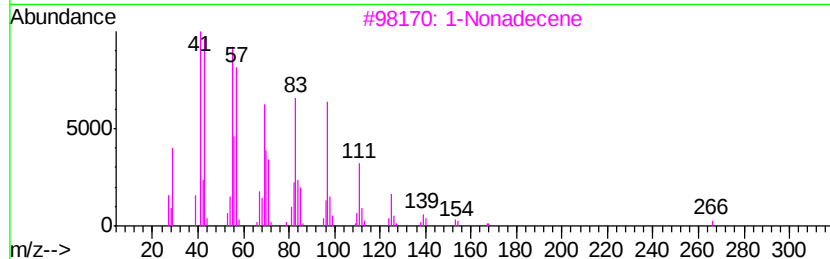
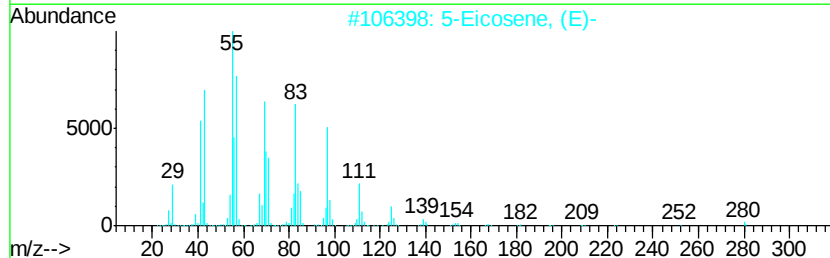
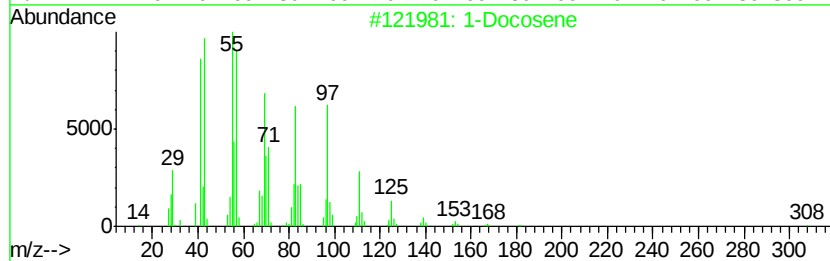
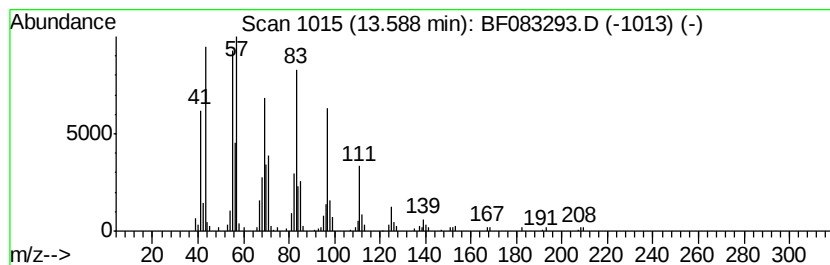
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1-Docosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.59	5.13 ng	284430	Chrysene-d12	13.71

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene		308	C22H44	001599-67-3	94
2	5-Eicosene, (E)-		280	C20H40	074685-30-6	91
3	1-Nonadecene		266	C19H38	018435-45-5	91
4	2-Bromopropionic acid, pentadec...		362	C18H35BrO2	1000292-44-2	91
5	11-Tricosene		322	C23H46	052078-56-5	91



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112515\
Data File : BF083293.D
Acq On : 26 Nov 2015 00:45
Operator : UM/IZ
Sample : G4569-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
IM-TOPSOIL-001

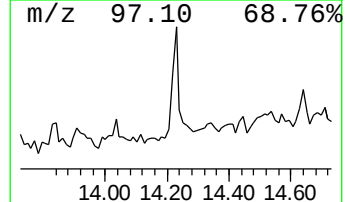
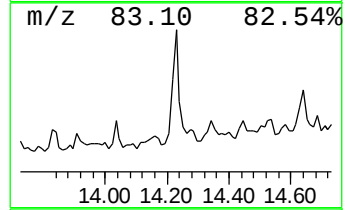
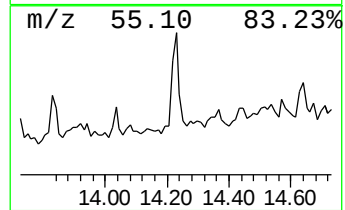
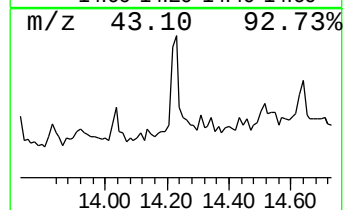
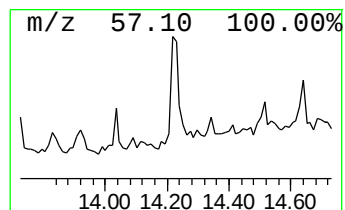
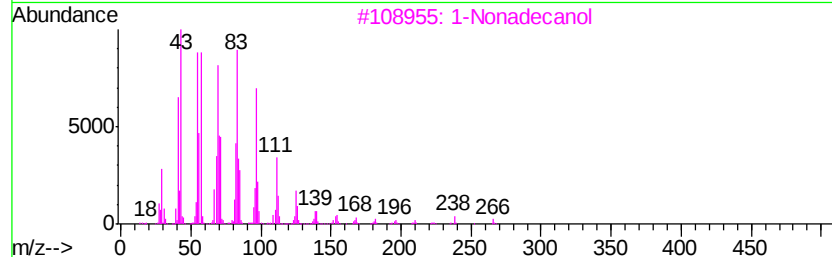
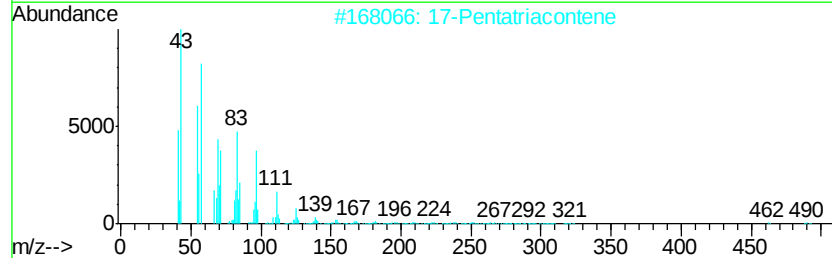
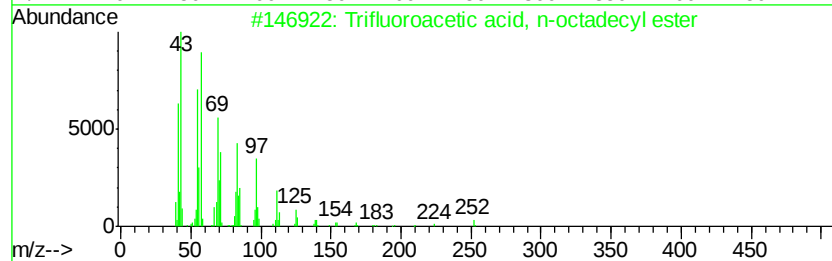
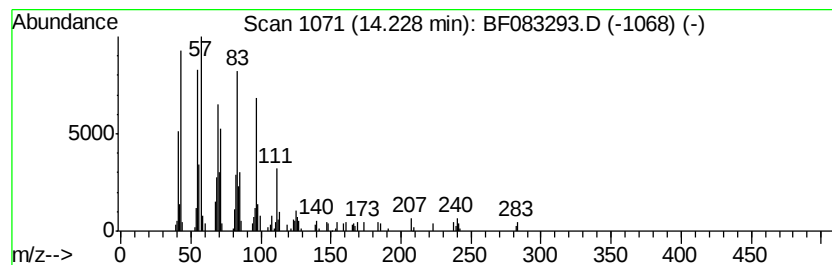
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Trifluoroacetic acid, n-oct... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.23	2.73 ng	151368	Chrysene-d12	13.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Trifluoroacetic acid, n-octadecv...	366	C20H37F3O2	079392-43-1	81
2		17-Pentatriacontene	491	C35H70	006971-40-0	80
3		1-Nonadecanol	284	C19H40O	001454-84-8	68
4		Heptafluorobutyric acid, n-octad...	466	C22H37F7O2	000400-57-7	68
5		Heptafluorobutyric acid, n-penta...	424	C19H31F7O2	1000216-79-3	64



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112515\
Data File : BF083293.D
Acq On : 26 Nov 2015 00:45
Operator : UM/IZ
Sample : G4569-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

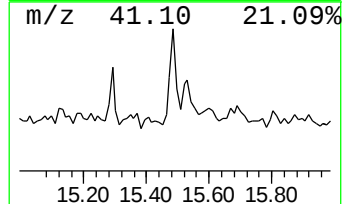
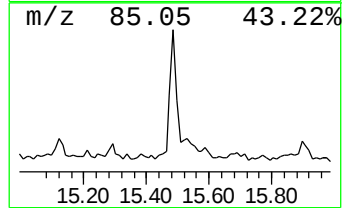
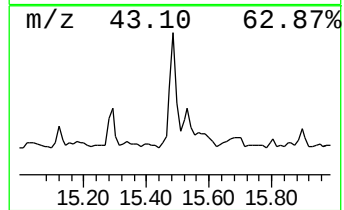
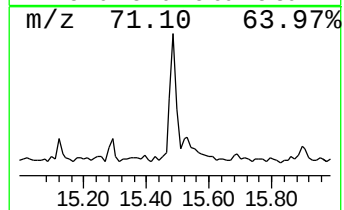
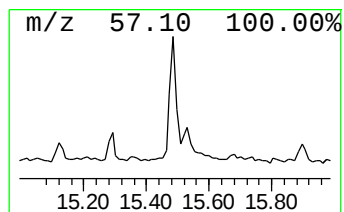
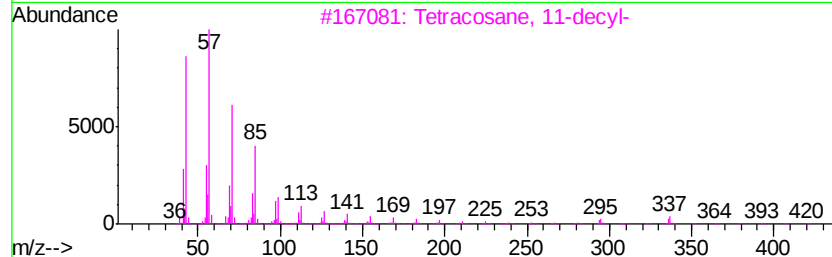
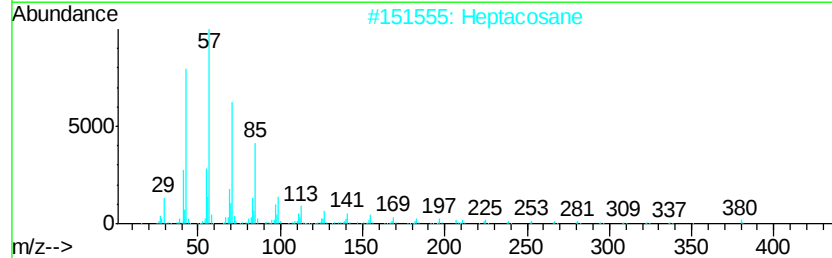
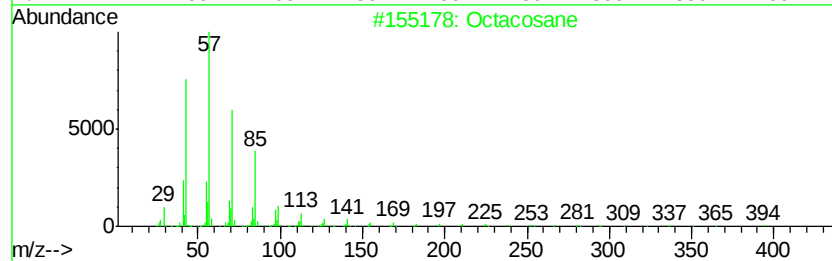
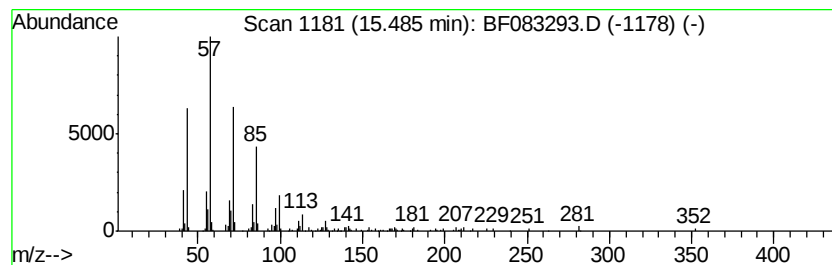
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ClientSampleId :
IM-TOPSOIL-001

Quant Method : Z:\HPCHEM1\BNA_F\Methods\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.49	4.21 ng	201002	Perylene-d12	15.07
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octacosane	394 C28H58	000630-02-4	90
2	Heptacosane	380 C27H56	000593-49-7	87
3	Tetracosane, 11-decyl-	479 C34H70	055429-84-0	86
4	Docosane, 9-butyl-	366 C26H54	055282-14-9	86
5	Nonacosane	408 C29H60	000630-03-5	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112515\
Data File : BF083293.D
Acq On : 26 Nov 2015 00:45
Operator : UM/IZ
Sample : G4569-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
IM-TOPSOIL-001

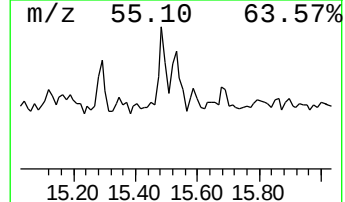
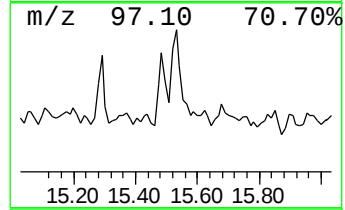
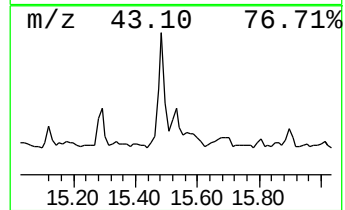
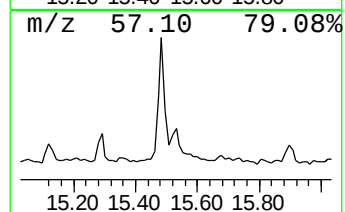
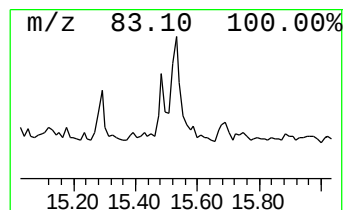
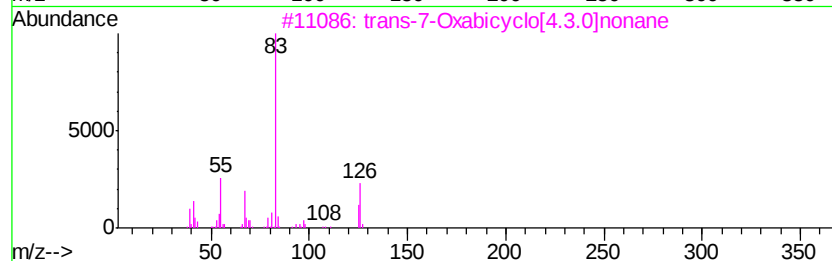
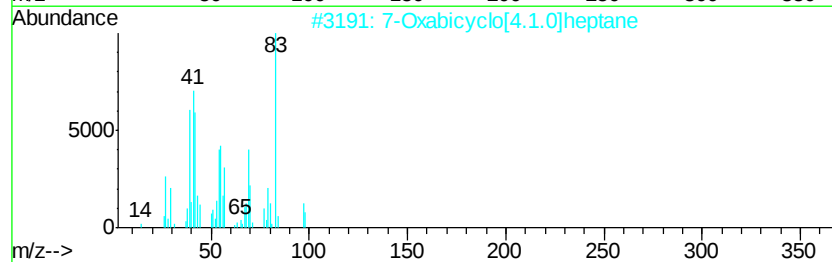
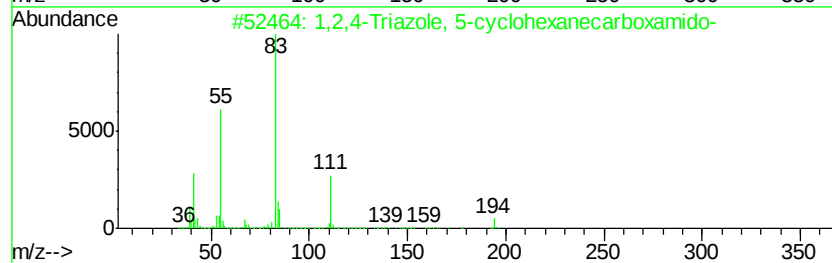
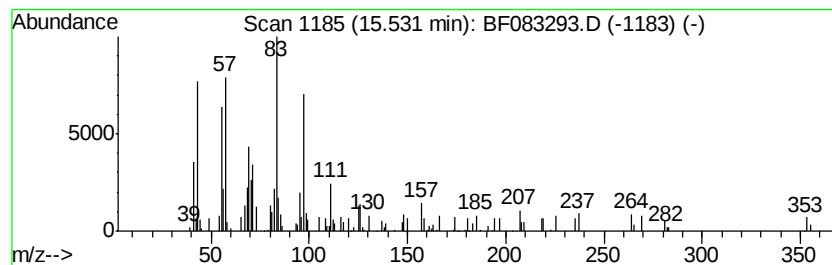
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown15.53 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.53	2.17 ng	103665	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,2,4-Triazole, 5-cvclohexanecar...	194	C9H14N4O	021051-17-2	43
2		7-Oxabicyclo[4.1.0]heptane	98	C6H10O	000286-20-4	38
3		trans-7-Oxabicyclo[4.3.0]nonane	126	C8H14O	027345-70-6	38
4		1-Tetracosanol	354	C24H50O	000506-51-4	30
5		Ethanone, 1-(1-methylcyclopentyl)-	126	C8H14O	013388-93-7	30



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112515\
Data File : BF083293.D
Acq On : 26 Nov 2015 00:45
Operator : UM/IZ
Sample : G4569-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

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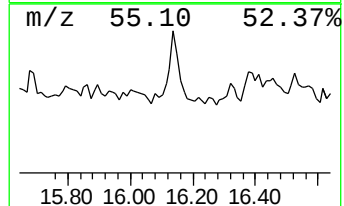
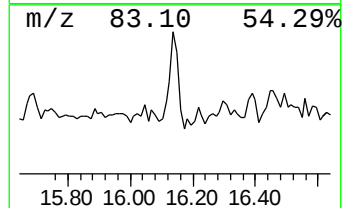
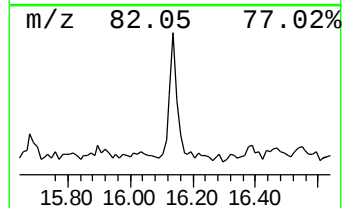
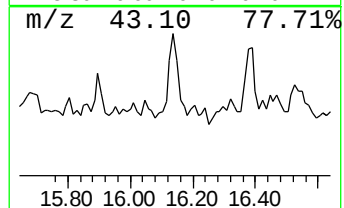
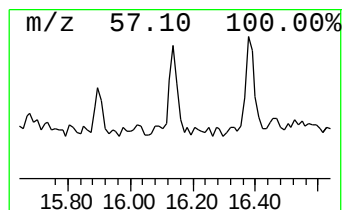
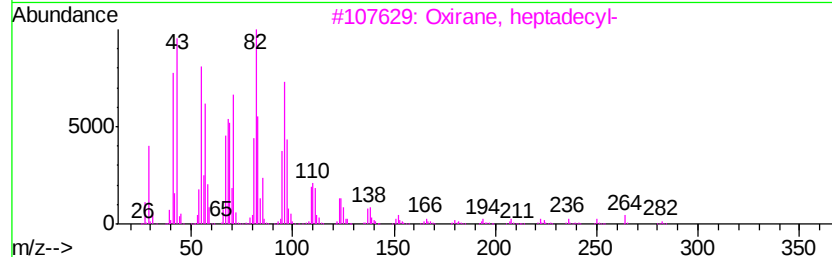
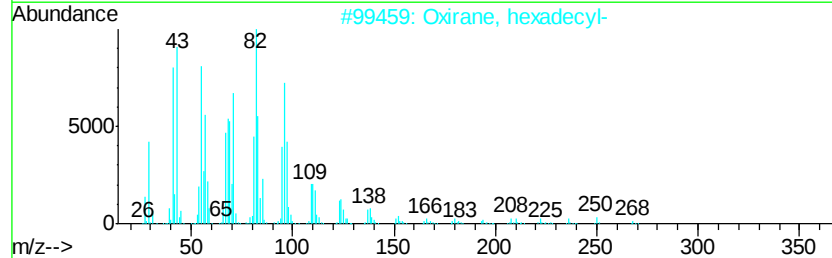
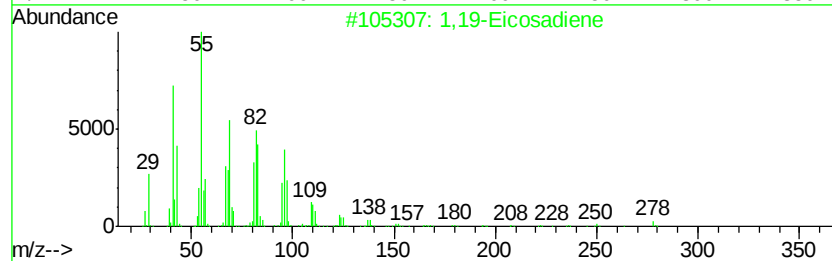
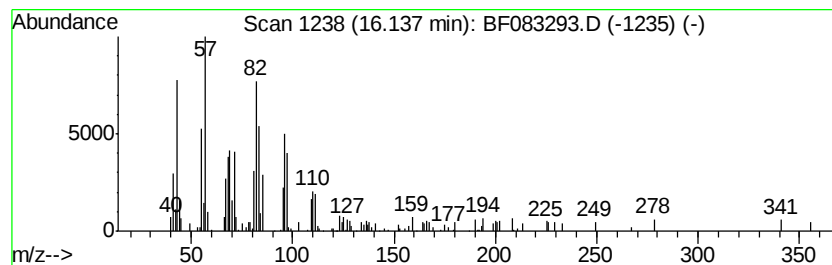
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 1,19-Eicosadiene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.14	2.63 ng	125566	Perylene-d12	15.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,19-Eicosadiene	278	C20H38	014811-95-1	76
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	72
3		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	72
4		Octadecanal	268	C18H36O	000638-66-4	72
5		Z-2-Octadecen-1-ol	268	C18H36O	1000131-11-0	53



Data Path : Z:\HPCHEM1\BNA F\DATA\BF112515\
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Acq On : 26 Nov 2015 00:45
Operator : UM/IZ
Sample : G4569-06
Misc :
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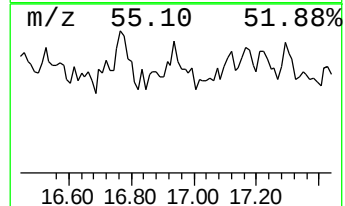
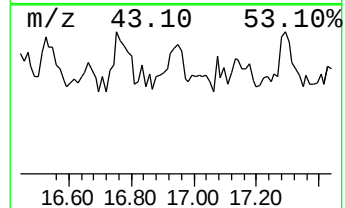
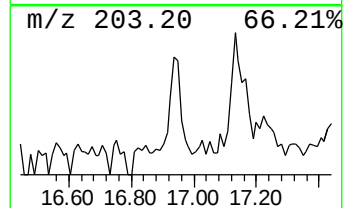
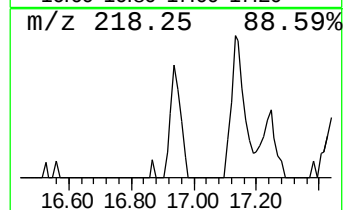
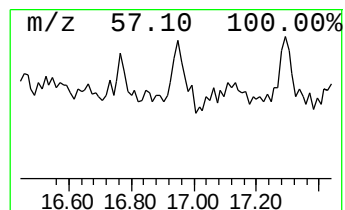
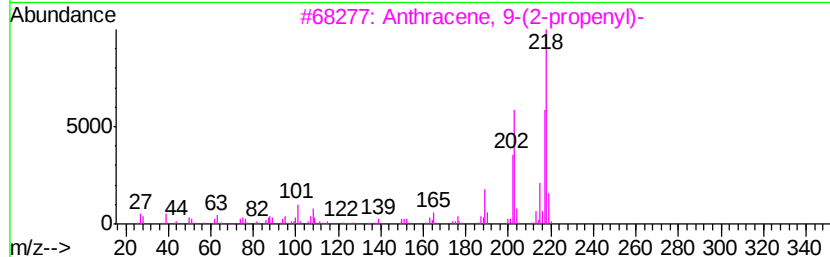
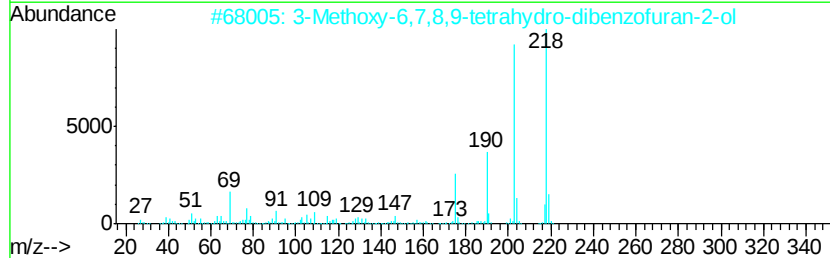
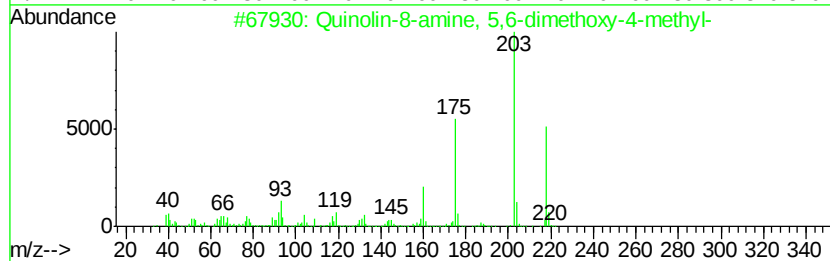
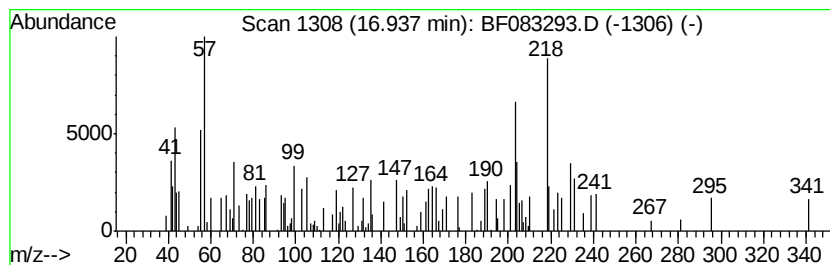
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 unknown16.94 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.94	2.03 ng	97112	Perylene-d12		15.07	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Quinolin-8-amine, 5,6-dimethoxy-	218	C12H14N2O2	064992-95-6	46
2		3-Methoxy-6,7,8,9-tetrahydro-dibenz	218	C13H14O3	1000186-57-9	43
3		Anthracene, 9-(2-propenyl)-	218	C17H14	023707-65-5	43
4		Pyrrolidino[3,2-b]benzo[e]-2H-1,4-	218	C11H10N2OS	017163-68-7	38
5		2,3-Diethoxyquinoxaline	218	C12H14N2O2	057050-66-5	37



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112515\
Data File : BF083293.D
Acq On : 26 Nov 2015 00:45
Operator : UM/IZ
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Misc :
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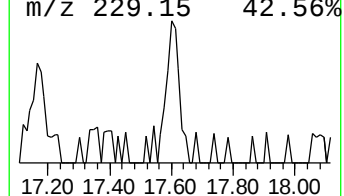
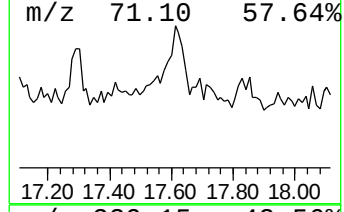
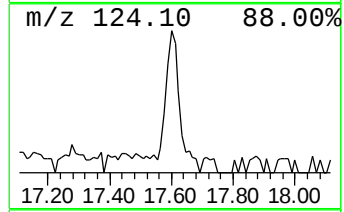
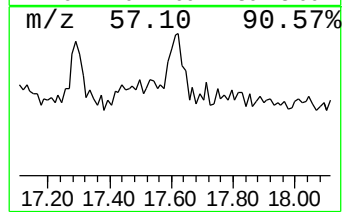
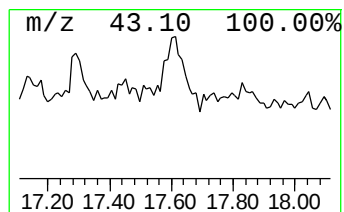
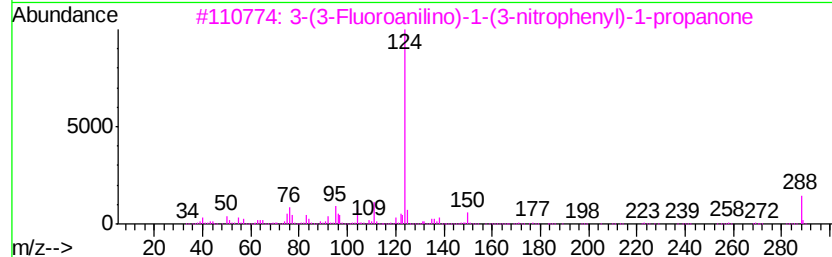
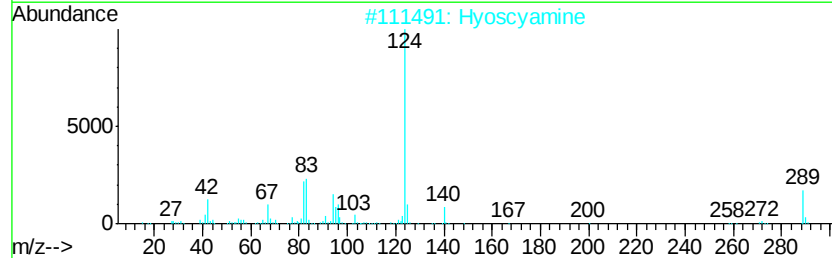
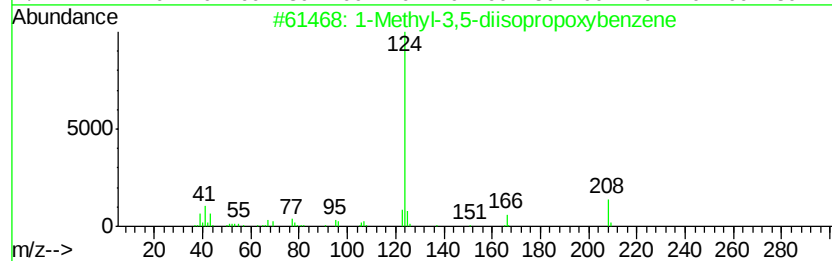
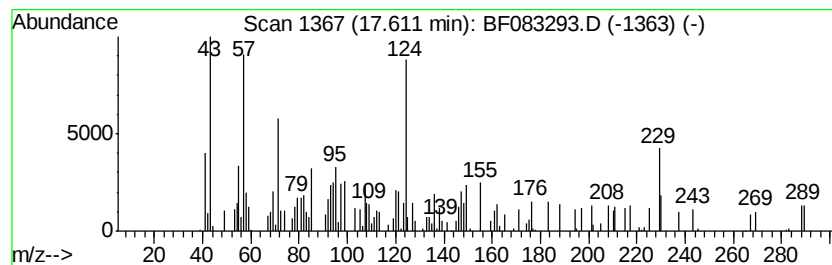
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown17.61 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.61	3.08 ng	147183	Perylene-d12	15.07

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methyl-3,5-diisopropoxybenzene	208	C13H20O2	1000281-22-2	38
2			Hyoscyamine	289	C17H23NO3	000101-31-5	38
3			3-(3-Fluoroanilino)-1-(3-nitroph...	288	C15H13FN2O3	300726-64-1	35
4			Atropine	289	C17H23NO3	000051-55-8	35
5			1,4-Benzenediol, 2-methyl-	124	C7H8O2	000095-71-6	30



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF112515\
Data File : BF083293.D
Acq On : 26 Nov 2015 00:45
Operator : UM/IZ
Sample : G4569-06
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_F\Methods\8270-BF112115.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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2-Pentanone, 4-hy...	4.72	39.4 ng		1968340	1	6.58	998181	20.0
Eicosane	10.08	2.6 ng		168404	3	9.62	1283570	20.0
n-Hexadecanoic acid	11.66	3.7 ng		207573	4	11.10	1120680	20.0
1-Docosene	13.59	5.1 ng		284430	5	13.71	1109110	20.0
Trifluoroacetic a...	14.23	2.7 ng		151368	5	13.71	1109110	20.0
Octacosane	15.49	4.2 ng		201002	6	15.07	954725	20.0
unknown15.53	15.53	2.2 ng		103665	6	15.07	954725	20.0
1,19-Eicosadiene	16.14	2.6 ng		125566	6	15.07	954725	20.0
unknown16.94	16.94	2.0 ng		97112	6	15.07	954725	20.0
unknown17.61	17.61	3.1 ng		147183	6	15.07	954725	20.0