

Data Path : Z:\HPCHEM1\BNA F\Data\BF040618\
 Data File : BF104320.D
 Acq On : 6 Apr 2018 21:56
 Operator : JU/SJ
 Sample : J2232-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CV-RB26667RT

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-BF040618.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	5.022	489	499	507	rVB	131315	182135	5.21%	0.556%
2	5.416	560	566	569	rBV	3083329	3383335	96.70%	10.324%
3	6.434	732	739	742	rBV	3079414	3444893	98.46%	10.512%
4	6.569	757	762	766	rBV	3146883	3480104	99.47%	10.620%
5	6.804	798	802	805	rVB	1155118	926191	26.47%	2.826%
6	6.963	824	829	832	rVB	2753605	2663573	76.13%	8.128%
7	7.369	892	898	901	rBV	2167990	2127722	60.81%	6.493%
8	8.086	1013	1020	1023	rBV	1559000	1287849	36.81%	3.930%
9	8.628	1108	1112	1114	rBV3	87989	122182	3.49%	0.373%
10	8.710	1123	1126	1128	rBV4	106165	117122	3.35%	0.357%
11	8.833	1143	1147	1149	rBV4	111038	158029	4.52%	0.482%
12	8.969	1167	1170	1172	rBV3	160529	179623	5.13%	0.548%
13	9.080	1187	1189	1191	rVB2	113235	71668	2.05%	0.219%
14	9.110	1191	1194	1198	rBV	200553	195271	5.58%	0.596%
15	9.163	1198	1203	1207	rBV	3243317	3498775	100.00%	10.677%
16	9.245	1213	1217	1219	rBV3	241974	331331	9.47%	1.011%
17	9.557	1267	1270	1274	rBV2	456319	600368	17.16%	1.832%
18	9.592	1274	1276	1280	rBV4	271285	405538	11.59%	1.238%
19	9.710	1293	1296	1299	rBV4	332150	540191	15.44%	1.648%
20	9.845	1316	1319	1323	rVB2	2012239	2116941	60.51%	6.460%
21	12.933	1841	1844	1853	rVB	1948699	2367091	67.65%	7.223%
22	13.810	1990	1993	2007	rVB	667397	780571	22.31%	2.382%
23	13.974	2016	2021	2024	rBV	993297	993706	28.40%	3.032%
24	13.998	2024	2025	2033	rVB	141147	99661	2.85%	0.304%
25	14.457	2099	2103	2110	rBV	484566	529075	15.12%	1.614%
26	14.904	2176	2179	2186	rVB4	48981	58797	1.68%	0.179%
27	15.004	2193	2196	2204	rVB	89362	166180	4.75%	0.507%
28	15.092	2207	2211	2213	rBV	91583	118896	3.40%	0.363%
29	15.115	2213	2215	2219	rVB	130941	128113	3.66%	0.391%
30	15.304	2243	2247	2252	rVB	52839	70255	2.01%	0.214%
31	15.362	2254	2257	2260	rVB	72825	76437	2.18%	0.233%
32	15.427	2263	2268	2272	rBV	703365	877781	25.09%	2.679%
33	15.898	2344	2348	2352	rBV	71195	110845	3.17%	0.338%
34	15.945	2352	2356	2366	rVB3	146120	237602	6.79%	0.725%

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Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	16.686	2477	2482	2490	rVB5	42235	96961	2.77%	0.296%
36	17.451	2606	2612	2618	rVB5	71056	149886	4.28%	0.457%
37	18.250	2743	2748	2754	rVB8	35538	75918	2.17%	0.232%

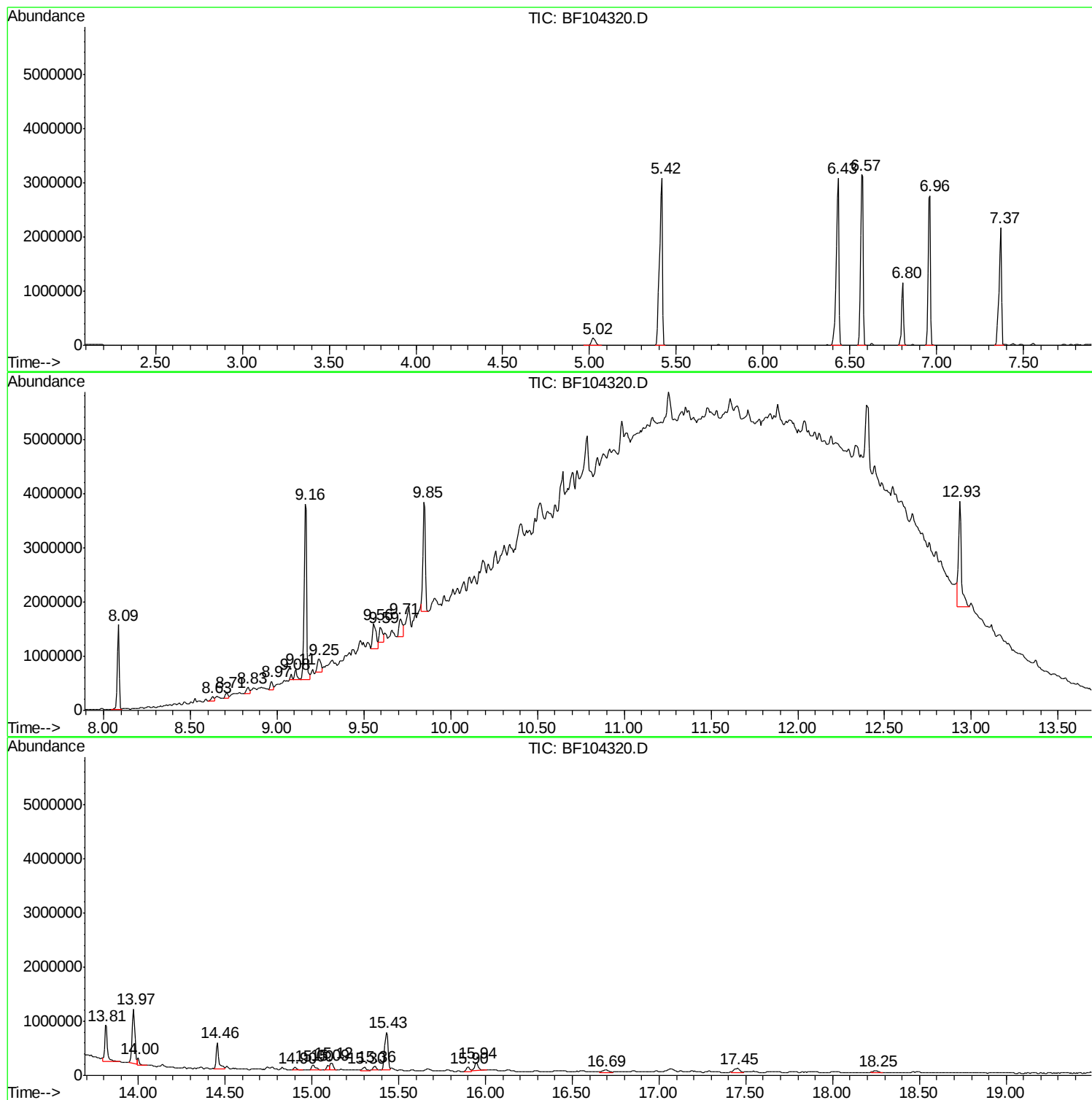
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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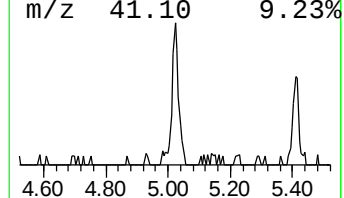
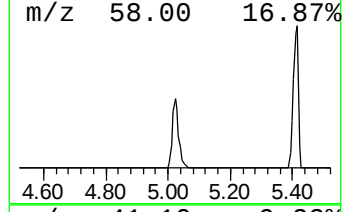
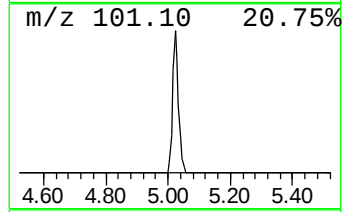
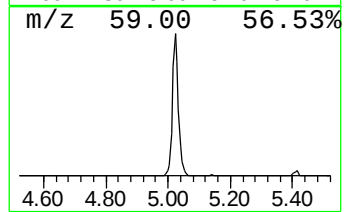
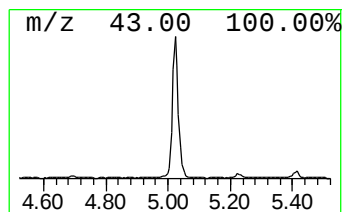
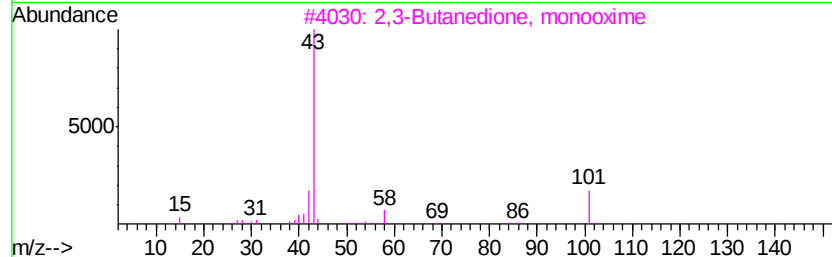
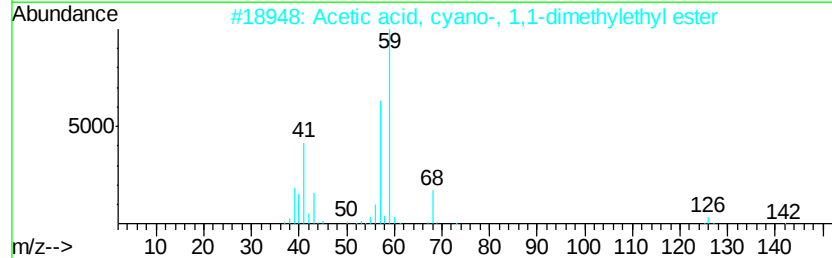
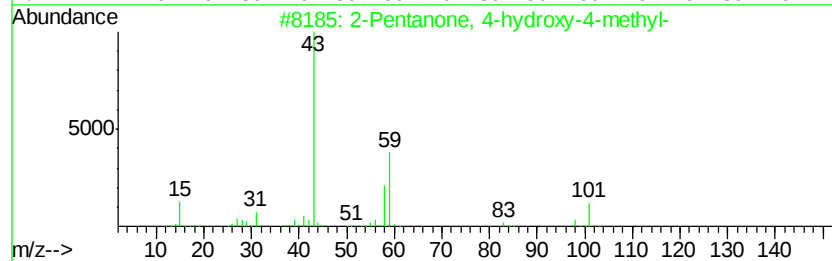
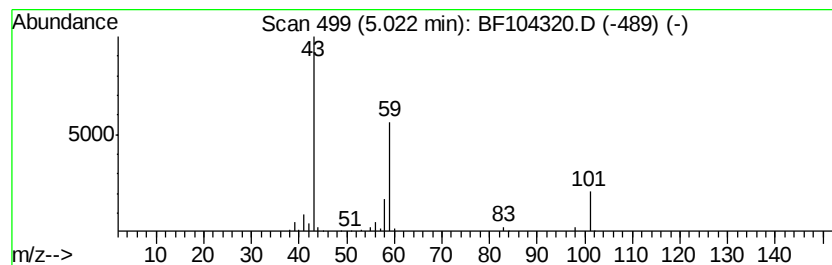
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	3.93 ng	182135	1,4-Dichlorobenzene-d4	6.80

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2		000123-42-2	50
2	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2		001116-98-9	32
3	2,3-Butanedione, monooxime	101	C4H7NO2		000057-71-6	25
4	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO		005780-40-5	17
5	5-Hexen-2-one	98	C6H10O		000109-49-9	9



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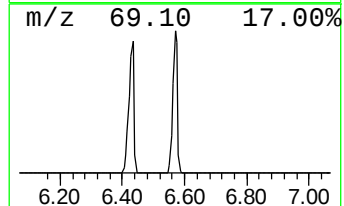
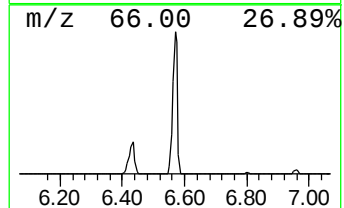
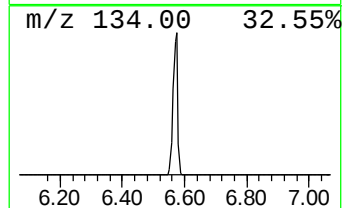
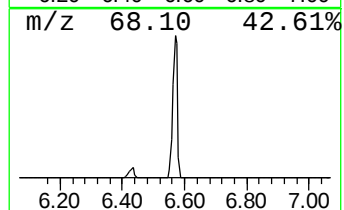
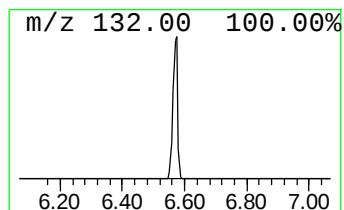
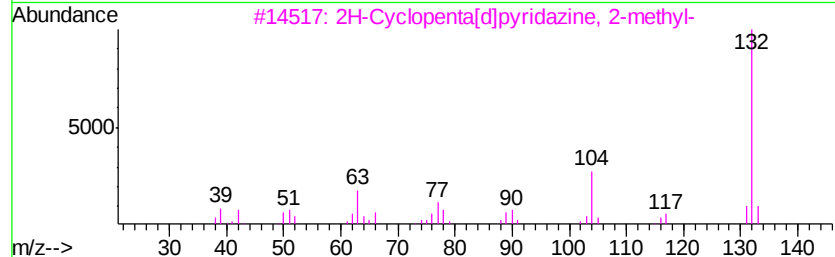
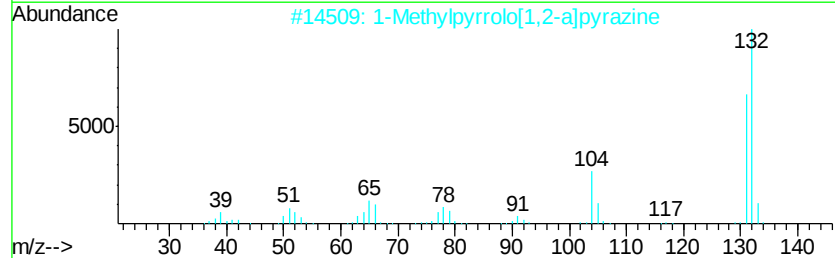
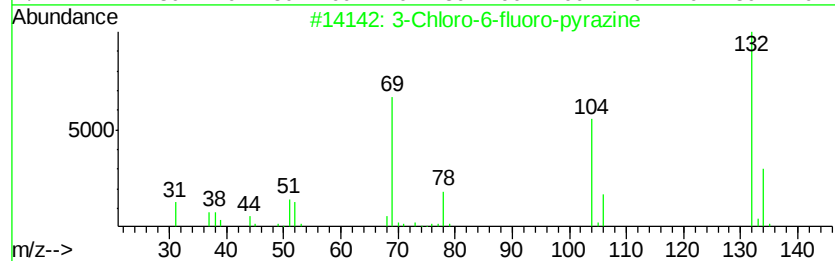
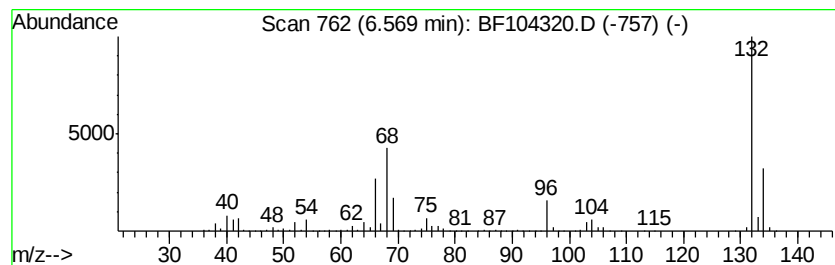
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown6.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	75.15 ng	3480100	1,4-Dichlorobenzene-d4	6.80

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
3		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	9



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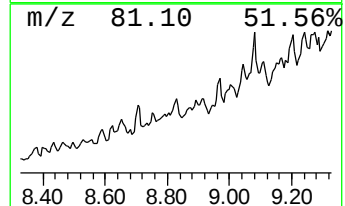
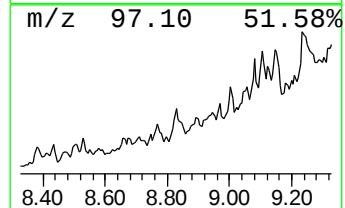
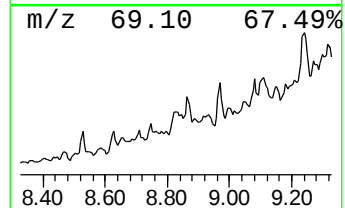
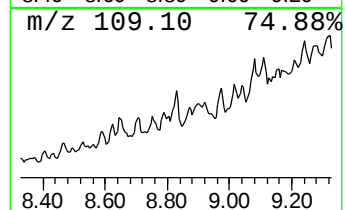
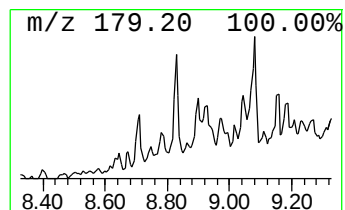
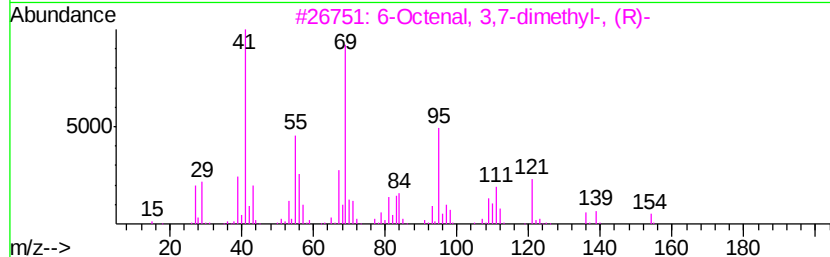
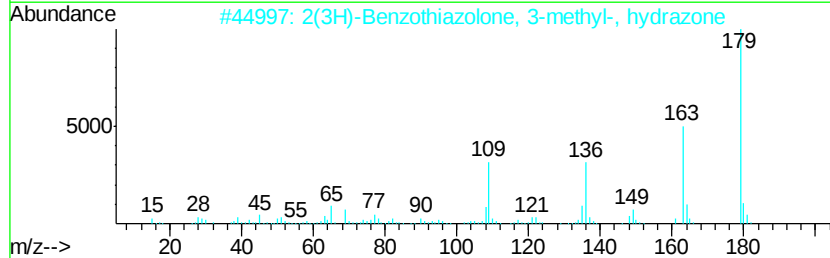
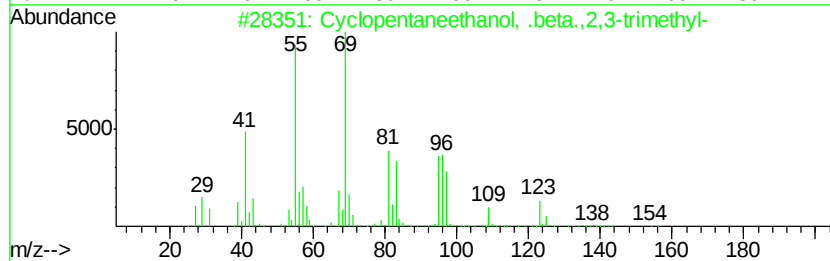
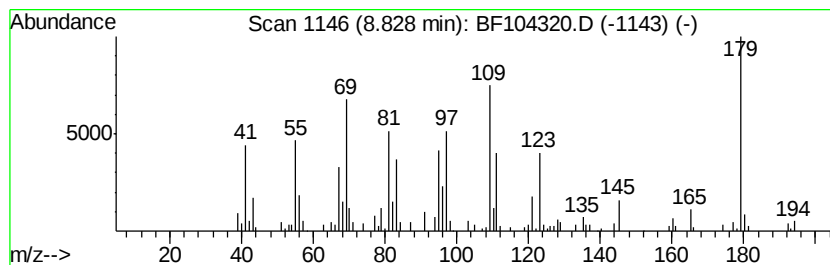
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown8.83 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.83	2.45 ng	158029	Naphthalene-d8	8.09

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentaneethanol, .beta.,2,3-...	156	C10H20O	021721-18-6	35
2		2(3H)-Benzothiazolone, 3-methyl-...	179	C8H9N3S	001128-67-2	22
3		6-Octenal, 3,7-dimethyl-, (R)-	154	C10H18O	002385-77-5	18
4		(2-Methylbutyl)cyclohexane	154	C11H22	054105-77-0	14
5		3-Oxatricyclo[4.1.1.0(2,4)]octan...	152	C10H16O	001686-14-2	14



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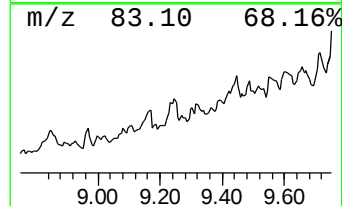
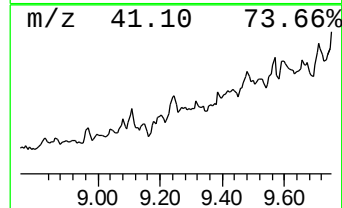
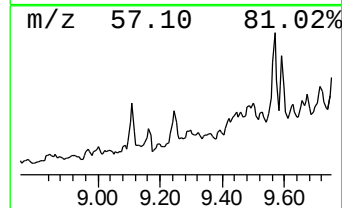
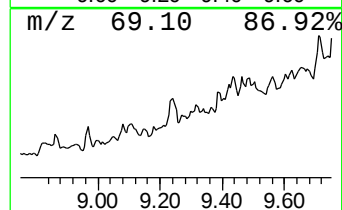
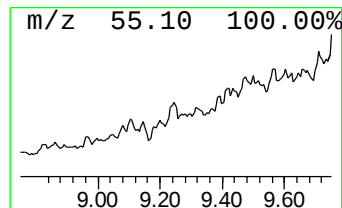
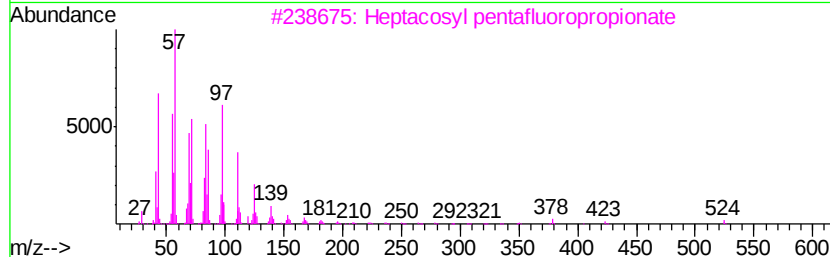
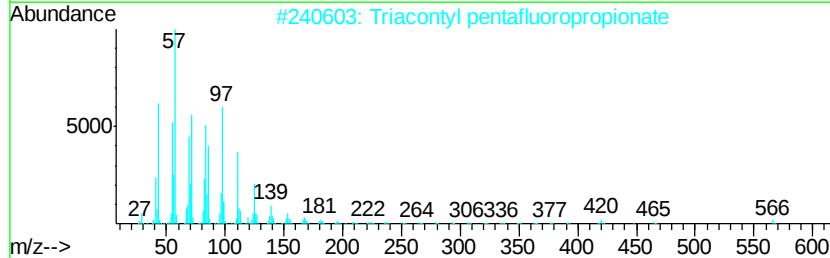
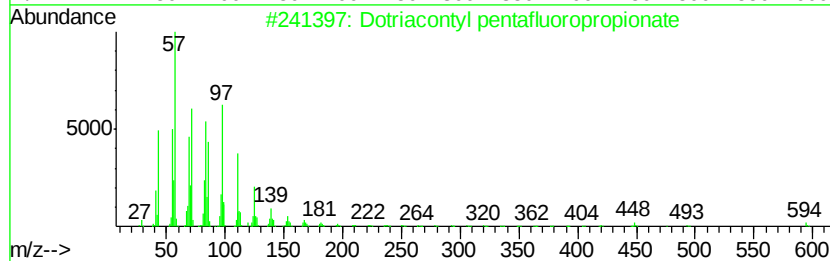
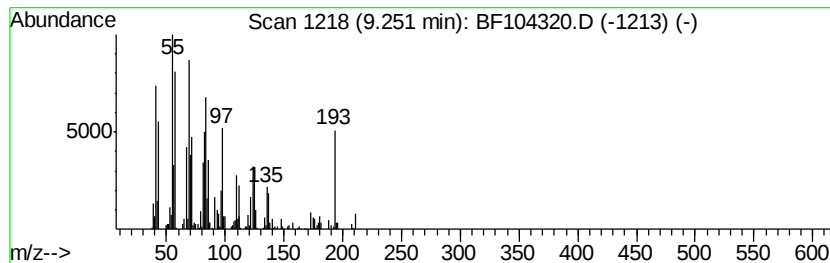
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 unknown9.25 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.25	3.13 ng	331331	Acenaphthene-d10	9.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dotriacontyl pentafluoropropionate	612	C35H65F5O2	1000351-81-4	30
2		Triacetyl pentafluoropropionate	584	C33H61F5O2	1000351-80-0	25
3		Heptacosyl pentafluoropropionate	542	C30H55F5O2	1000351-81-6	25
4		Triacetyl trifluoroacetate	534	C32H61F3O2	1000351-75-7	25
5		1,1'-Bicyclohexyl, 2-ethyl-, cis-	194	C14H26	050991-12-3	25



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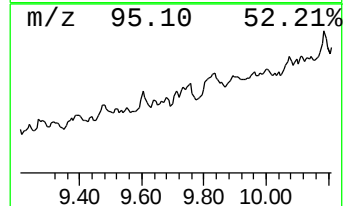
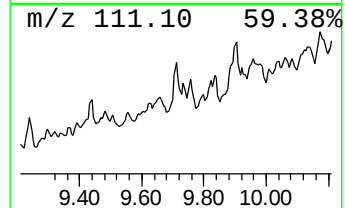
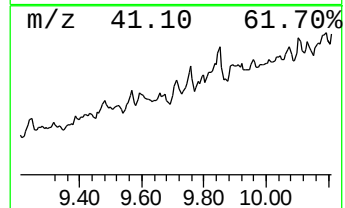
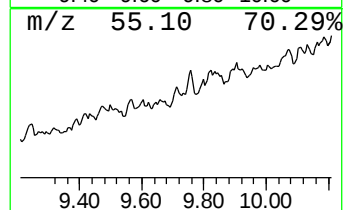
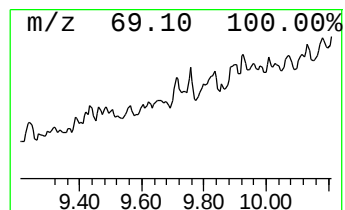
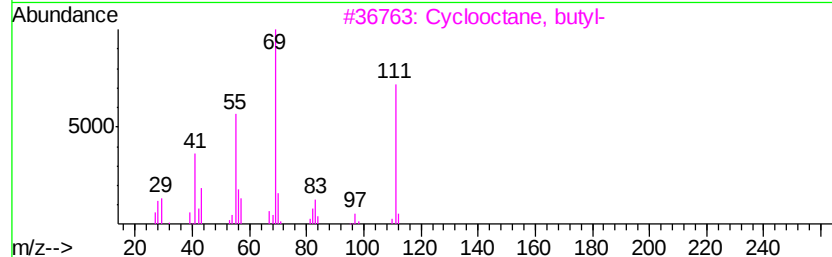
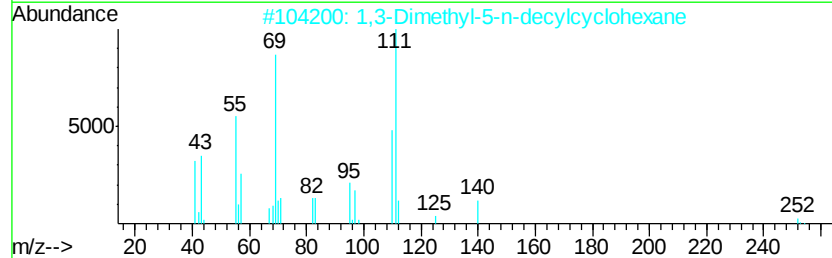
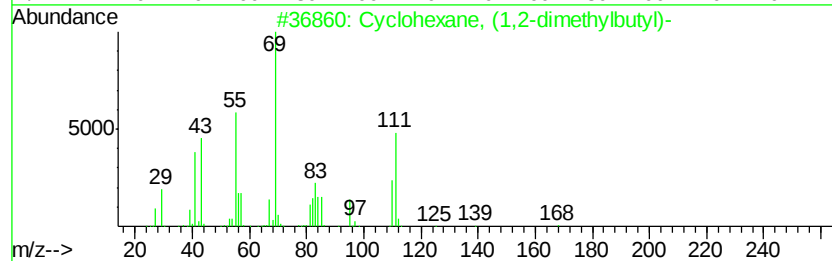
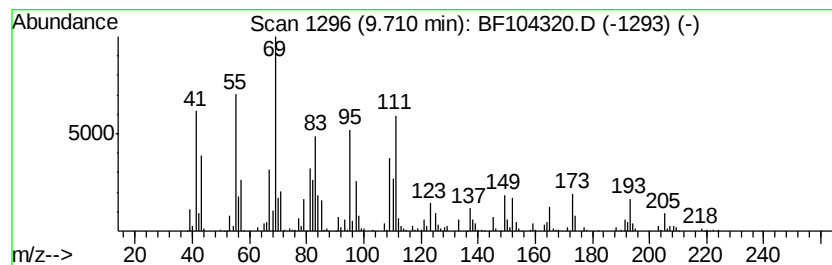
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Peak Number 6 Cyclohexane, (1,2-dimethylb... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.71	5.10 ng	540191	Acenaphthene-d10	9.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (1,2-dimethylbutyl)-	168	C12H24	061142-37-8	55
2			1,3-Dimethyl-5-n-decylcyclohexane	252	C18H36	086710-36-3	43
3			Cyclooctane, butyl-	168	C12H24	016538-93-5	35
4			1-Pentyne, 3-methyl-3-(1-methyle...	140	C9H16O	053966-55-5	35
5			2-Hexanone, 3-cyclohexylidene-4-...	208	C14H24O	1000150-19-2	27



Data Path : Z:\HPCHEM1\BNA F\Data\BF040618\
Data File : BF104320.D
Acq On : 6 Apr 2018 21:56
Operator : JU/SJ
Sample : J2232-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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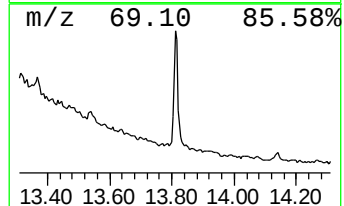
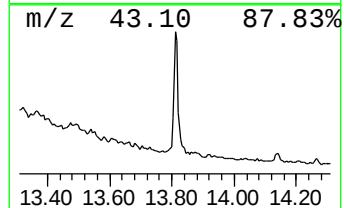
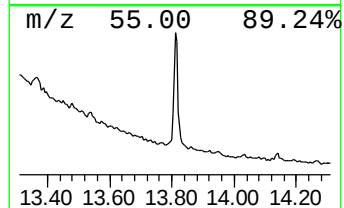
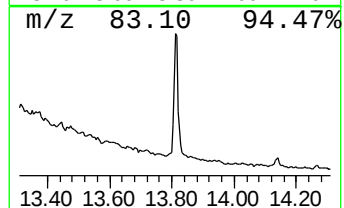
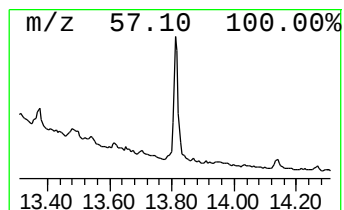
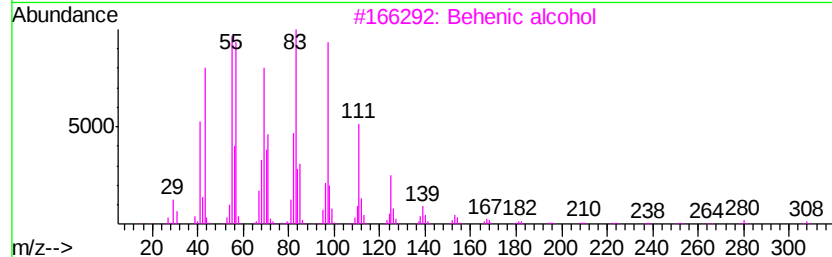
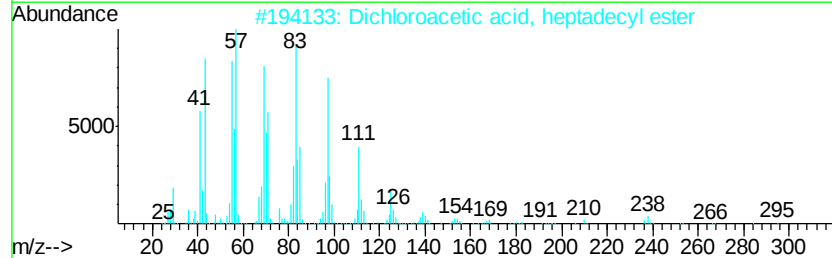
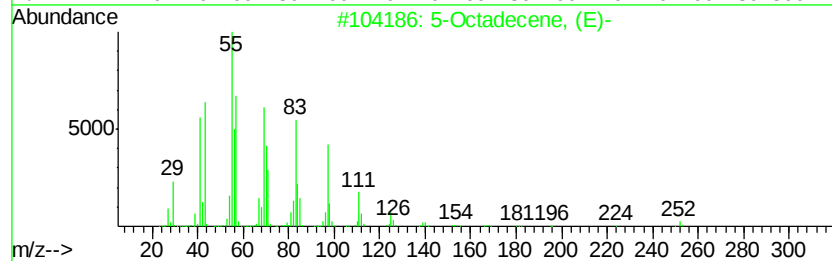
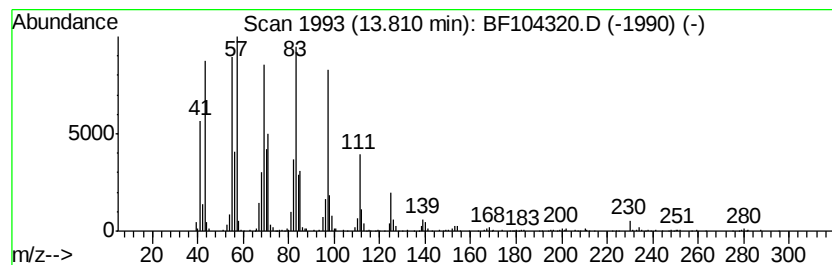
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 5-Octadecene, (E)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.81	15.71 ng	780571	Chrysene-d12	13.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Octadecene, (E)-	252	C18H36	007206-21-5	94
2		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
3		Behenic alcohol	326	C22H46O	000661-19-8	94
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5		1-Heneicosanol	312	C21H44O	015594-90-8	94



Data Path : Z:\HPCHEM1\BNA F\Data\BF040618\
 Data File : BF104320.D
 Acq On : 6 Apr 2018 21:56
 Operator : JU/SJ
 Sample : J2232-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CV-RB26667RT

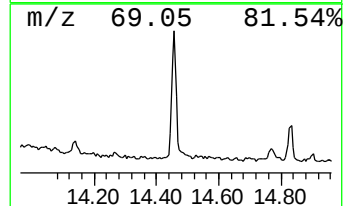
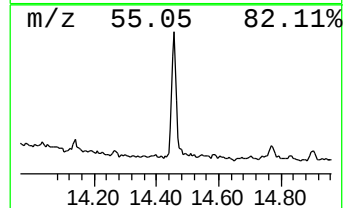
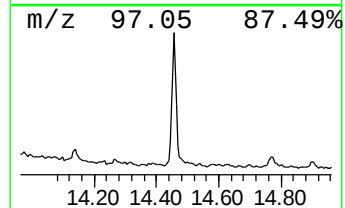
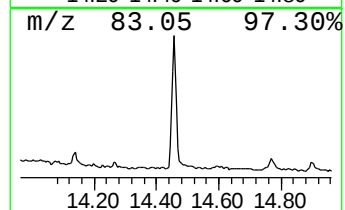
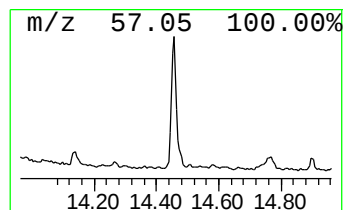
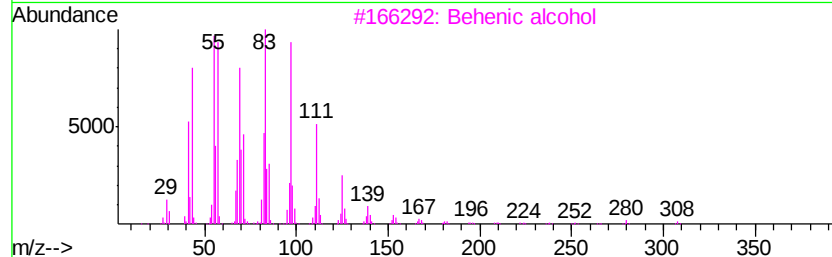
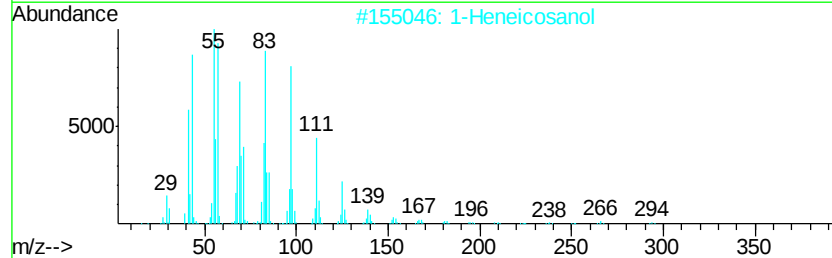
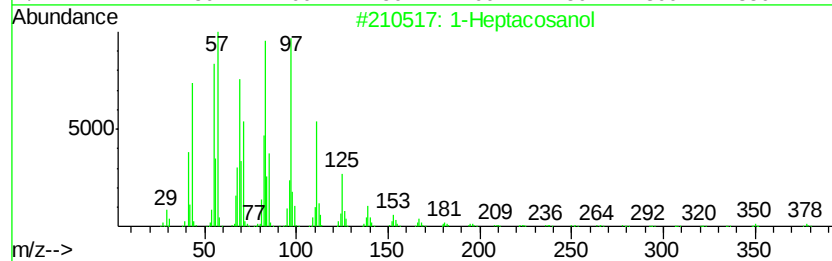
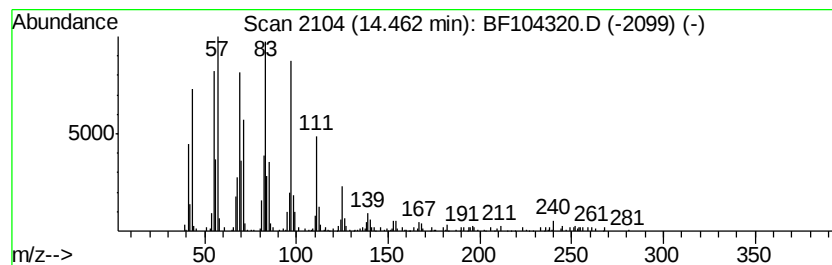
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 8 1-Heptacosanol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.46	10.65 ng	529075	Chrysene-d12	13.97

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heptacosanol	396	C27H56O	002004-39-9	95
2		1-Heneicosanol	312	C21H44O	015594-90-8	93
3		Behenic alcohol	326	C22H46O	000661-19-8	93
4		n-Tetracosanol-1	354	C24H50O	000506-51-4	93
5		Octacosanol	410	C28H58O	000557-61-9	92



Data Path : Z:\HPCHEM1\BNA F\Data\BF040618\
Data File : BF104320.D
Acq On : 6 Apr 2018 21:56
Operator : JU/SJ
Sample : J2232-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

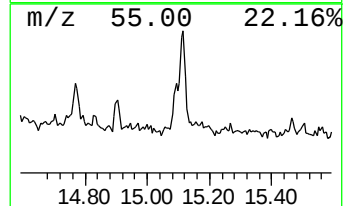
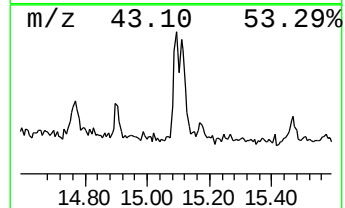
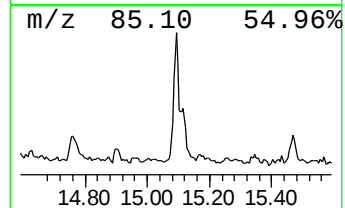
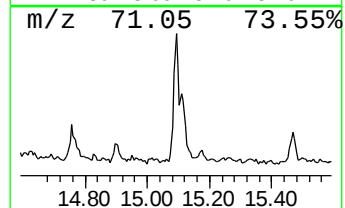
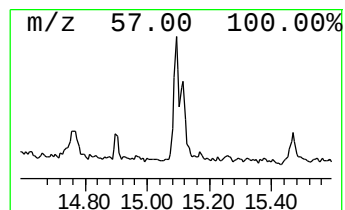
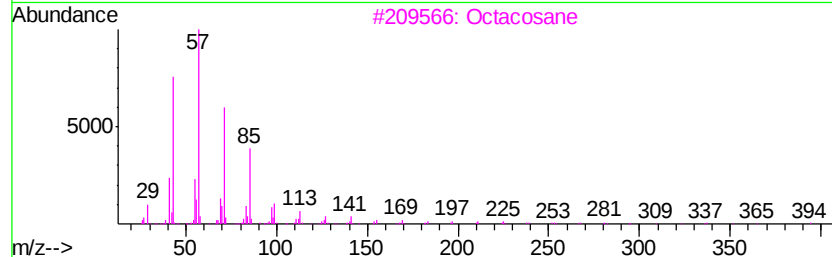
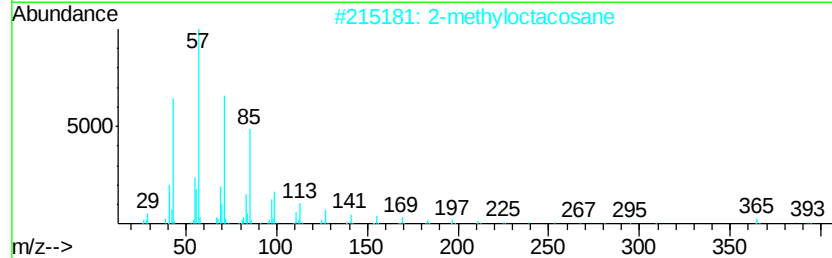
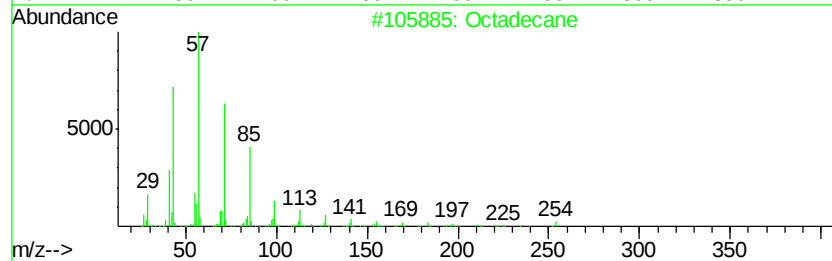
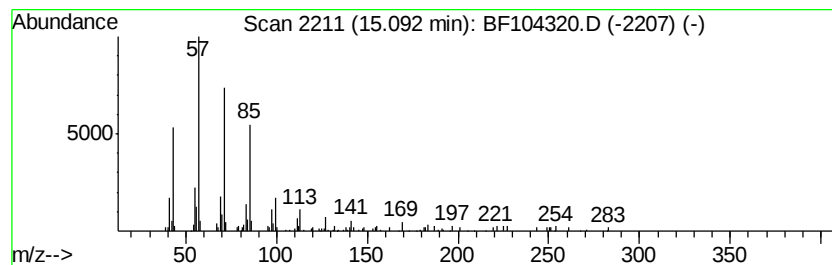
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Octadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.09	2.71 ng	118896	Perylene-d12	15.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octadecane	254 C18H38	000593-45-3	93
2	2-methyloctacosane	408 C29H60	1000376-72-8	91
3	Octacosane	394 C28H58	000630-02-4	91
4	Heptadecane, 9-octyl-	352 C25H52	007225-64-1	91
5	triacontane	422 C30H62	000638-68-6	91



Data Path : Z:\HPCHEM1\BNA F\Data\BF040618\
Data File : BF104320.D
Acq On : 6 Apr 2018 21:56
Operator : JU/SJ
Sample : J2232-01
Misc :
ALS Vial : 16 Sample Multiplier: 1

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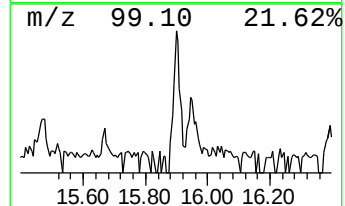
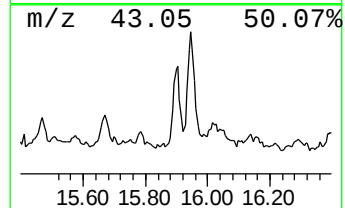
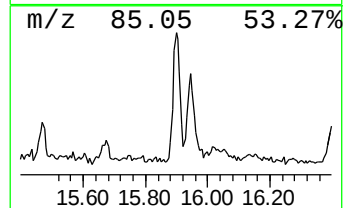
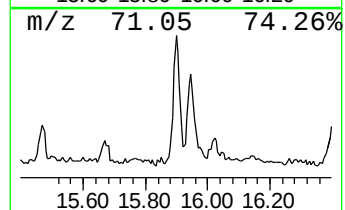
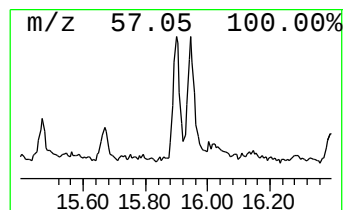
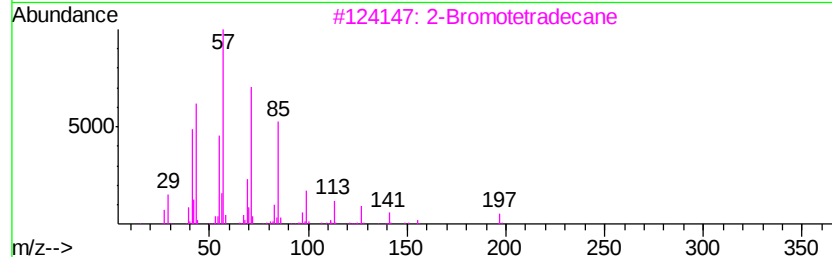
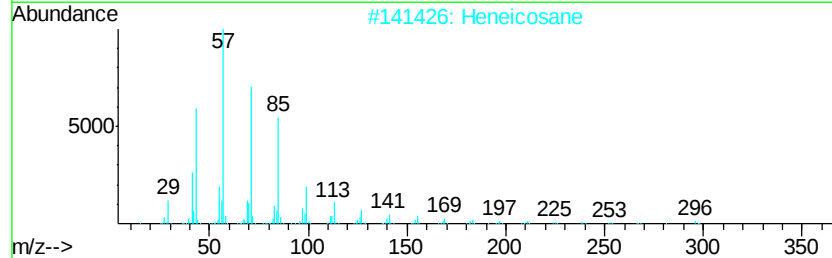
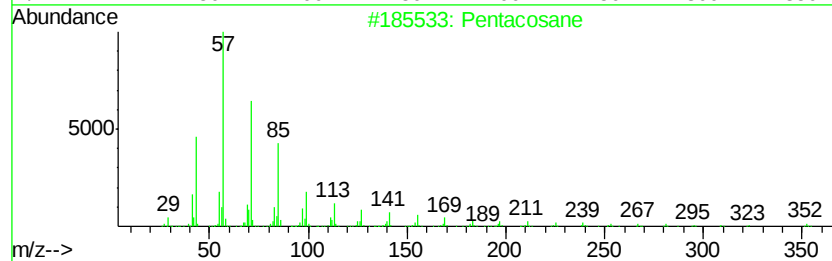
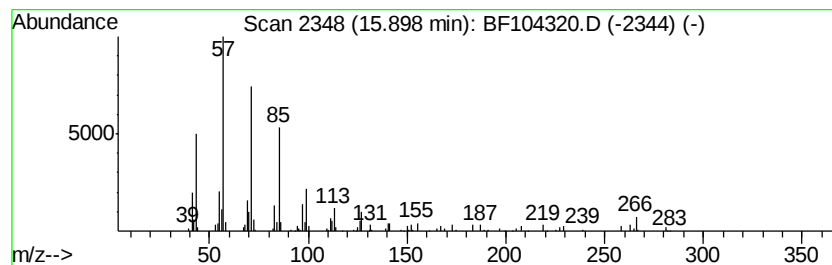
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Pentacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.90	2.53 ng	110845	Perylene-d12	15.43

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane		352	C25H52	000629-99-2	90
2	Heneicosane		296	C21H44	000629-94-7	87
3	2-Bromotetradecane		276	C14H29Br	074036-95-6	86
4	Nonadecane, 2-methyl-		282	C20H42	001560-86-7	86
5	5-Ethyl-5-methylnonadecane		310	C22H46	1000360-42-7	80



Data Path : Z:\HPCHEM1\BNA F\Data\BF040618\
Data File : BF104320.D
Acq On : 6 Apr 2018 21:56
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Sample : J2232-01
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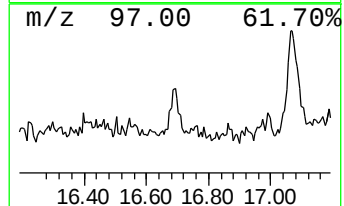
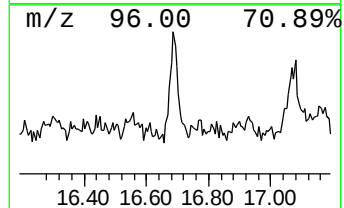
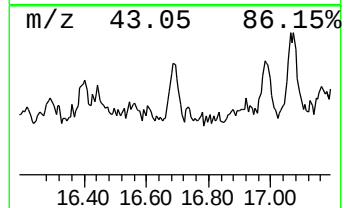
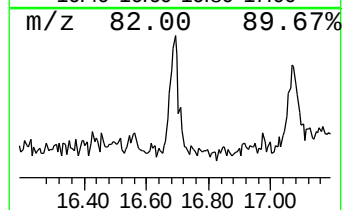
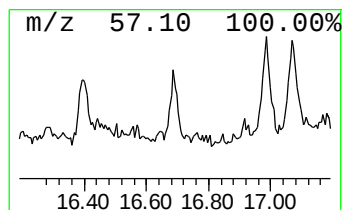
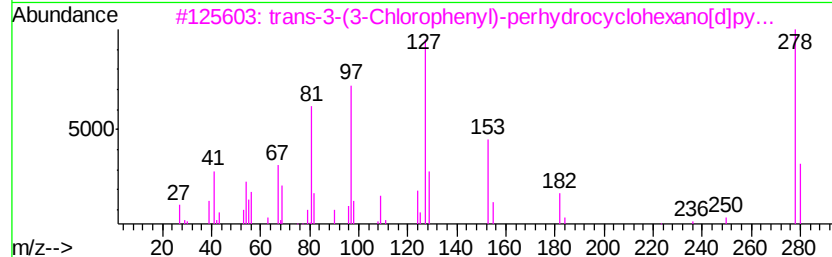
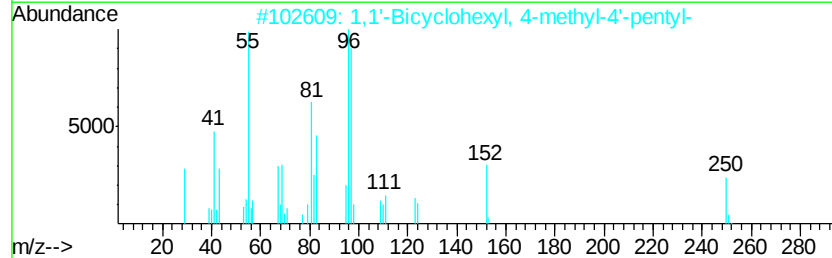
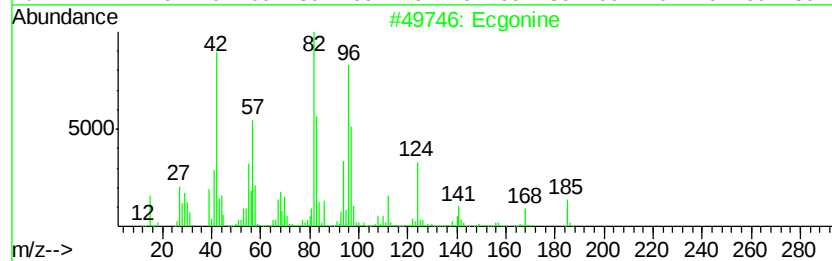
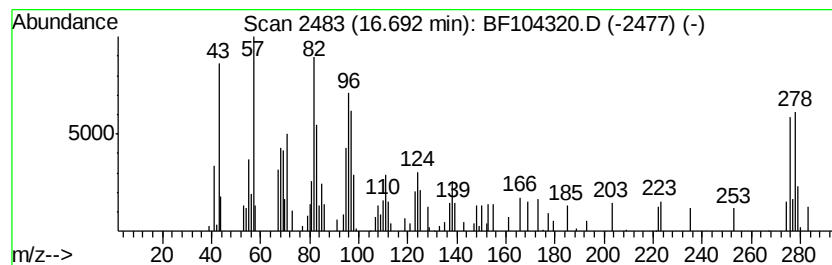
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown16.69 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.69	2.21 ng	96961	Perylene-d12	15.43
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Ecgonine	185 C9H15NO3	000481-37-8 38
2		1,1'-Bicyclohexyl, 4-methyl-4'-p...	250 C18H34	092263-42-8 38
3		trans-3-(3-Chlorophenyl)-perhydr...	278 C14H15ClN2O2	1000148-52-4 30
4		1,1'-Bicyclohexyl, 4,4'-dimethyl-	194 C14H26	054823-99-3 27
5		11H-Dibenzo[c,f][1,2]diazepine, ...	278 C13H8Cl2N2O	000958-71-4 22



Data Path : Z:\HPCHEM1\BNA F\Data\BF040618\
Data File : BF104320.D
Acq On : 6 Apr 2018 21:56
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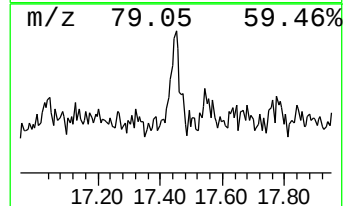
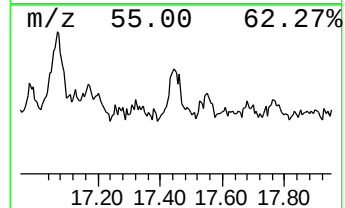
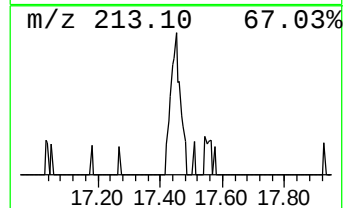
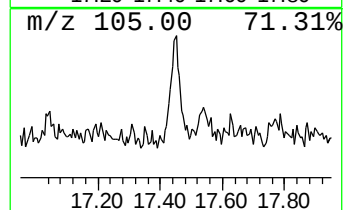
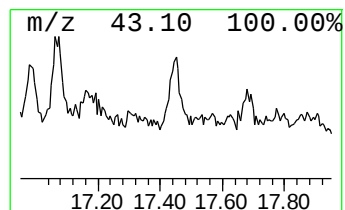
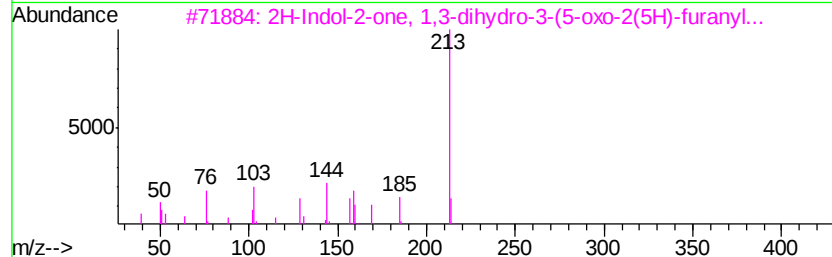
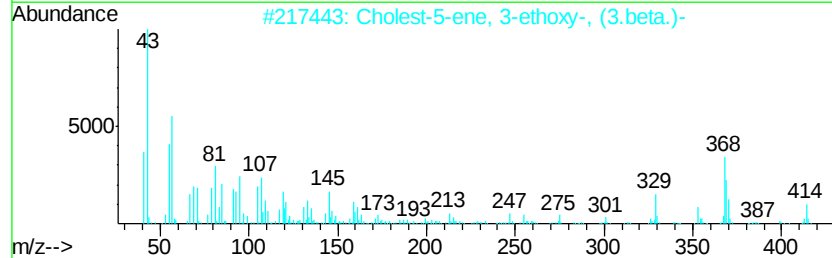
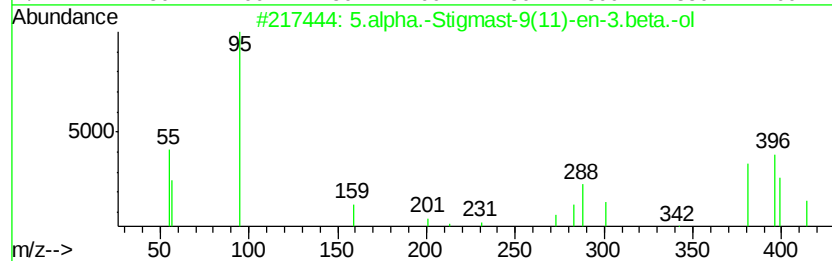
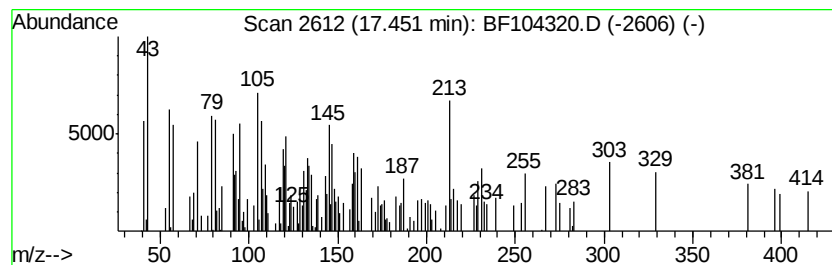
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 unknown17.45 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.45	3.42 ng	149886	Perylene-d12	15.43

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5.alpha.-Stigmast-9(11)-en-3.beta...	414	C29H50O	077794-81-1	14
2		Cholest-5-ene, 3-ethoxy-, (3.beta...	414	C29H50O	000986-19-6	11
3		2H-Indol-2-one, 1,3-dihydro-3-(5...	213	C12H7NO3	013191-62-3	10
4		2,6-Diamino-9-[2,6-dichlorobenzyl...	308	C12H10Cl2N6	1000212-70-6	9
5		1-Pentyne, 1-(trimethylsilyl)-4-...	270	C14H30OSi2	167390-04-7	9



Data Path : Z:\HPCHEM1\BNA_F\Data\BF040618\
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Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
CV-RB26667RT

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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unknown6.57	6.57	75.2 ng		3480100	1	6.80	926191	20.0
unknown8.83	8.83	2.5 ng		158029	2	8.09	1287850	20.0
unknown9.25	9.25	3.1 ng		331331	3	9.85	2116940	20.0
Cyclohexane, (1,2...	9.71	5.1 ng		540191	3	9.85	2116940	20.0
5-Octadecene, (E)-	13.81	15.7 ng		780571	5	13.97	993706	20.0
1-Heptacosanol	14.46	10.7 ng		529075	5	13.97	993706	20.0
Octadecane	15.09	2.7 ng		118896	6	15.43	877781	20.0
Pentacosane	15.90	2.5 ng		110845	6	15.43	877781	20.0
unknown16.69	16.69	2.2 ng		96961	6	15.43	877781	20.0
unknown17.45	17.45	3.4 ng		149886	6	15.43	877781	20.0