

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122015\
 Data File : BF083958.D
 Acq On : 19 Dec 2015 16:57
 Operator : UM/SJ
 Sample : G4875-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SWD-1

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-BF121815.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	5.013	254	256	261	rVB	122805	170259	6.26%	1.090%
2	5.459	292	295	298	rBV	542940	496286	18.25%	3.177%
3	6.487	383	385	394	rVB	403451	347792	12.79%	2.226%
4	6.613	394	396	399	rBV	787200	1007271	37.03%	6.448%
5	6.842	414	416	419	rBV	346617	446561	16.42%	2.859%
6	7.002	428	430	433	rBV	1333803	1480185	54.42%	9.475%
7	7.413	463	466	468	rVB	1091326	1050644	38.63%	6.726%
8	7.676	486	489	494	rVB	22543	35975	1.32%	0.230%
9	7.996	512	517	520	rBV3	53388	146356	5.38%	0.937%
10	8.133	527	529	531	rBV	609417	475184	17.47%	3.042%
11	8.270	537	541	543	rBV3	33118	62068	2.28%	0.397%
12	8.305	543	544	547	rVV	44503	63686	2.34%	0.408%
13	8.385	547	551	552	rVV2	24006	53752	1.98%	0.344%
14	8.442	552	556	559	rVV4	53045	127716	4.70%	0.818%
15	8.545	559	565	568	rVV2	135482	348758	12.82%	2.233%
16	8.590	568	569	573	rVV2	104373	157033	5.77%	1.005%
17	8.682	573	577	579	rVB3	20816	41009	1.51%	0.263%
18	9.208	620	623	626	rBV	2022996	2110340	77.59%	13.509%
19	9.333	632	634	637	rBV	33552	32112	1.18%	0.206%
20	9.596	655	657	660	rBV	109899	139849	5.14%	0.895%
21	9.756	667	671	673	rBV2	29777	49624	1.82%	0.318%
22	9.813	673	676	678	rVB3	15009	28453	1.05%	0.182%
23	9.882	680	682	685	rVV	638319	629383	23.14%	4.029%
24	10.270	713	716	717	rBV	94902	101556	3.73%	0.650%
25	10.305	717	719	722	rVB3	21805	34821	1.28%	0.223%
26	10.671	748	751	754	rBV	978328	944884	34.74%	6.049%
27	10.796	758	762	766	rBV2	23721	54525	2.00%	0.349%
28	11.242	798	801	802	rBV	27654	29404	1.08%	0.188%
29	11.368	809	812	814	rBV	752428	655119	24.09%	4.194%
30	11.905	857	859	863	rBV2	52038	85076	3.13%	0.545%
31	12.682	925	927	931	rBV2	39431	46435	1.71%	0.297%
32	12.956	948	951	953	rBV	2456751	2719882	100.00%	17.411%
33	13.848	1027	1029	1032	rBV	67481	82616	3.04%	0.529%
34	13.997	1040	1042	1045	rBV	540370	672098	24.71%	4.302%

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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-BF121815.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.625	1095	1097	1099	rBV	113511	114792	4.22%	0.735%
36	14.842	1114	1116	1122	rVB	23833	39576	1.46%	0.253%
37	15.437	1165	1168	1178	rVB	377266	540236	19.86%	3.458%

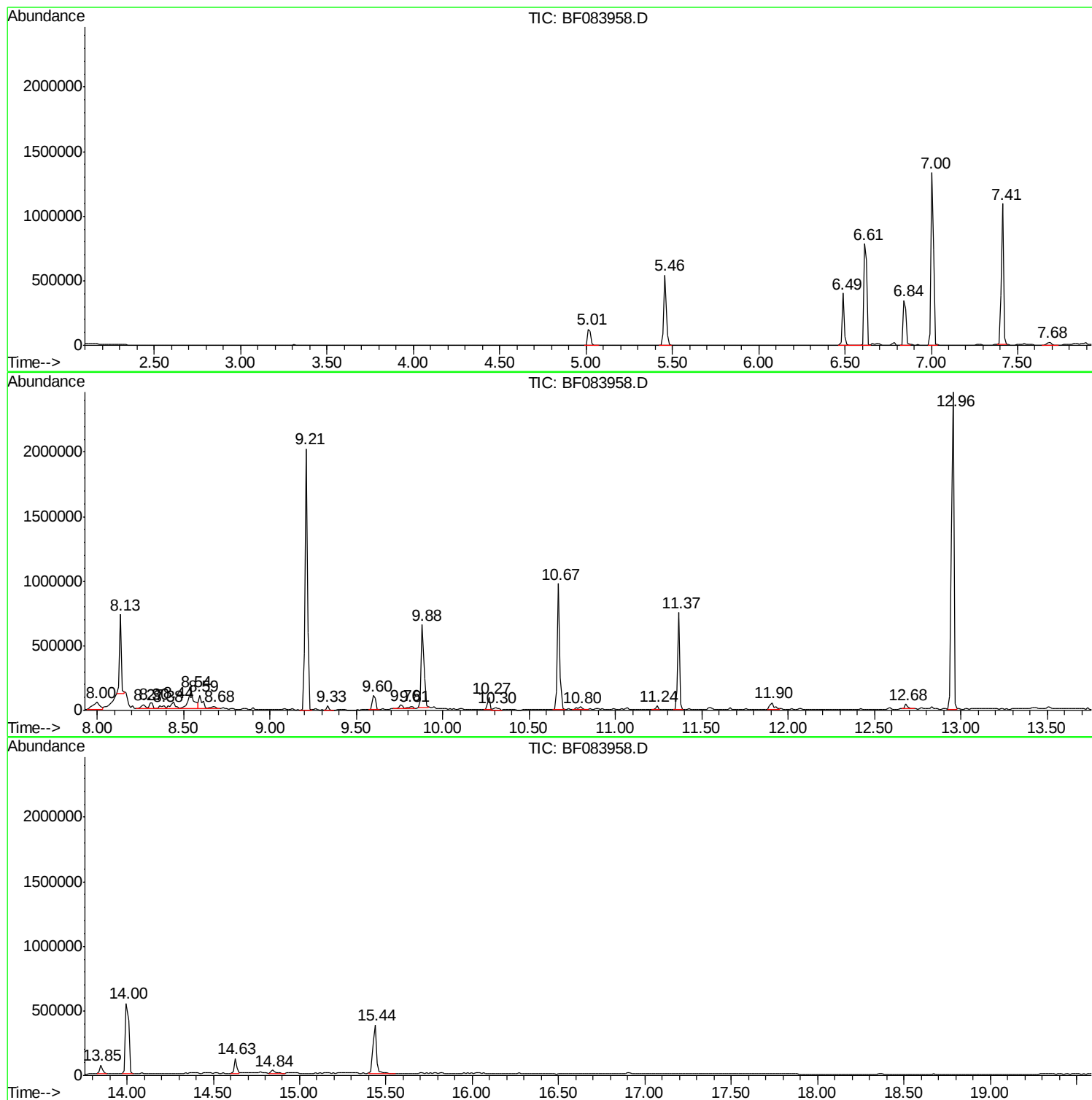
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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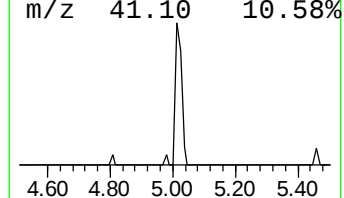
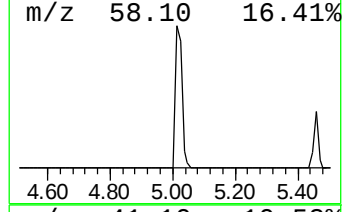
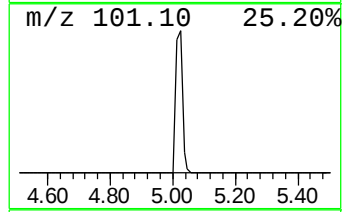
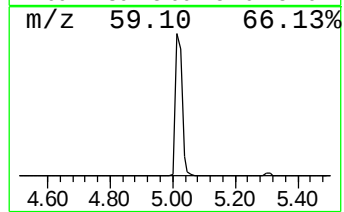
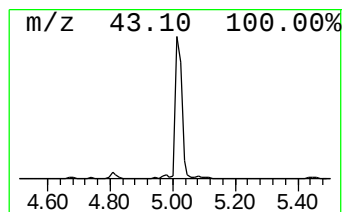
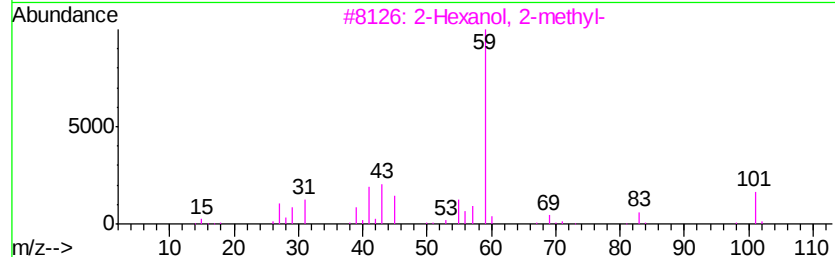
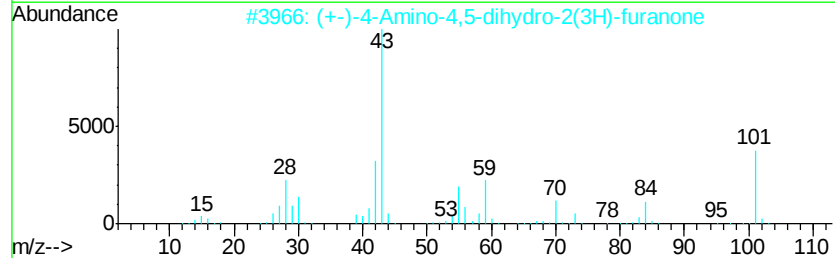
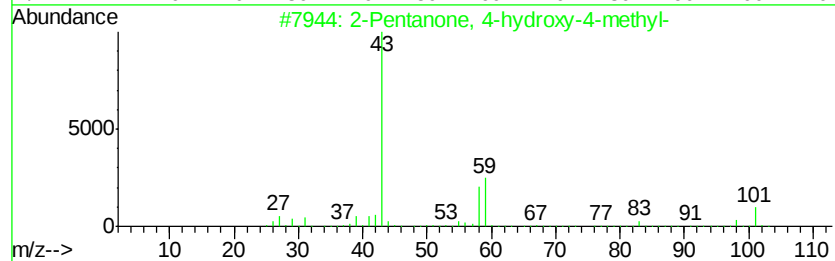
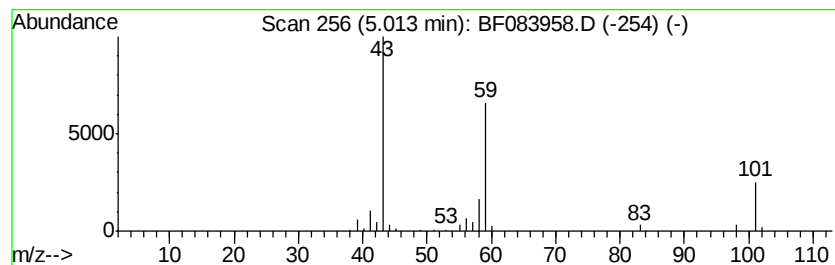
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	7.63 ng	170259	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	47
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	40
4		3-Hydroxy-3-methyl-2-butanone	102	C5H10O2	000115-22-0	25
5		1,2-Butanediol, (.+/-.)-	90	C4H10O2	026171-83-5	25



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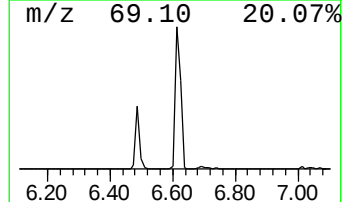
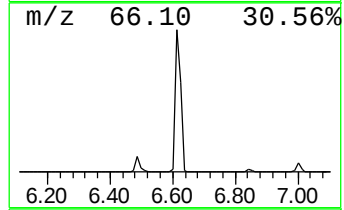
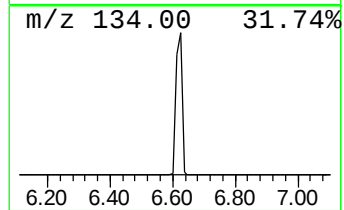
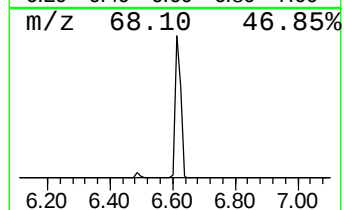
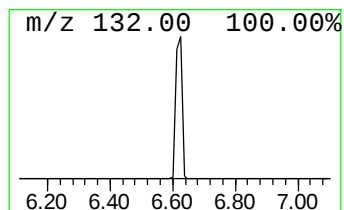
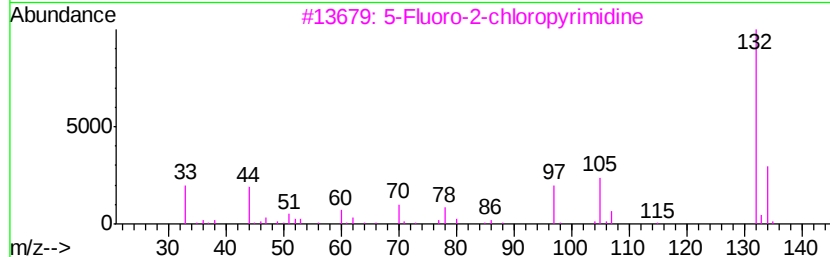
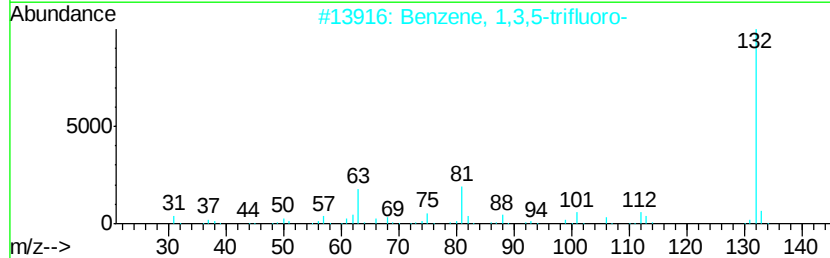
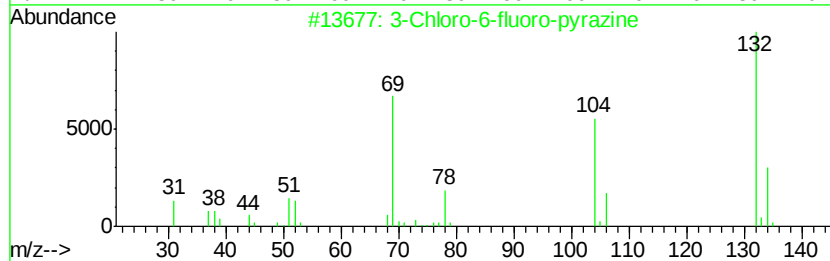
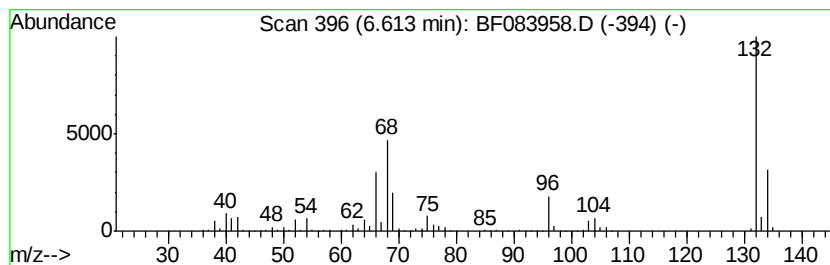
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown6.61 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.61	45.11 ng	1007270	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	14
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10



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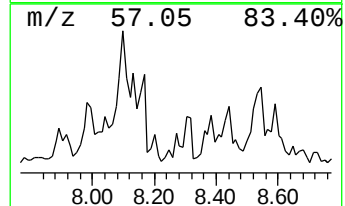
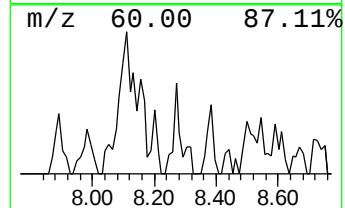
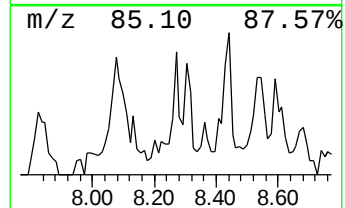
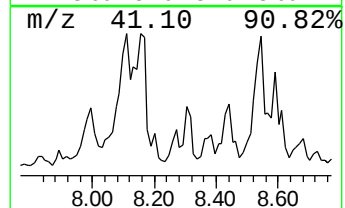
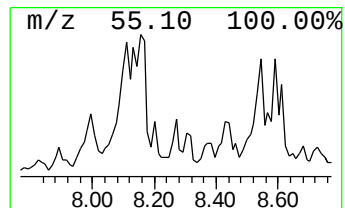
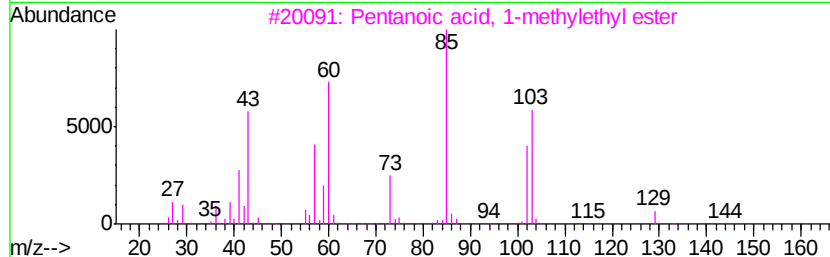
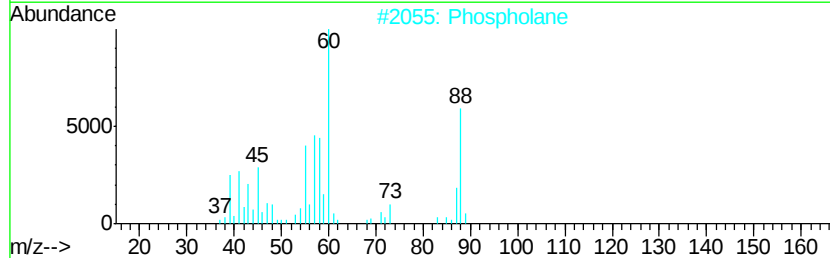
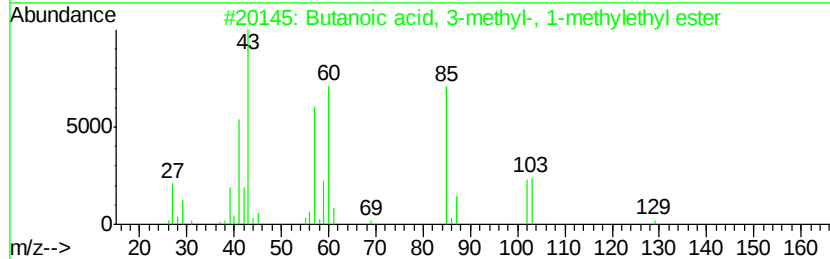
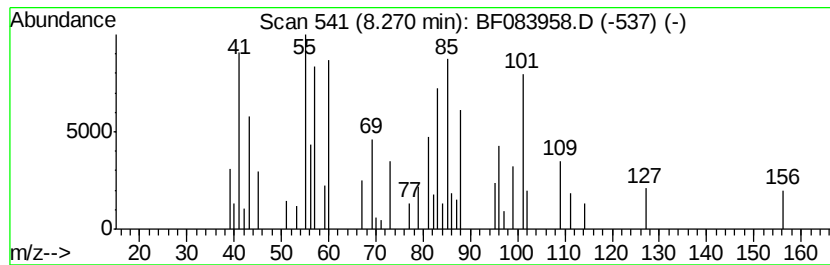
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown8.27 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.27	2.61 ng	62068	Naphthalene-d8	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butanoic acid, 3-methyl-, 1-meth...	144	C8H16O2	032665-23-9	27
2		Phospholane	88	C4H9P	003466-00-0	18
3		Pentanoic acid, 1-methylethyl ester	144	C8H16O2	018362-97-5	16
4		Heptanedioic acid	160	C7H12O4	000111-16-0	14
5		Thiophene, tetrahydro-	88	C4H8S	000110-01-0	14



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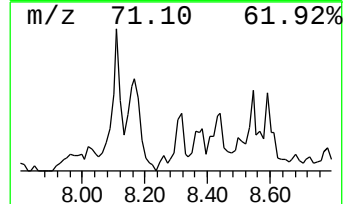
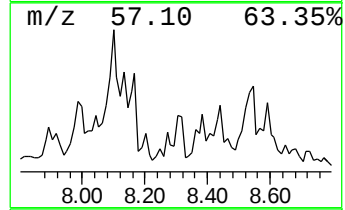
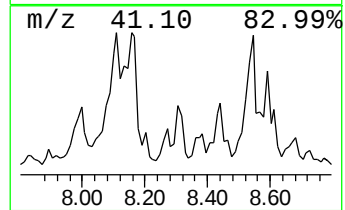
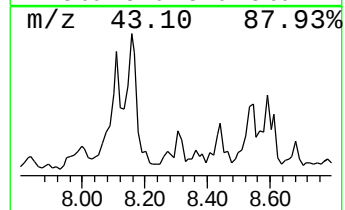
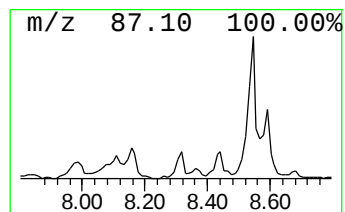
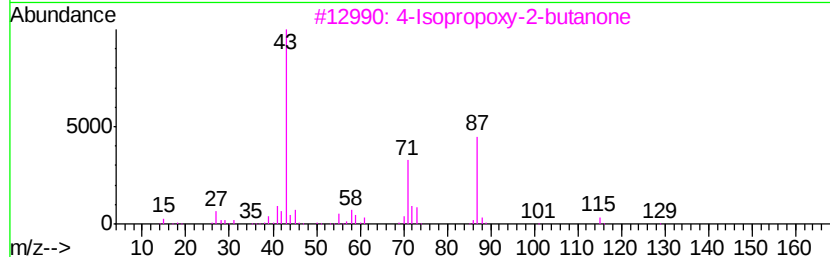
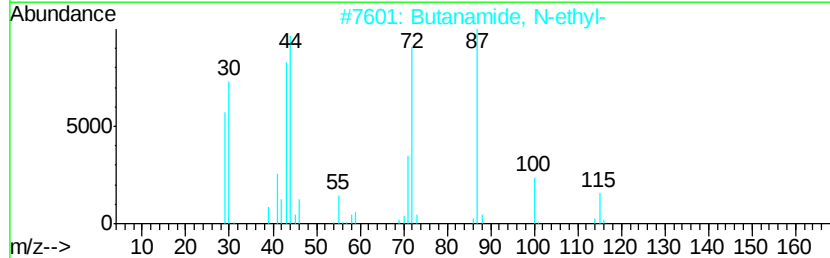
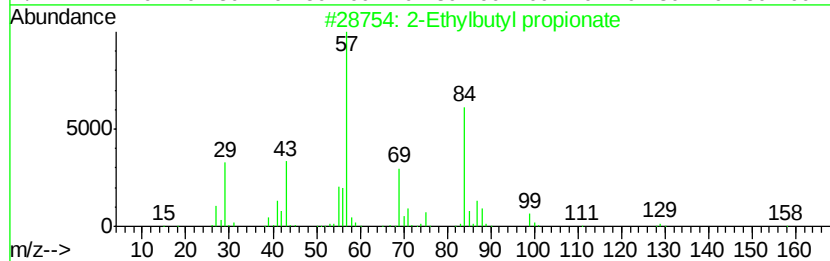
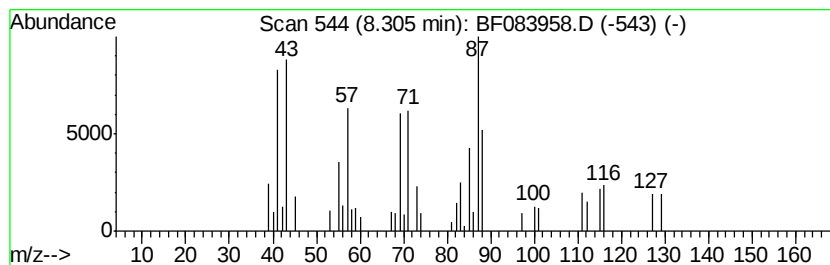
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown8.30 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.30	2.68 ng	63686	Naphthalene-d8	8.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Ethylbutyl	propionate	158	C9H18O2	090677-00-2	40	
2	Butanamide, N-ethyl-	115	C6H13NO	013091-16-2	27		
3	4-Isopropoxy-2-butanone	130	C7H14O2	032541-58-5	25		
4	2-t-Butylpentanoic acid, methyl ...	172	C10H20O2	1000202-22-5	22		
5	3-Mercaptohexyl	butanoate	204	C10H20O2S	136954-21-7	22	



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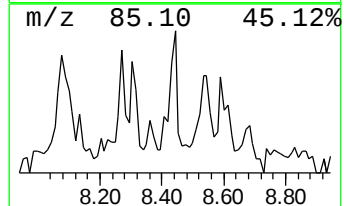
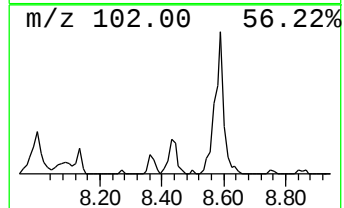
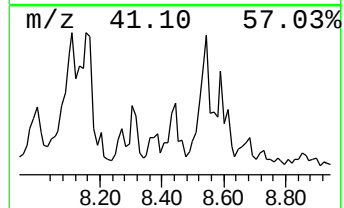
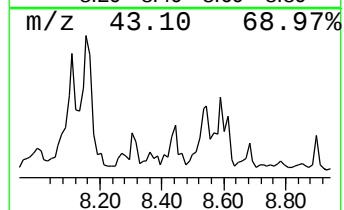
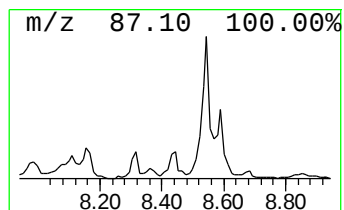
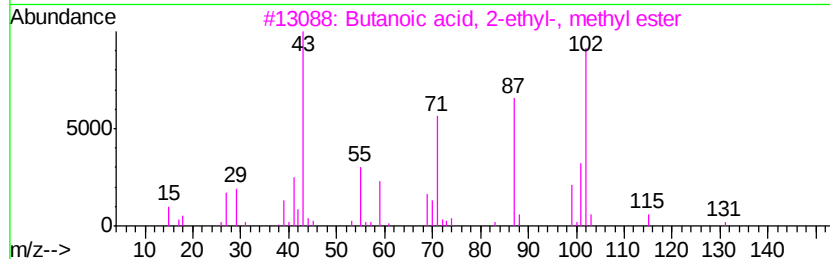
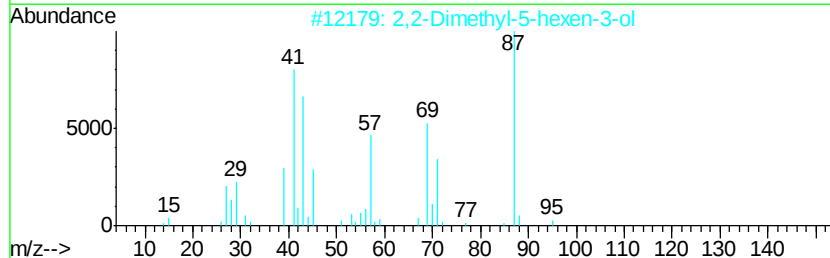
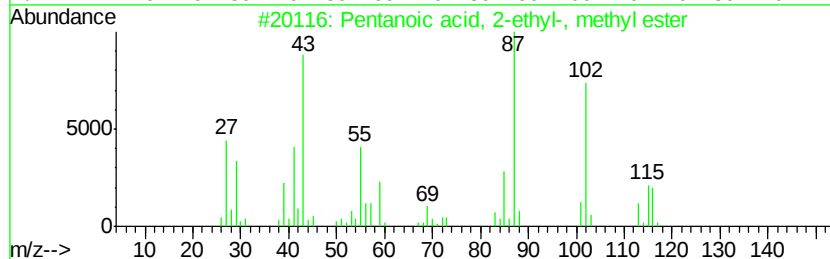
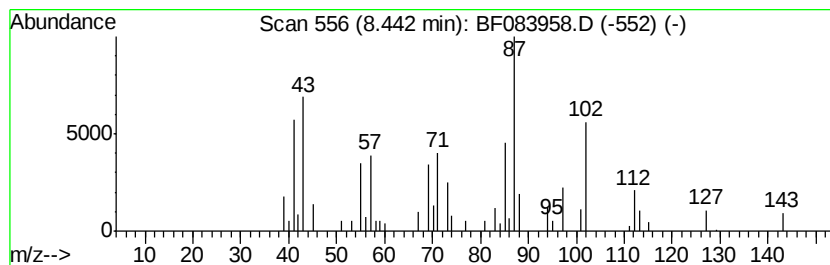
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown8.44 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.44	5.38 ng	127716	Naphthalene-d8	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentanoic acid, 2-ethyl-, methyl...	144	C8H16O2	000816-16-0	32
2		2,2-Dimethyl-5-hexen-3-ol	128	C8H16O	019550-89-1	12
3		Butanoic acid, 2-ethyl-, methyl ...	130	C7H14O2	000816-11-5	12
4		Thiophane, propyl-	130	C7H14S	001551-34-4	10
5		Butanamide, N,N,3,3-tetramethyl-	143	C8H17NO	026153-90-2	9



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Sample : G4875-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SWD-1

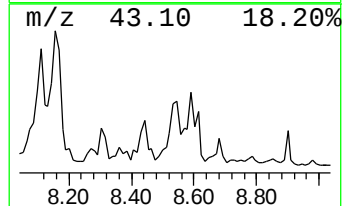
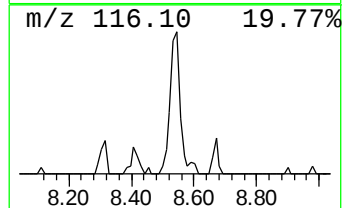
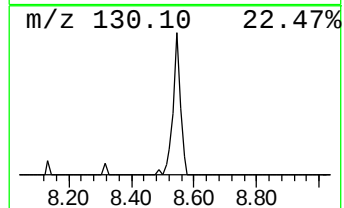
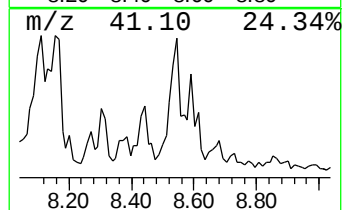
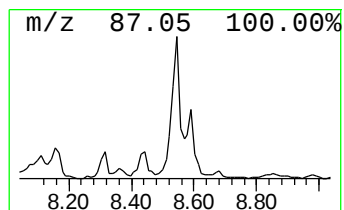
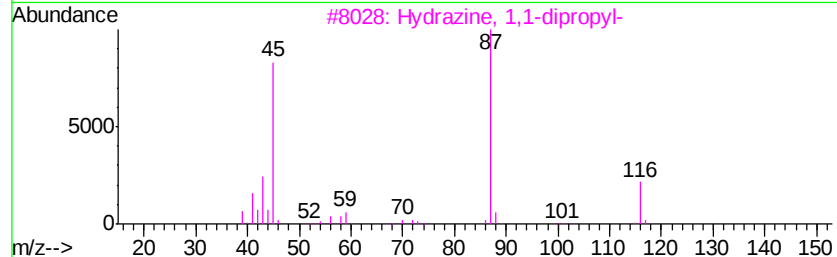
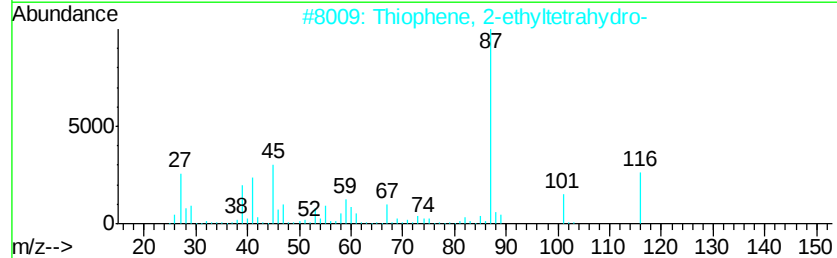
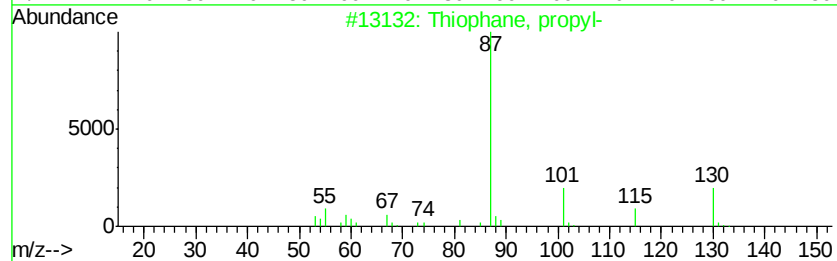
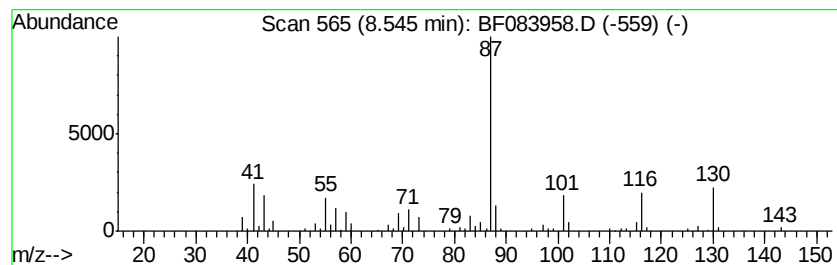
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Thiophane, propyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.54	14.68 ng	348758	Napthalene-d8	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Thiophane, propyl-	130	C7H14S	001551-34-4	64
2		Thiophene, 2-ethyltetrahydro-	116	C6H12S	001551-32-2	59
3		Hvdrazine, 1,1-dipropyl-	116	C6H16N2	004986-50-9	49
4		3,4-Di-O-methyl-L-arabinopyranose	178	C7H14O5	086049-20-9	42
5		.alpha.-D-Xylofuranoside, methyl...	192	C8H16O5	035007-52-4	40



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122015\
Data File : BF083958.D
Acq On : 19 Dec 2015 16:57
Operator : UM/SJ
Sample : G4875-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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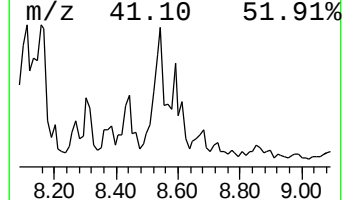
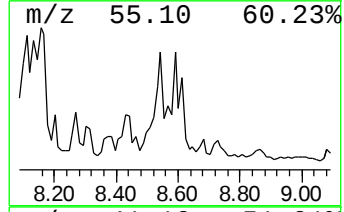
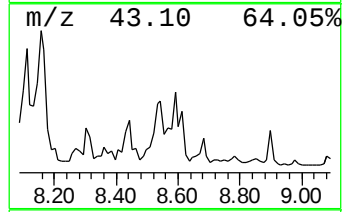
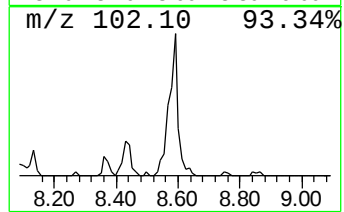
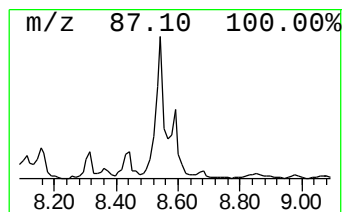
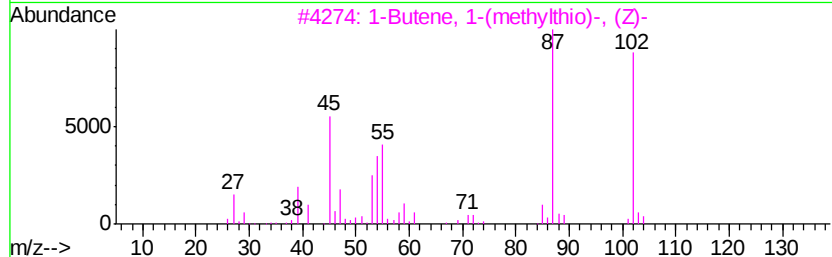
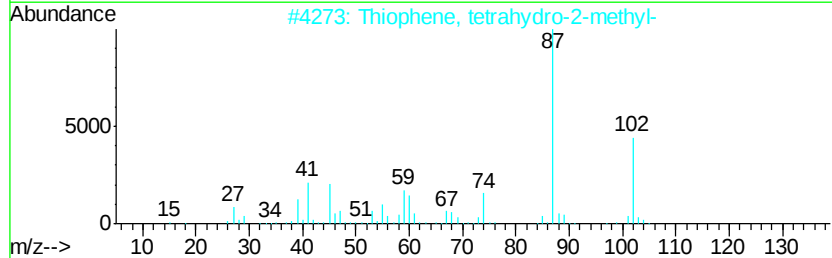
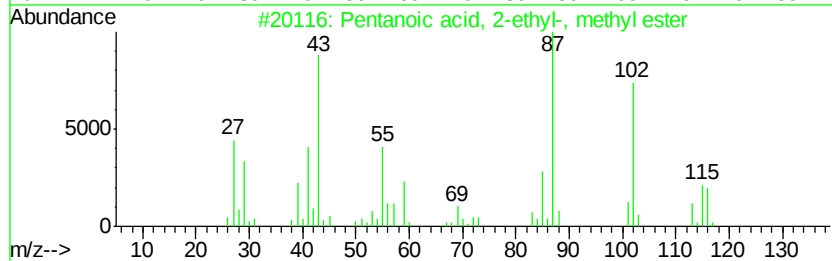
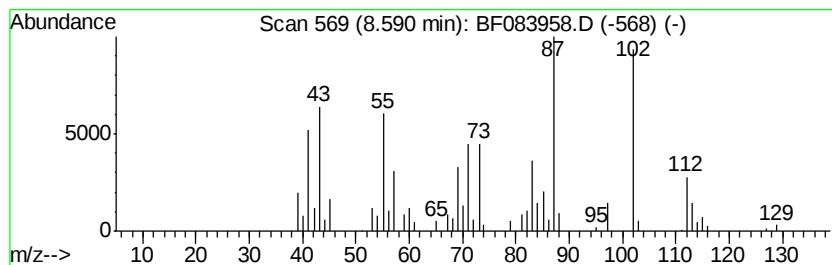
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Pentanoic acid, 2-ethyl-, m... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.59	6.61 ng	157033	Naphthalene-d8	8.13

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentanoic acid, 2-ethyl-, methyl...	144	C8H16O2	000816-16-0	50
2			Thiophene, tetrahydro-2-methyl-	102	C5H10S	001795-09-1	47
3			1-Butene, 1-(methylthio)-, (Z)-	102	C5H10S	017414-15-2	47
4			Butanoic acid, 2-ethyl-, methyl ...	130	C7H14O2	000816-11-5	42
5			1-Butene, 1-(methylthio)-, (E)-	102	C5H10S	017414-27-6	38



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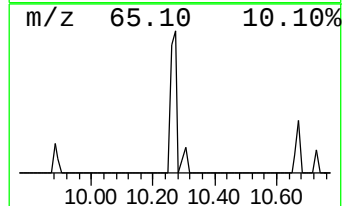
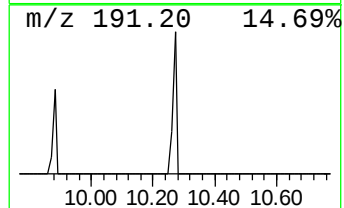
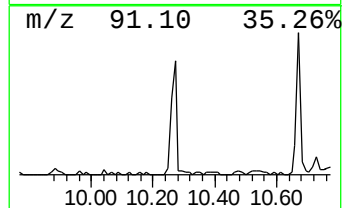
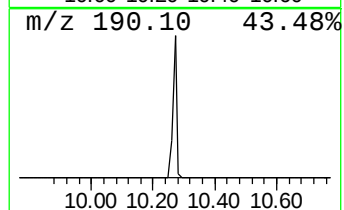
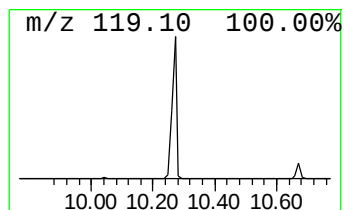
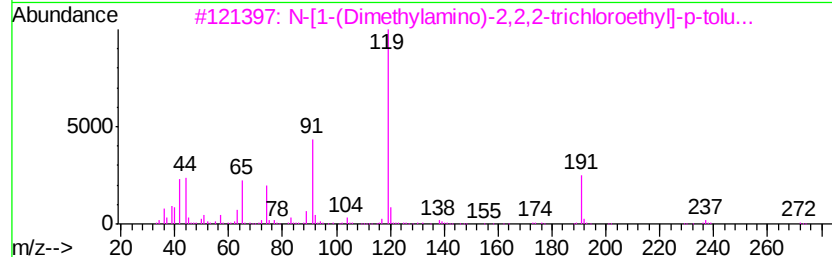
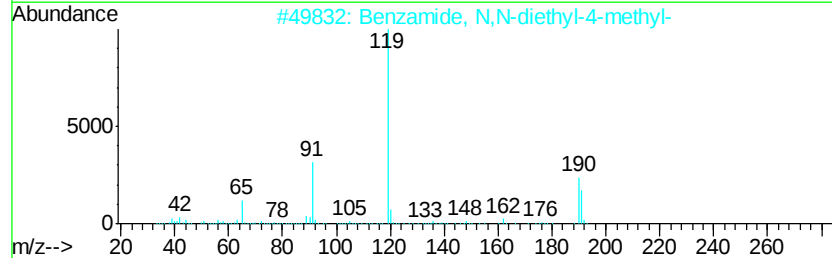
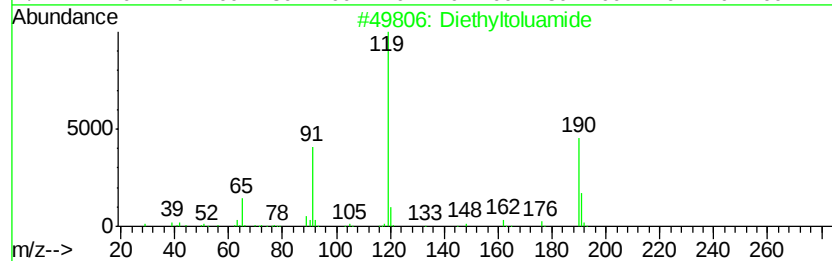
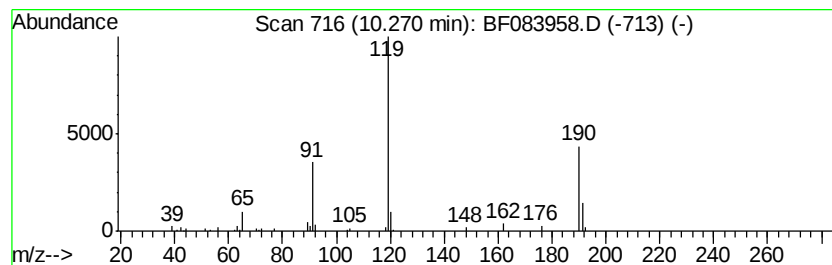
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Diethyltoluamide Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.27	3.23 ng	101556	Acenaphthene-d10	9.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Diethyltoluamide	191	C12H17NO	000134-62-3	96
2		Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	93
3		N-[1-(Dimethylamino)-2,2,2-trich...	308	C12H15Cl3N2O	300814-41-9	53
4		N-Isopropyl-p-toluamide	177	C11H15NO	002144-17-4	50
5		p-Toluric acid	193	C10H11NO3	027115-50-0	50



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Sample : G4875-01
Misc :
ALS Vial : 8 Sample Multiplier: 1

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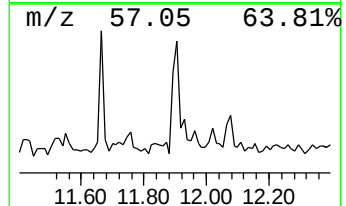
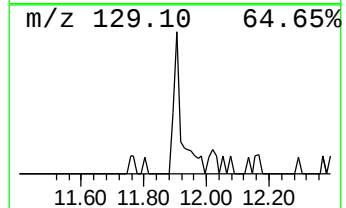
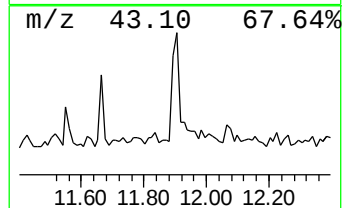
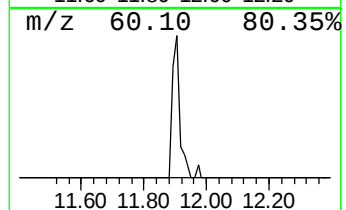
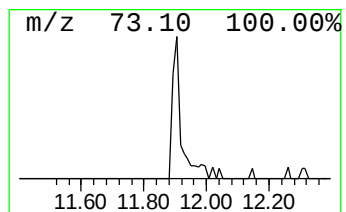
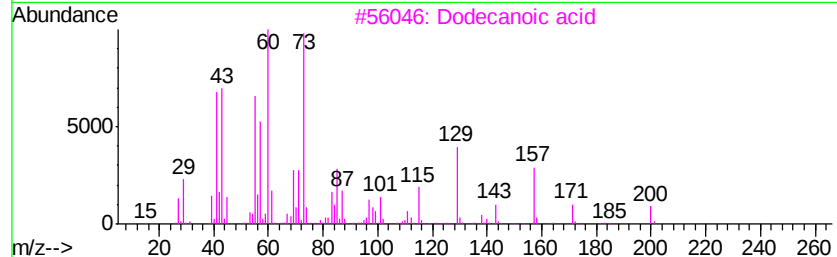
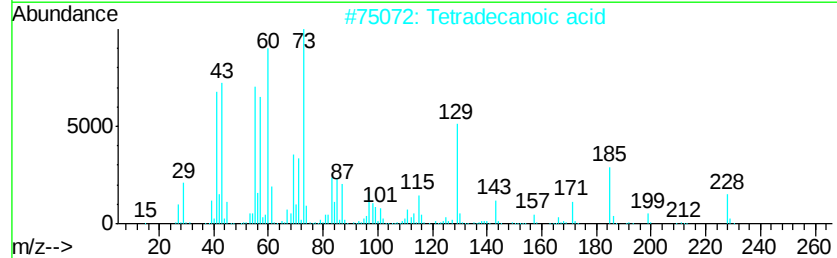
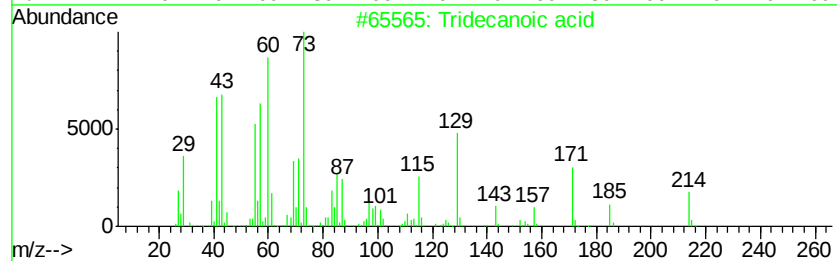
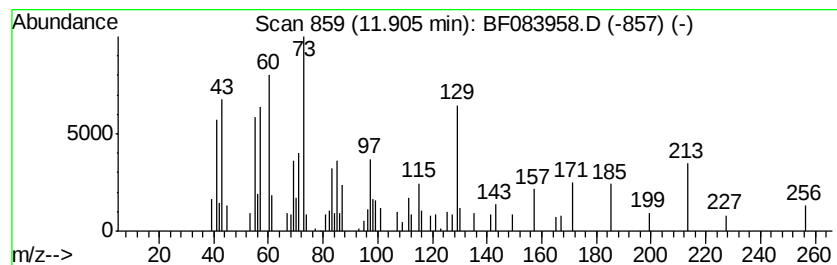
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Tridecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	2.60 ng	85076	Phenanthrene-d10	11.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tridecanoic acid	214	C13H26O2	000638-53-9	72
2		Tetradecanoic acid	228	C14H28O2	000544-63-8	64
3		Dodecanoic acid	200	C12H24O2	000143-07-7	50
4		n-Decanoic acid	172	C10H20O2	000334-48-5	43
5		.alpha.-D-Glucopyranose, 4-O-.be...	342	C12H22O11	014641-93-1	38



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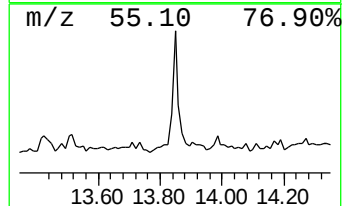
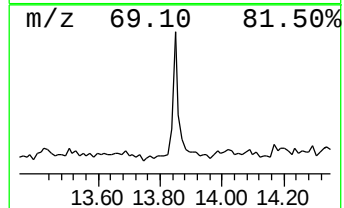
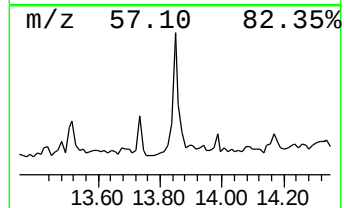
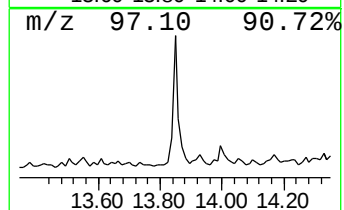
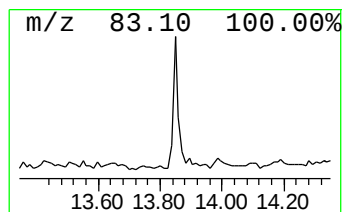
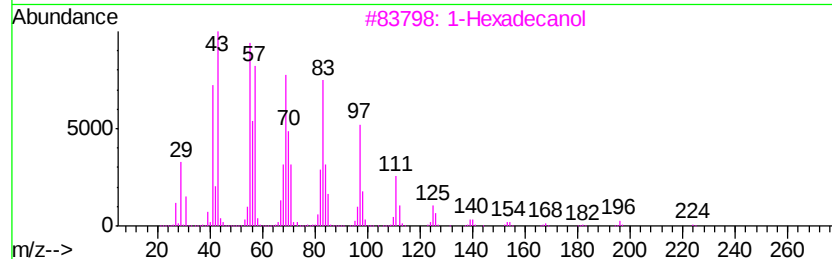
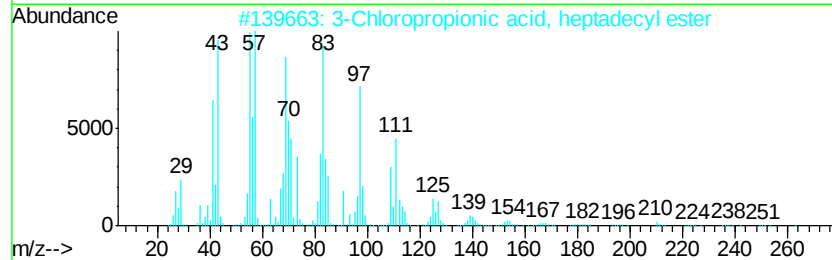
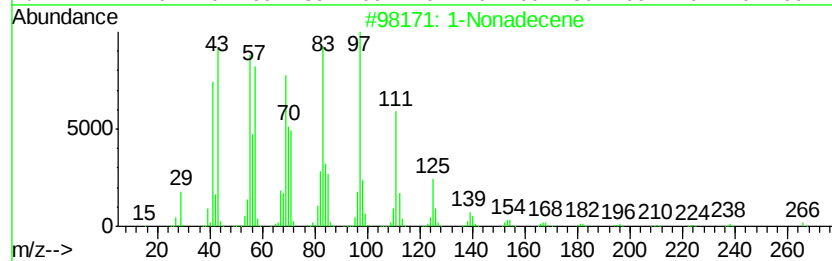
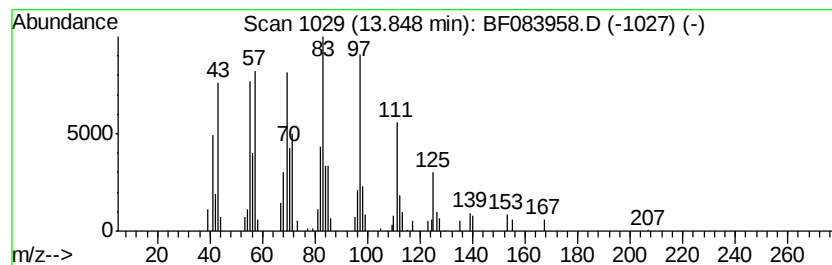
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 1-Nonadecene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.85	2.46 ng	82616	Chrysene-d12	14.00
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1-Nonadecene		266 C19H38	018435-45-5 95
2	3-Chloropropionic acid, heptadec...		346 C20H39ClO2	1000283-05-1 91
3	1-Hexadecanol		242 C16H34O	036653-82-4 91
4	Trifluoroacetic acid, n-heptadec...		324 C17H31F3O2	1000216-79-2 91
5	Trichloroacetic acid, tetradecyl...		358 C16H29Cl3O2	074339-52-9 90



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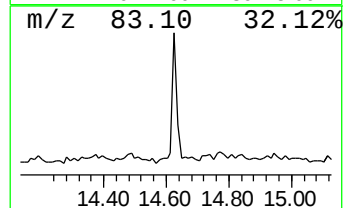
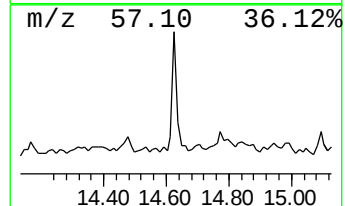
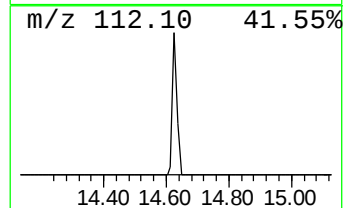
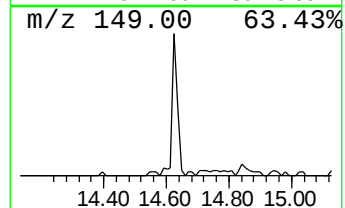
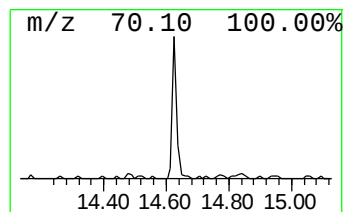
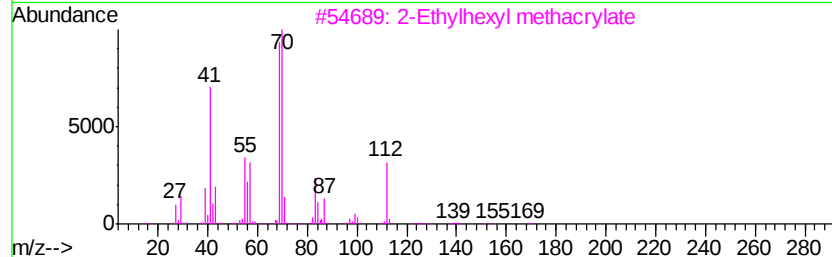
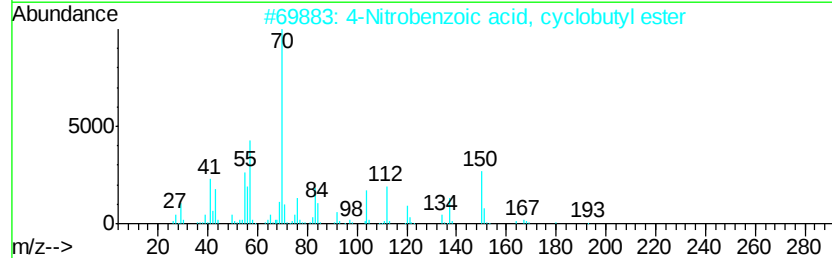
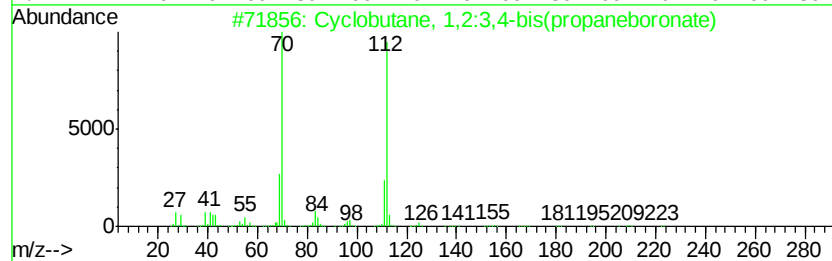
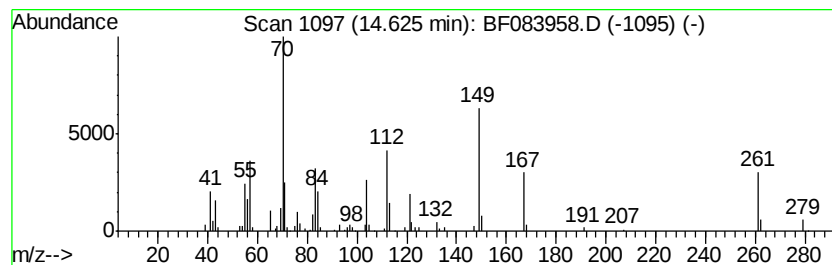
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 unknown14.63 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.63	3.42 ng	114792	Chrysene-d12	14.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclobutane, 1,2:3,4-bis(propene...	224	C10H18B2O4	1000159-70-4	32
2			4-Nitrobenzoic acid, cyclobutyl ...	221	C11H11N04	070335-00-1	32
3			2-Ethylhexyl methacrylate	198	C12H22O2	000688-84-6	32
4			Bromoacetic acid, 2-ethylhexyl e...	250	C10H19BrO2	068144-73-0	25
5			5-Pyrrolidino-2-pyrrolidone	154	C8H14N2O	076284-12-3	22



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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...	5.01	7.6 ng		170259	1	6.85	446561	20.0
unknown6.61	6.61	45.1 ng		1007270	1	6.85	446561	20.0
unknown8.27	8.27	2.6 ng		62068	2	8.13	475184	20.0
unknown8.30	8.30	2.7 ng		63686	2	8.13	475184	20.0
unknown8.44	8.44	5.4 ng		127716	2	8.13	475184	20.0
Thiophane, propyl-	8.54	14.7 ng		348758	2	8.13	475184	20.0
Pentanoic acid, 2...	8.59	6.6 ng		157033	2	8.13	475184	20.0
Diethyltoluamide	10.27	3.2 ng		101556	3	9.88	629383	20.0
Tridecanoic acid	11.90	2.6 ng		85076	4	11.37	655119	20.0
1-Nonadecene	13.85	2.5 ng		82616	5	14.00	672098	20.0
unknown14.63	14.63	3.4 ng		114792	5	14.00	672098	20.0