

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG091117\
 Data File : BG028658.D
 Acq On : 12 Sep 2017 2:24
 Operator : SJ/JU
 Sample : I5180-02
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 090717-TIDO-F-D-S

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_G\METHODS\8270-BG090717.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	5.445	308	316	327	rVB	43097	69680	1.79%	0.491%
2	5.909	388	395	406	rBV	221334	372901	9.60%	2.626%
3	7.489	657	664	675	rBV	124638	232455	5.98%	1.637%
4	7.866	721	728	739	rBV	387698	703600	18.11%	4.955%
5	8.336	802	808	817	rVB2	132300	237739	6.12%	1.674%
6	8.659	854	863	872	rBV	544762	991944	25.53%	6.986%
7	9.505	999	1007	1017	rBV	452279	848450	21.84%	5.975%
8	11.150	1279	1287	1298	rBV	170781	336129	8.65%	2.367%
9	13.565	1690	1698	1709	rBV	1397979	2168805	55.81%	15.274%
10	14.945	1925	1933	1942	rBV2	315513	528538	13.60%	3.722%
11	16.432	2179	2186	2201	rBV	774481	1276914	32.86%	8.993%
12	17.695	2393	2401	2407	rBV2	479148	736612	18.96%	5.188%
13	20.280	2834	2841	2850	rBV	2858624	3885704	100.00%	27.365%
14	22.025	3131	3138	3147	rBV	501766	902511	23.23%	6.356%
15	25.533	3716	3735	3748	rBV2	274843	907626	23.36%	6.392%

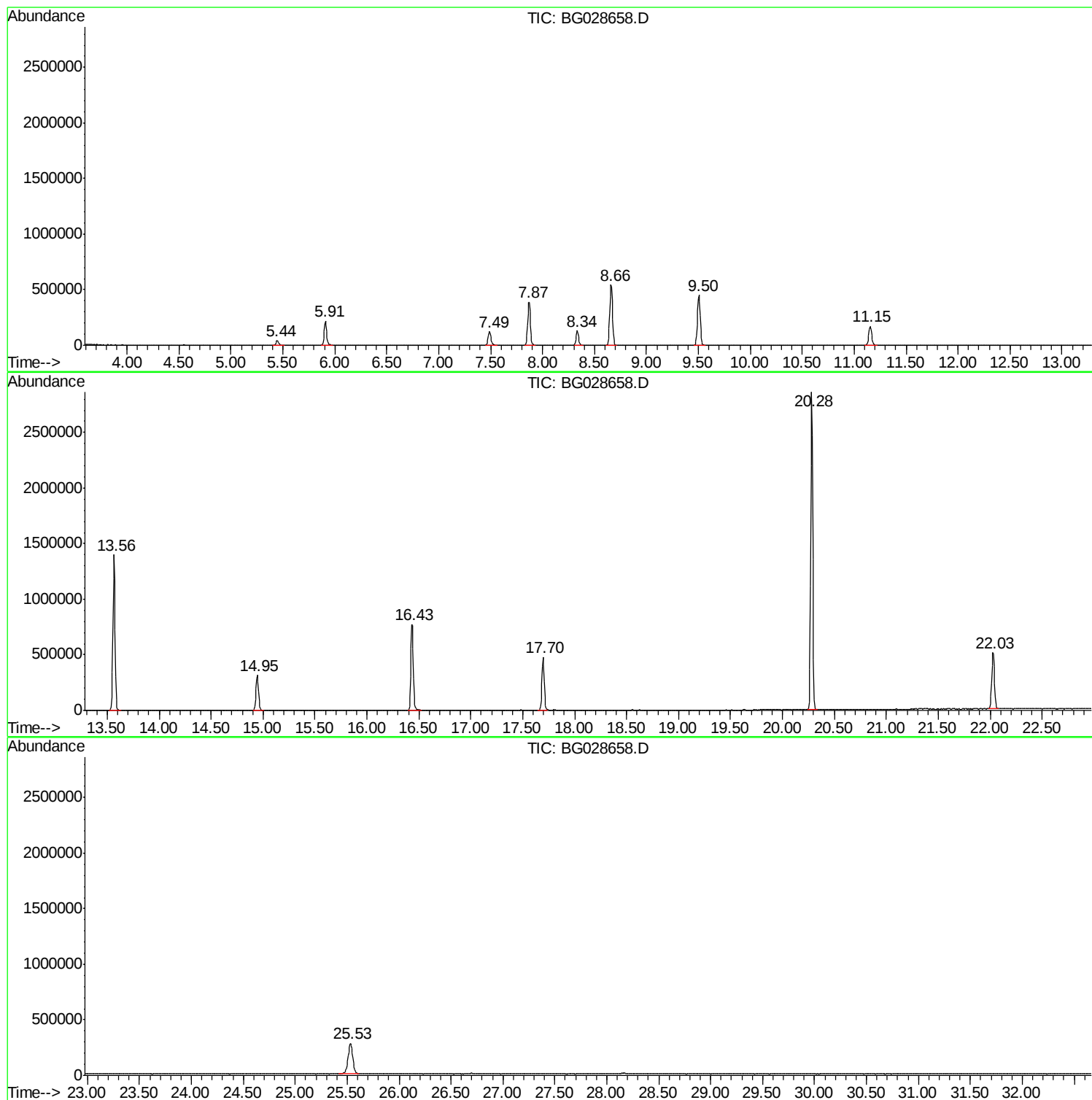
Sum of corrected areas: 14199608

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG090717.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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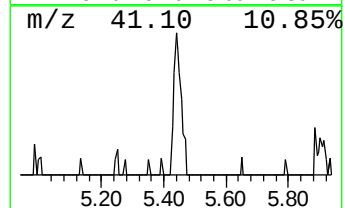
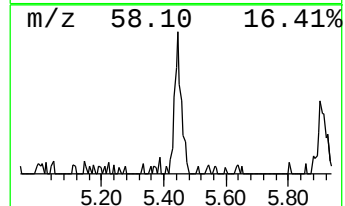
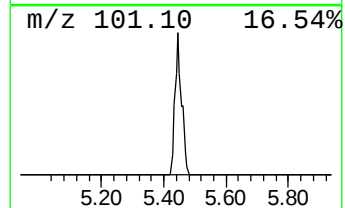
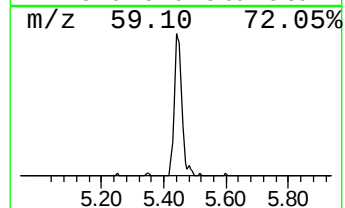
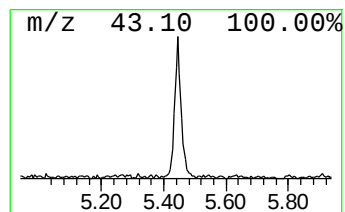
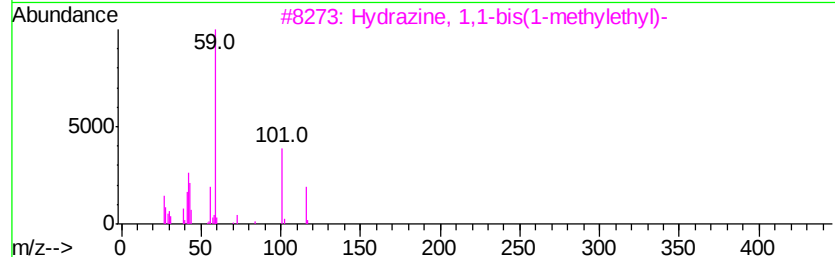
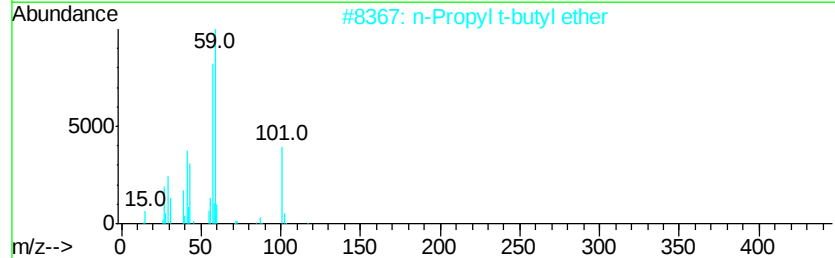
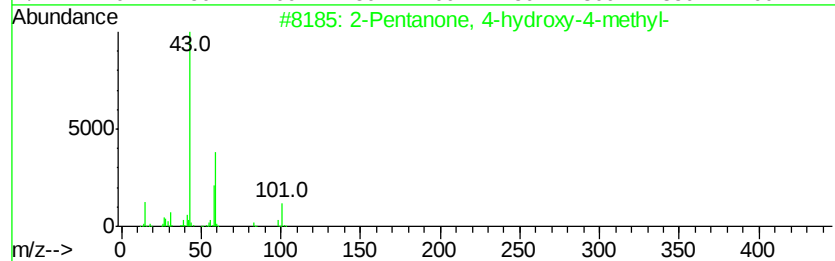
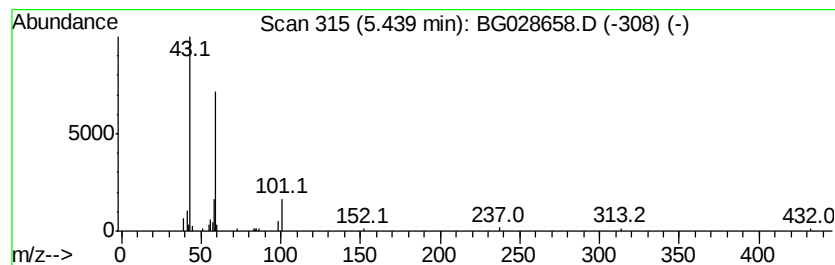
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.44	5.86 ng	69680	1,4-Dichlorobenzene-d4	8.34

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2			n-Propyl t-butyl ether	116	C7H16O	1000334-81-9	28
3			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	25
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
5			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17



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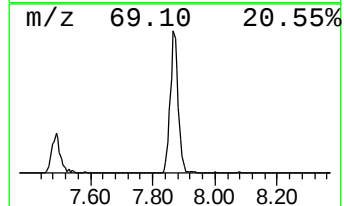
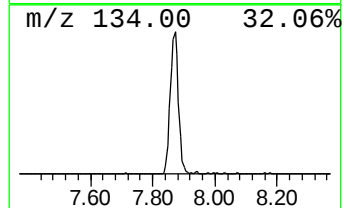
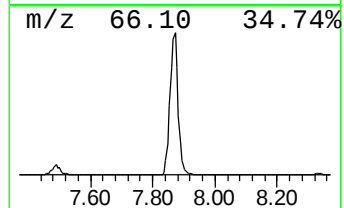
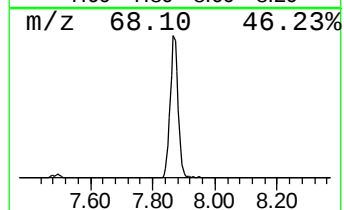
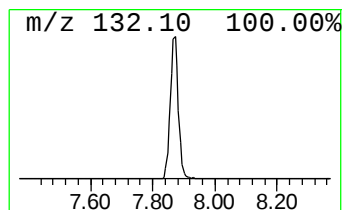
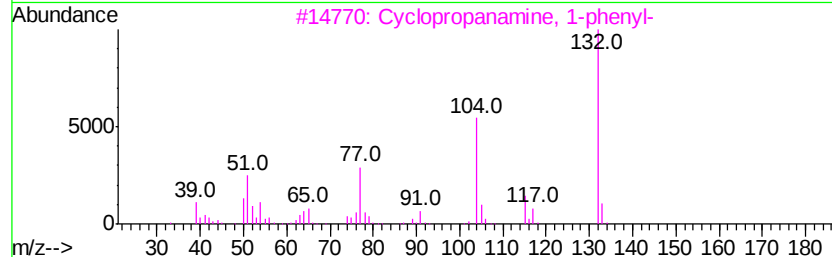
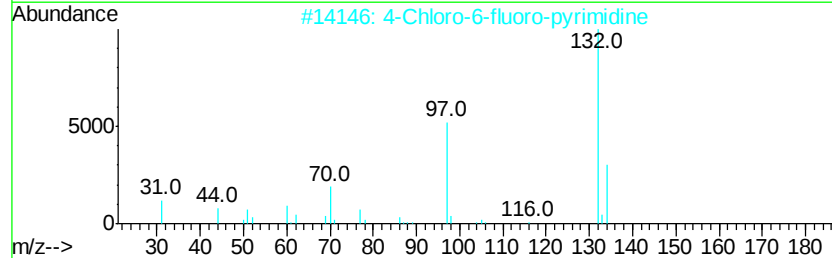
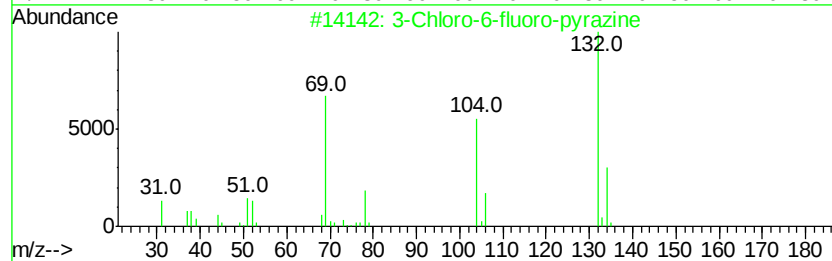
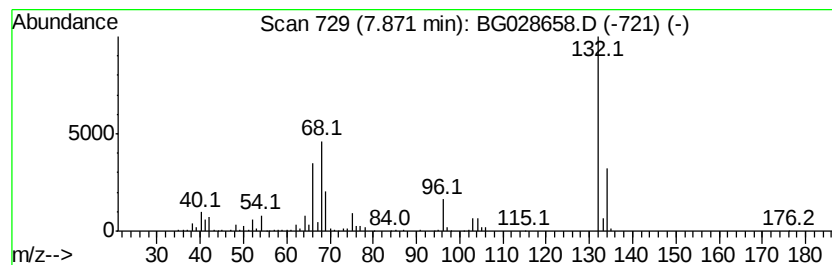
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Peak Number 2 unknown7.87 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.87	59.19 ng	703600	1,4-Dichlorobenzene-d4	8.34

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	22
4		Tranvlcypropamine-propionyl	189	C12H15NO	1000123-86-3	22
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.44	5.9	ng	69680	1	8.34	237739	20.0
unknown7.87	7.87	59.2	ng	703600	1	8.34	237739	20.0