

Data Path : Z:\HPCHEM1\BNA G\Data\BG042117\
 Data File : BG026689.D
 Acq On : 21 Apr 2017 16:52
 Operator : SJ/MA
 Sample : PB98386BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB98386BL

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-BG041817.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	5.127	341	347	359	rVB	250245	371625	2.94%	0.616%
2	5.615	423	430	442	rBV	2720024	4344045	34.34%	7.205%
3	7.219	694	703	717	rBV	2527937	4342262	34.33%	7.202%
4	7.571	756	763	780	rBV	2775716	4759553	37.63%	7.894%
5	8.029	833	841	850	rBV	532717	901173	7.12%	1.495%
6	8.353	887	896	906	rBV	2126876	3703769	29.28%	6.143%
7	9.187	1030	1038	1050	rBV	1513066	2673506	21.14%	4.434%
8	10.826	1308	1317	1329	rBV	740384	1362132	10.77%	2.259%
9	13.265	1724	1732	1745	rBV	4650363	7411864	58.60%	12.293%
10	14.639	1959	1966	1982	rBV2	1239520	1958093	15.48%	3.248%
11	16.120	2211	2218	2236	rBV	3925094	6344100	50.15%	10.522%
12	17.371	2424	2431	2441	rBV2	1858045	2754266	21.77%	4.568%
13	19.974	2867	2874	2885	rBV	9090461	12649289	100.00%	20.979%
14	21.619	3147	3154	3163	rBV	2090092	3348526	26.47%	5.554%
15	24.792	3683	3694	3706	rBV2	1186168	3370914	26.65%	5.591%

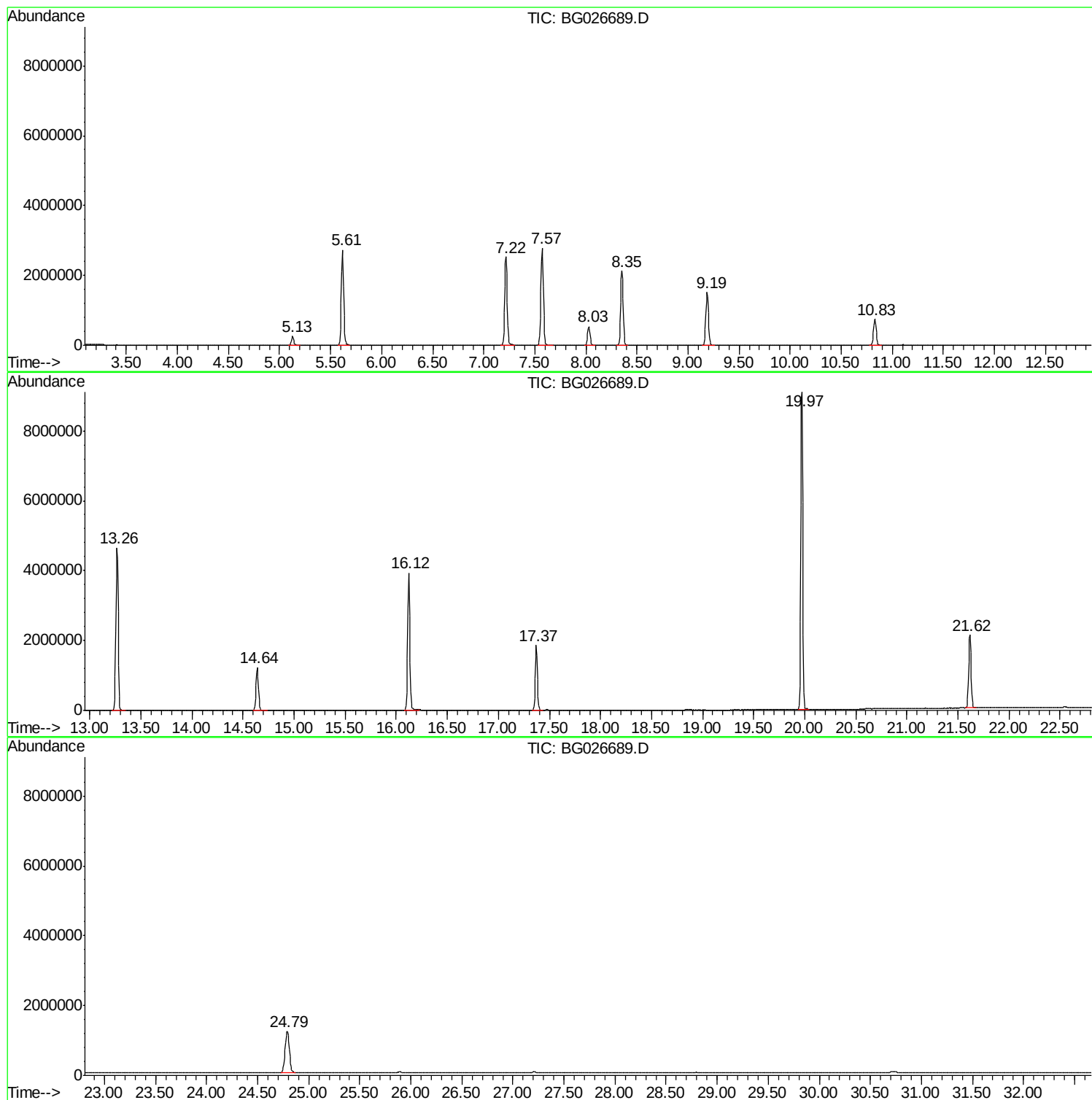
Sum of corrected areas: 60295117

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG041817.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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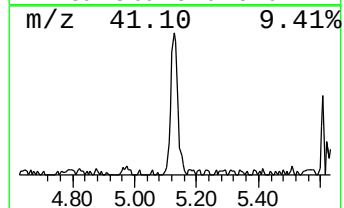
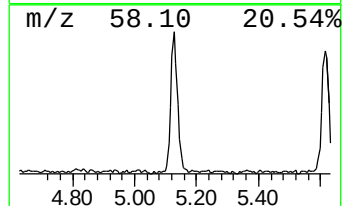
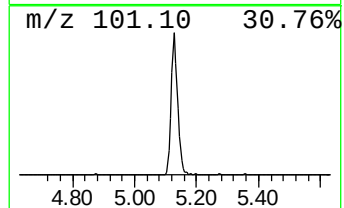
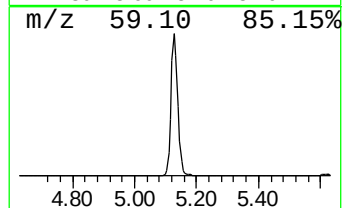
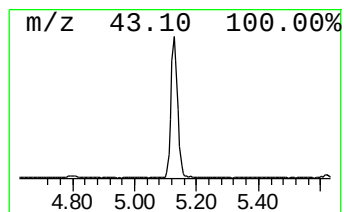
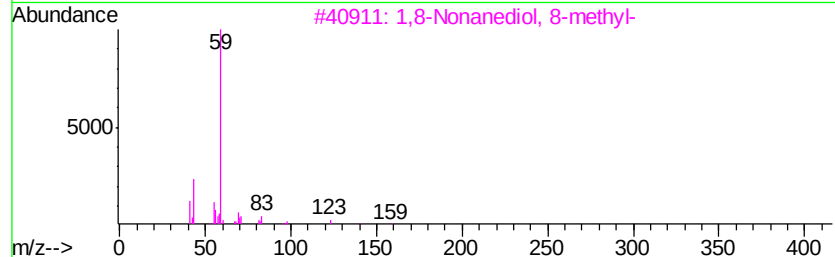
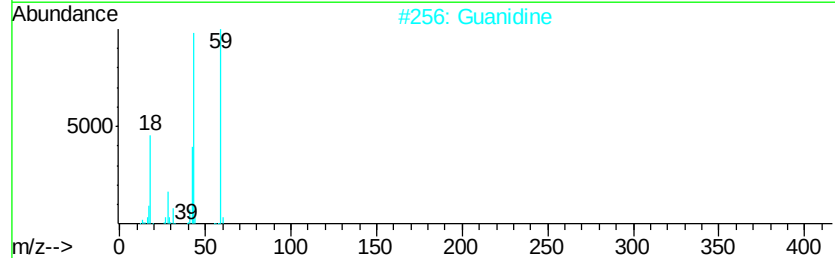
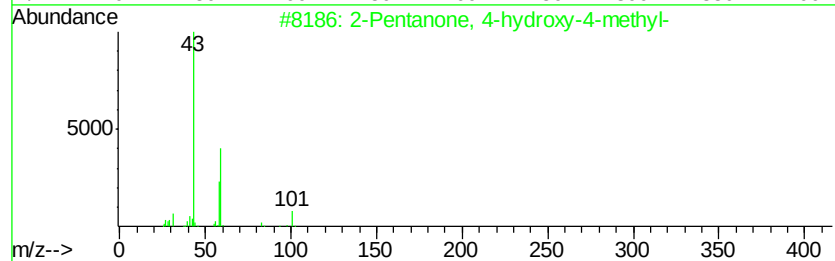
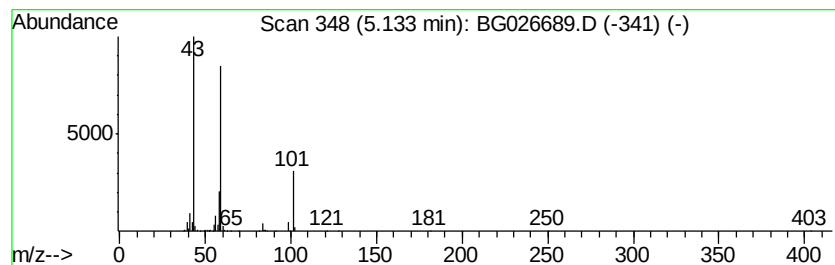
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	8.25 ng	371625	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Guanidine	59	CH5N3	000113-00-8	47
3		1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	28
4		(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	27
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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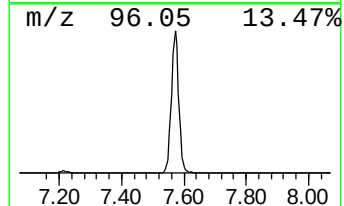
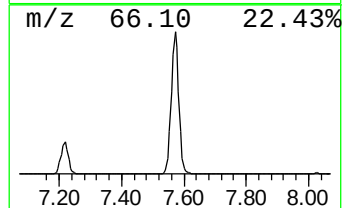
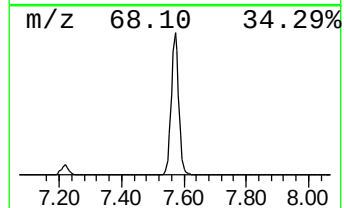
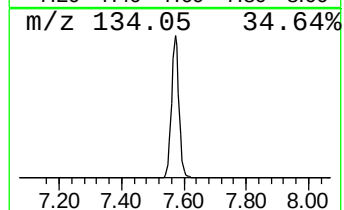
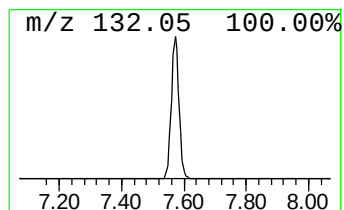
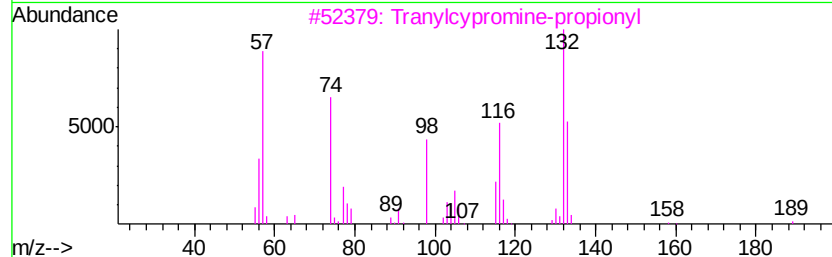
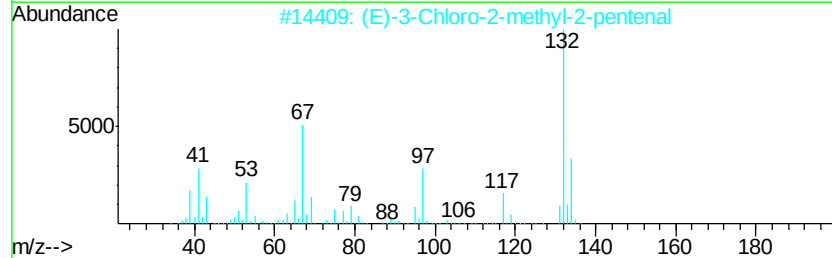
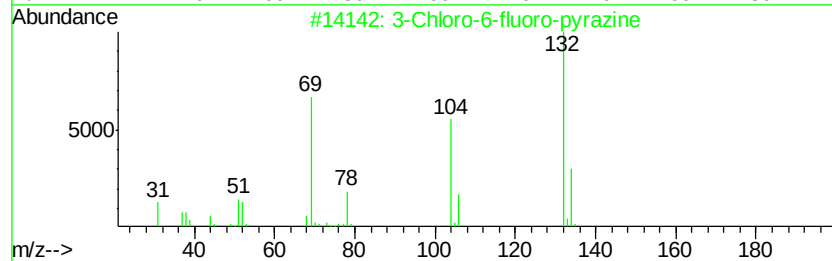
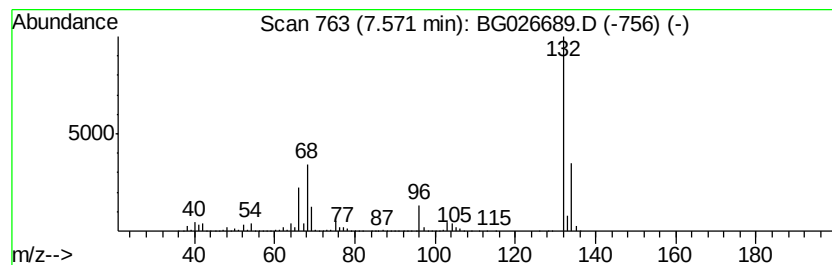
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	105.63 ng	4759550	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	12
4		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	8.3	ng	371625	1	8.03	901173	20.0
unknown7.57	7.57	105.6	ng	4759550	1	8.03	901173	20.0