

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022717\
 Data File : BG025966.D
 Acq On : 27 Feb 2017 16:33
 Operator : SJ/MA
 Sample : PB97190BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB97190BL

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-BG021717.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	5.176	306	313	324	rBV	545975	801859	9.68%	2.157%
2	5.646	385	393	404	rBV	1628867	2528959	30.52%	6.802%
3	7.232	655	663	677	rBV	1607079	2672330	32.25%	7.188%
4	7.608	719	727	738	rBV	1555849	2638717	31.84%	7.097%
5	8.078	800	807	814	rVB	277592	457091	5.52%	1.229%
6	8.401	854	862	872	rVB	1092871	1871998	22.59%	5.035%
7	9.230	995	1003	1012	rBV	895940	1558606	18.81%	4.192%
8	10.875	1275	1283	1293	rVB	416713	751291	9.07%	2.021%
9	13.313	1690	1698	1705	rBV	2679946	4296482	51.85%	11.556%
10	14.676	1922	1930	1939	rBV3	710664	1176610	14.20%	3.165%
11	16.157	2175	2182	2198	rBV	2656200	4045642	48.82%	10.881%
12	17.409	2388	2395	2407	rVB2	1207936	1806111	21.80%	4.858%
13	20.005	2831	2837	2843	rBV	6299065	8286637	100.00%	22.288%
14	21.656	3111	3118	3126	rVB	1306338	2131994	25.73%	5.734%
15	24.853	3651	3662	3676	rBV2	753681	2155311	26.01%	5.797%

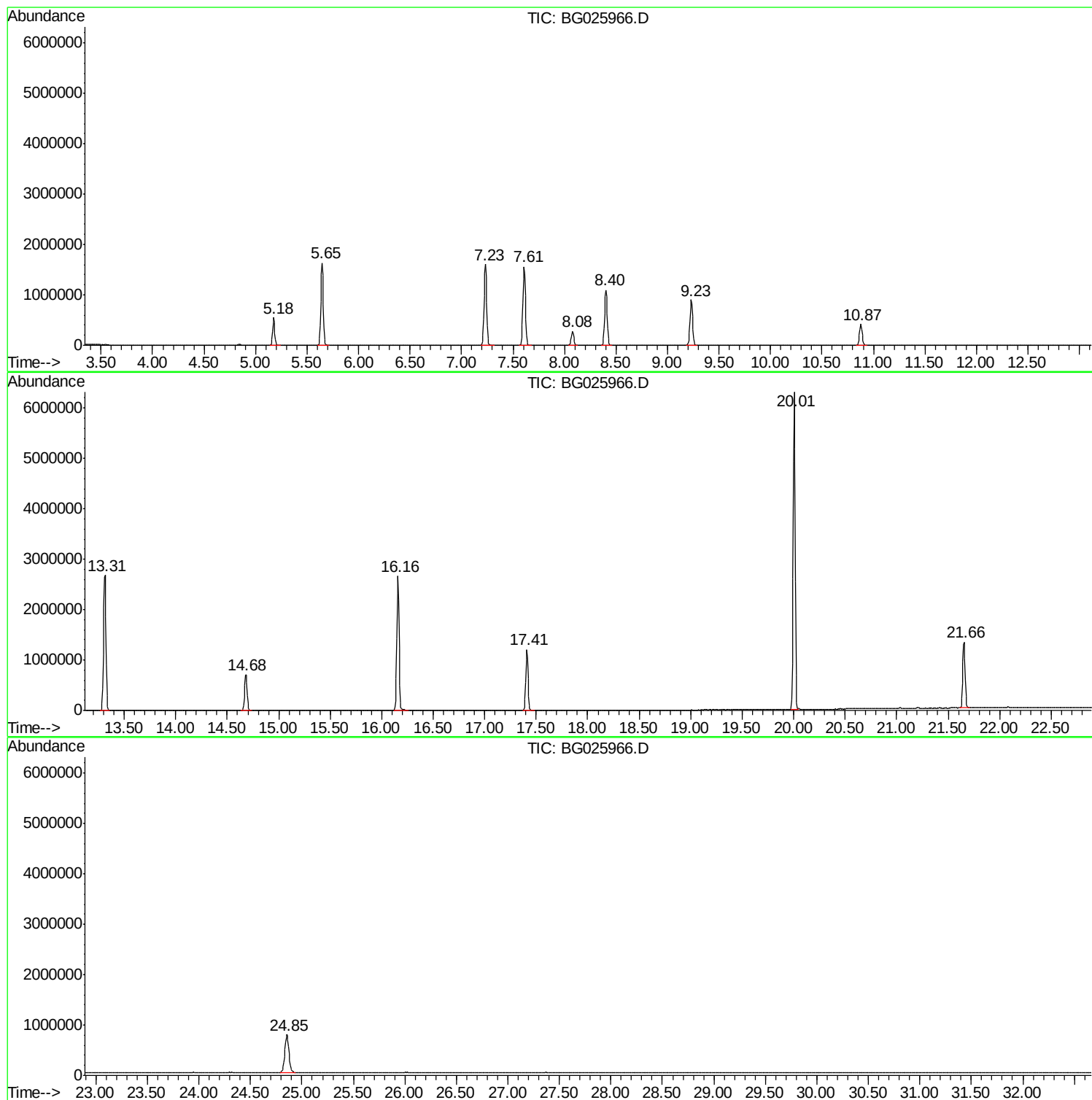
Sum of corrected areas: 37179638

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021717.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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Instrument :
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ClientSampled :
PB97190BL

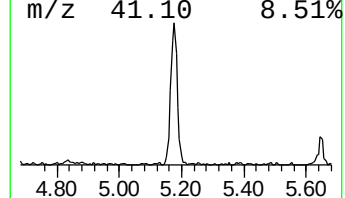
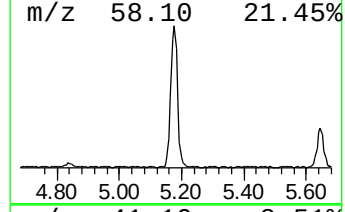
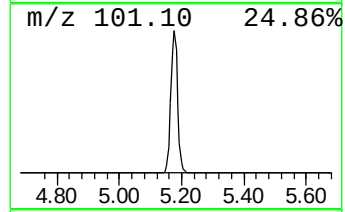
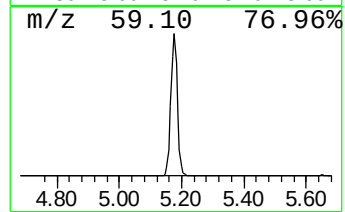
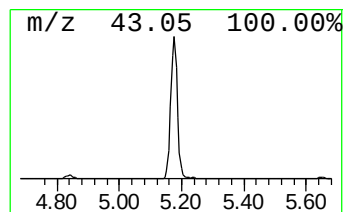
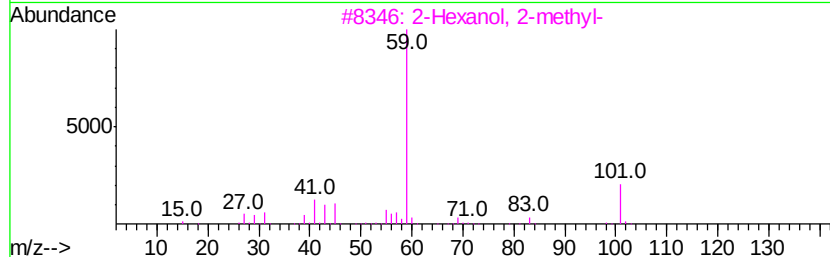
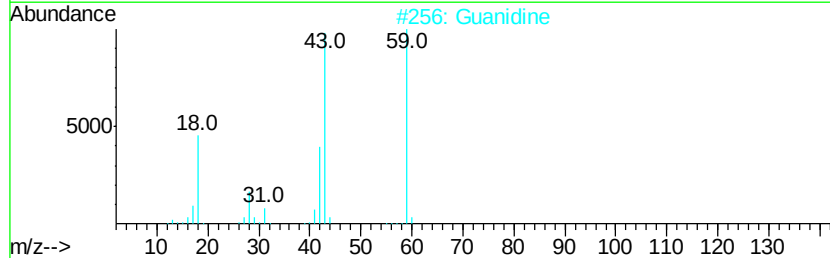
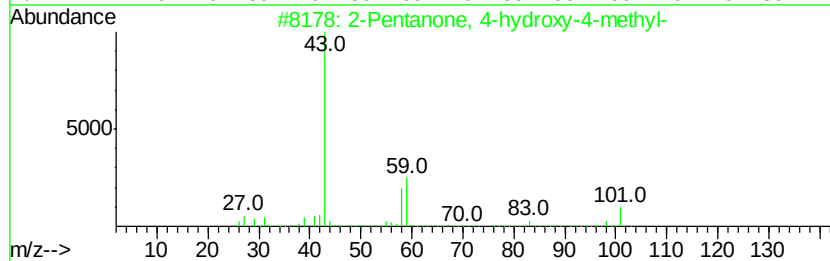
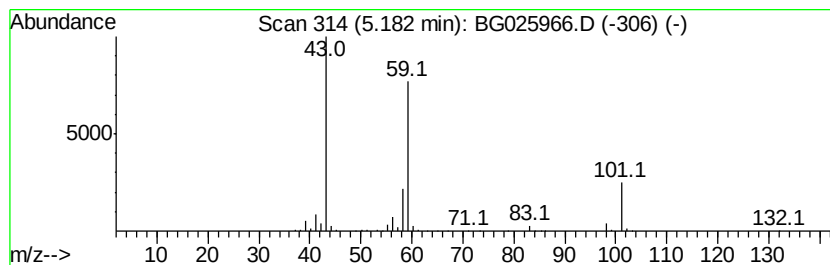
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	35.09 ng	801859	1,4-Dichlorobenzene-d4	8.08

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Guanidine	59	CH5N3	000113-00-8	47
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	36
4		Propane, 2-methyl-2-(1-methylethyl)-	116	C7H16O	017348-59-3	36
5		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	25



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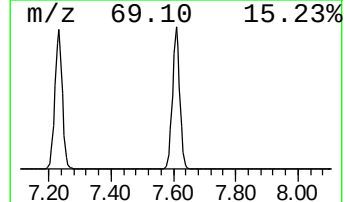
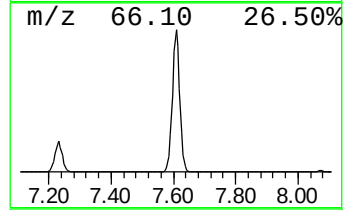
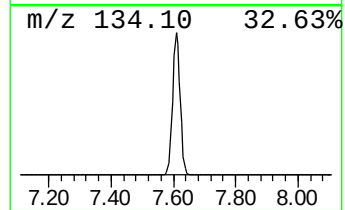
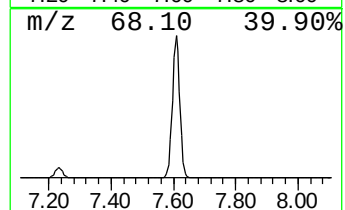
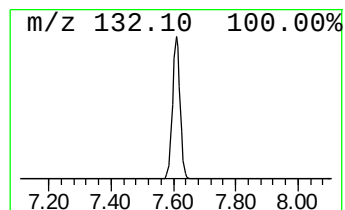
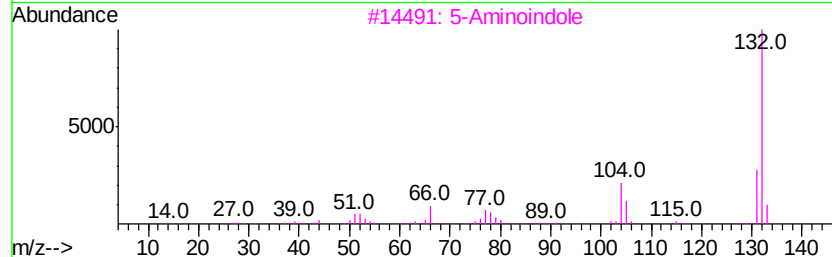
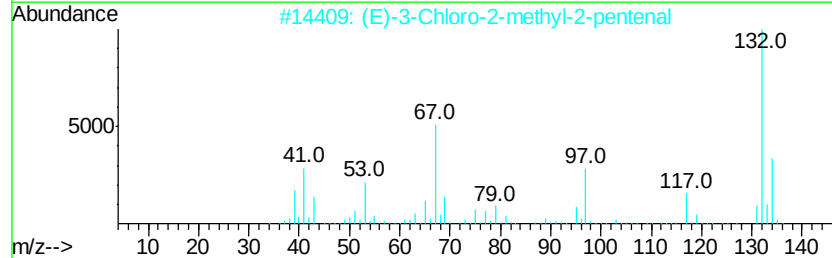
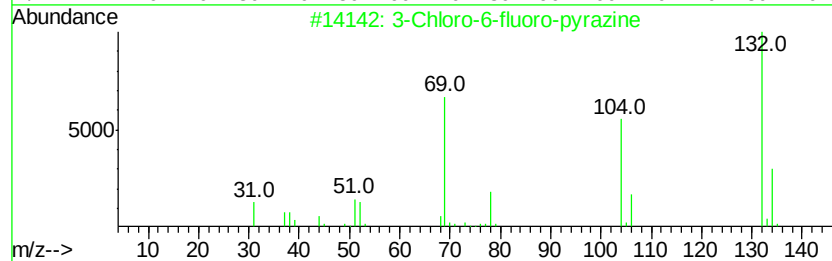
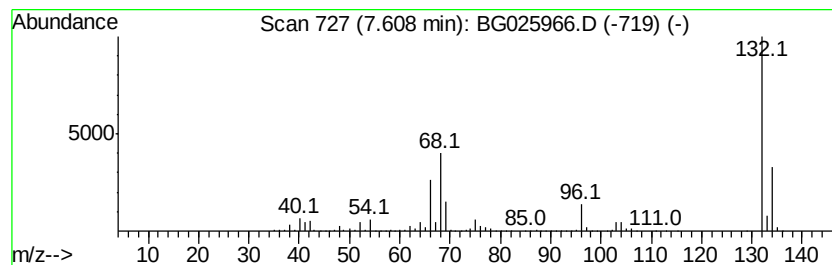
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	115.46 ng	2638720	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	35.1	ng	801859	1	8.08	457091	20.0
unknown7.61	7.61	115.5	ng	2638720	1	8.08	457091	20.0