

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022417\
 Data File : BG025940.D
 Acq On : 24 Feb 2017 20:19
 Operator : SJ/MA
 Sample : I1957-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FB20170222

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	5.180	306	313	321	rBV	195661	287231	3.03%	0.790%
2	5.650	385	393	402	rVB	752209	1183668	12.47%	3.257%
3	7.231	655	662	670	rBV	449628	756816	7.98%	2.083%
4	7.612	719	727	736	rBV	1361994	2312381	24.37%	6.363%
5	8.083	800	807	814	rBV	289598	497338	5.24%	1.369%
6	8.406	853	862	870	rBV	1264715	2159627	22.76%	5.943%
7	9.234	994	1003	1015	rBV	1023981	1818007	19.16%	5.003%
8	10.879	1274	1283	1293	rVB	443007	796227	8.39%	2.191%
9	13.312	1688	1697	1705	rBV	3230026	5057989	53.30%	13.918%
10	14.687	1922	1931	1939	rVB2	753675	1252495	13.20%	3.447%
11	16.161	2174	2182	2191	rBV	2969155	4531351	47.75%	12.469%
12	17.413	2388	2395	2404	rVB2	1211324	1845829	19.45%	5.079%
13	20.010	2831	2837	2844	rBV	6904789	9489150	100.00%	26.111%
14	21.661	3111	3118	3124	rBV	1383676	2184625	23.02%	6.011%
15	24.869	3654	3664	3676	rVB2	757435	2168232	22.85%	5.966%

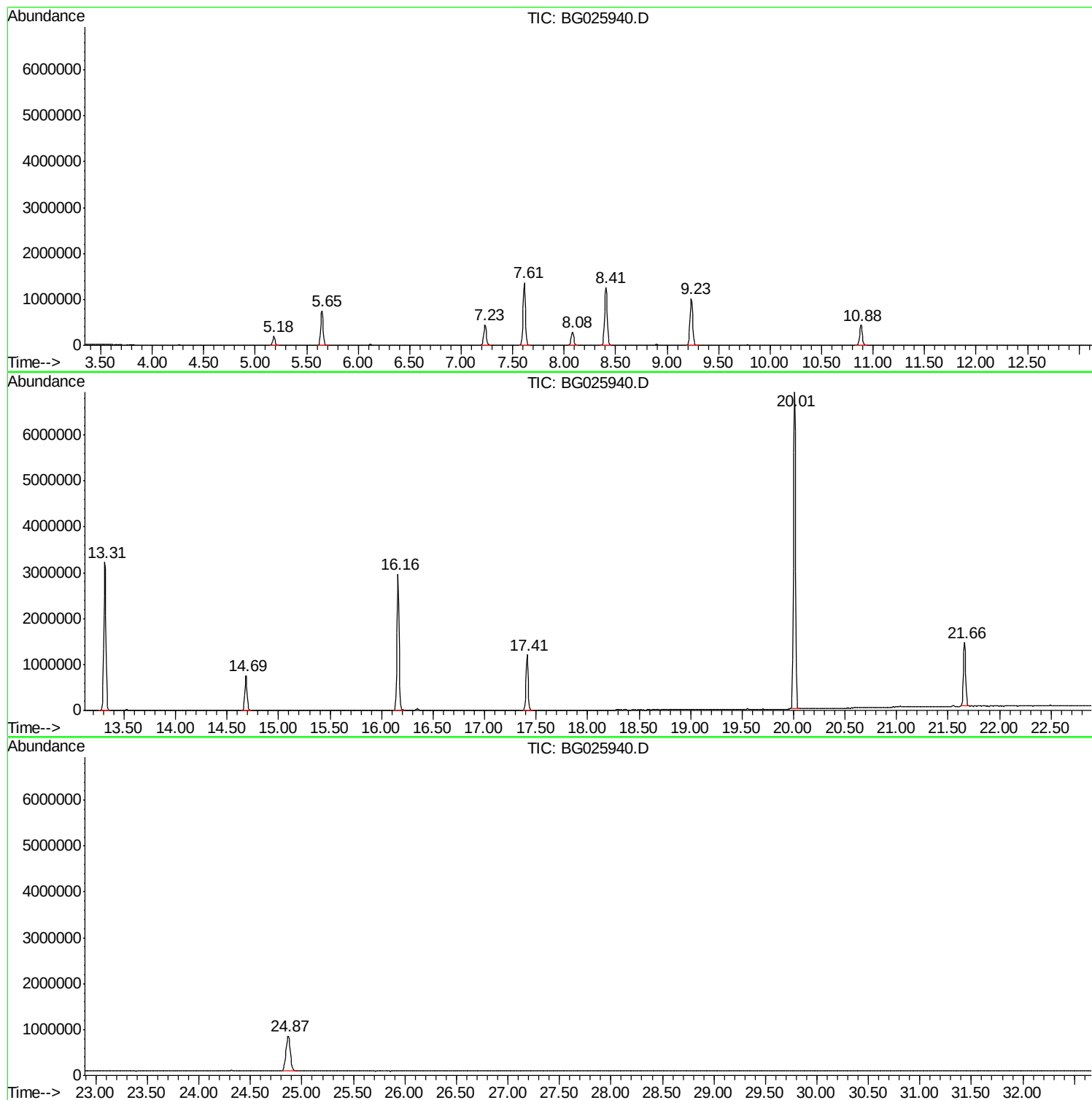
Sum of corrected areas: 36340966

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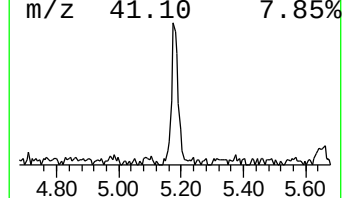
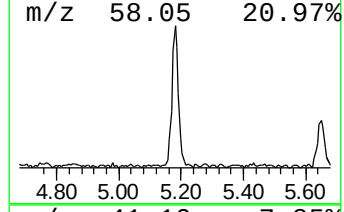
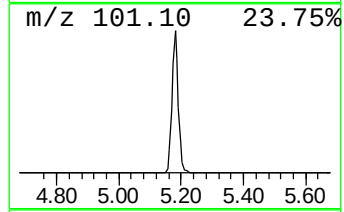
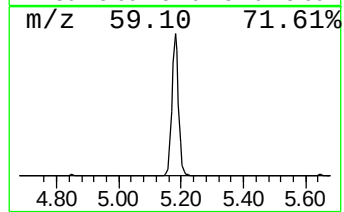
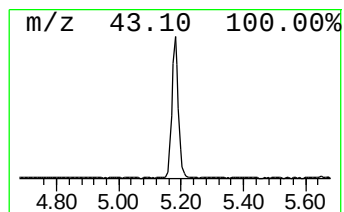
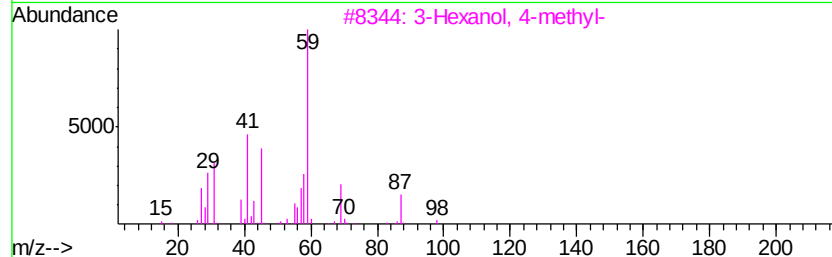
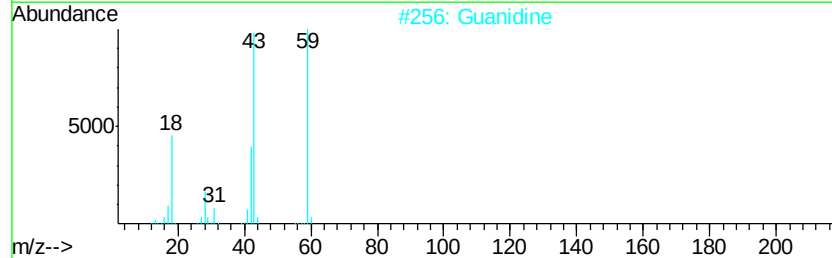
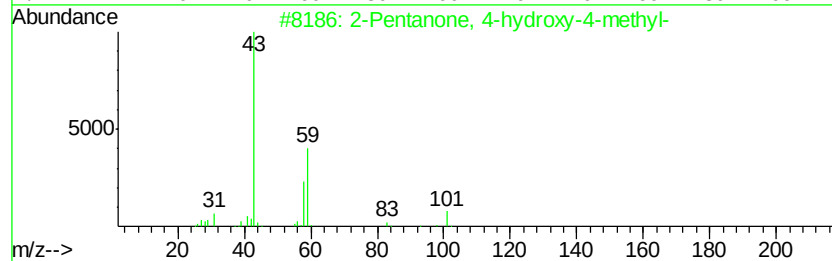
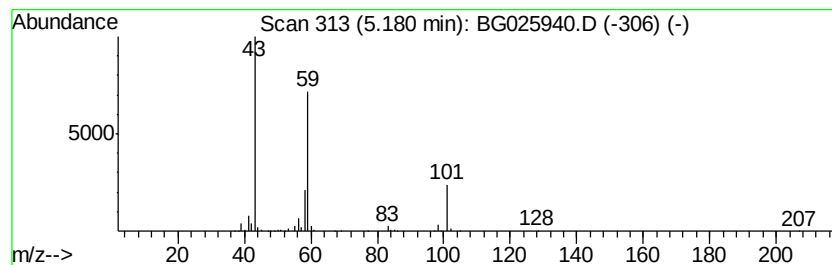
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	11.55 ng	287231	1,4-Dichlorobenzene-d4	8.08

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Guanidine	59	CH5N3	000113-00-8	32
3		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28
4		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	12
5		Acetone	58	C3H6O	000067-64-1	9



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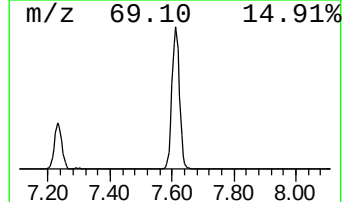
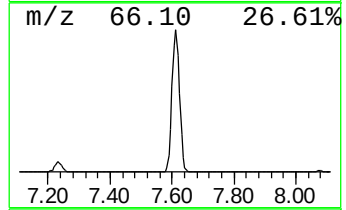
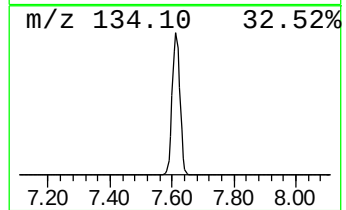
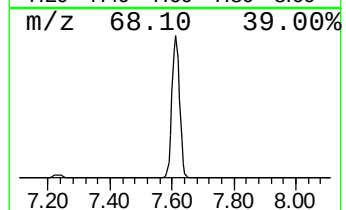
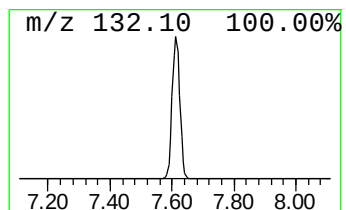
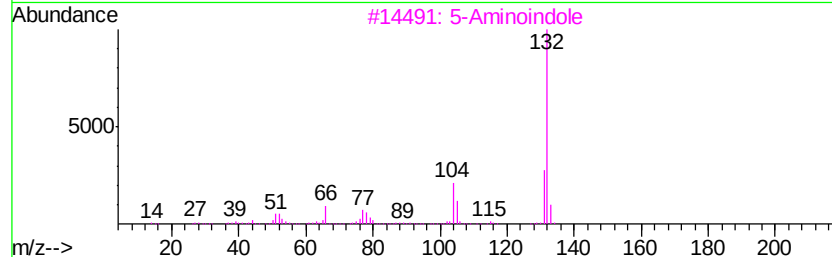
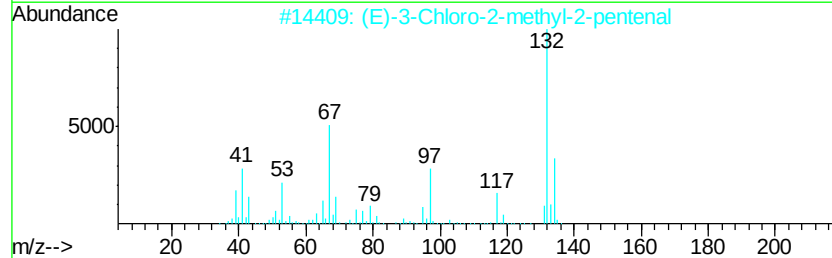
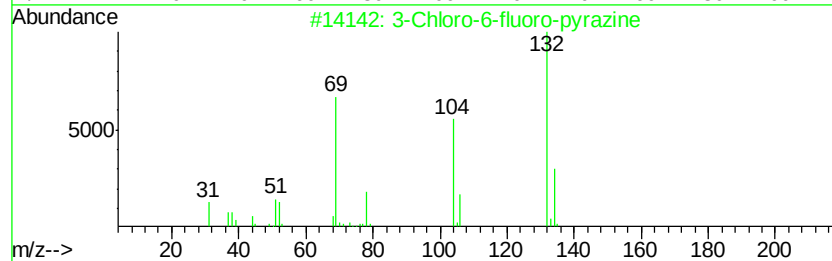
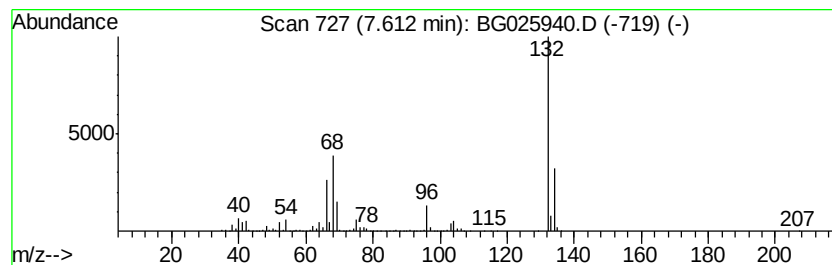
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Peak Number 2 unknown7.61 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.61	92.99 ng	2312380	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	16
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	12
3		5-Aminoindole	132	C8H8N2	005192-03-0	12
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	11.6	ng	287231	1	8.08	497338	20.0
unknown7.61	7.61	93.0	ng	2312380	1	8.08	497338	20.0