

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060616\
 Data File : BG022181.D
 Acq On : 6 Jun 2016 23:07
 Operator : UM/SJ
 Sample : PB91112BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB91112BL

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-BG060616.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	5.190	393	400	408	rBV	26270	50101	2.39%	0.571%
2	5.666	474	481	501	rBV	294608	565145	27.00%	6.437%
3	7.276	748	755	771	rBV	309183	611178	29.20%	6.961%
4	7.646	810	818	834	rBV	334094	664369	31.74%	7.567%
5	8.116	892	898	908	rBV	62603	128233	6.13%	1.461%
6	8.445	944	954	974	rBV	263922	535810	25.60%	6.103%
7	9.291	1090	1098	1112	rBV	155707	344110	16.44%	3.919%
8	10.942	1371	1379	1397	rBV	88250	201614	9.63%	2.296%
9	13.369	1784	1792	1807	rBV	617964	1104016	52.74%	12.575%
10	14.750	2020	2027	2037	rBV2	174449	318812	15.23%	3.631%
11	16.236	2273	2280	2295	rBV	442756	788775	37.68%	8.984%
12	17.494	2486	2494	2505	rBV	229254	403234	19.26%	4.593%
13	20.085	2928	2935	2946	rBV	1415751	2093282	100.00%	23.842%
14	21.765	3214	3221	3241	rBV	244336	496748	23.73%	5.658%
15	25.055	3769	3781	3799	rBV	144744	474335	22.66%	5.403%

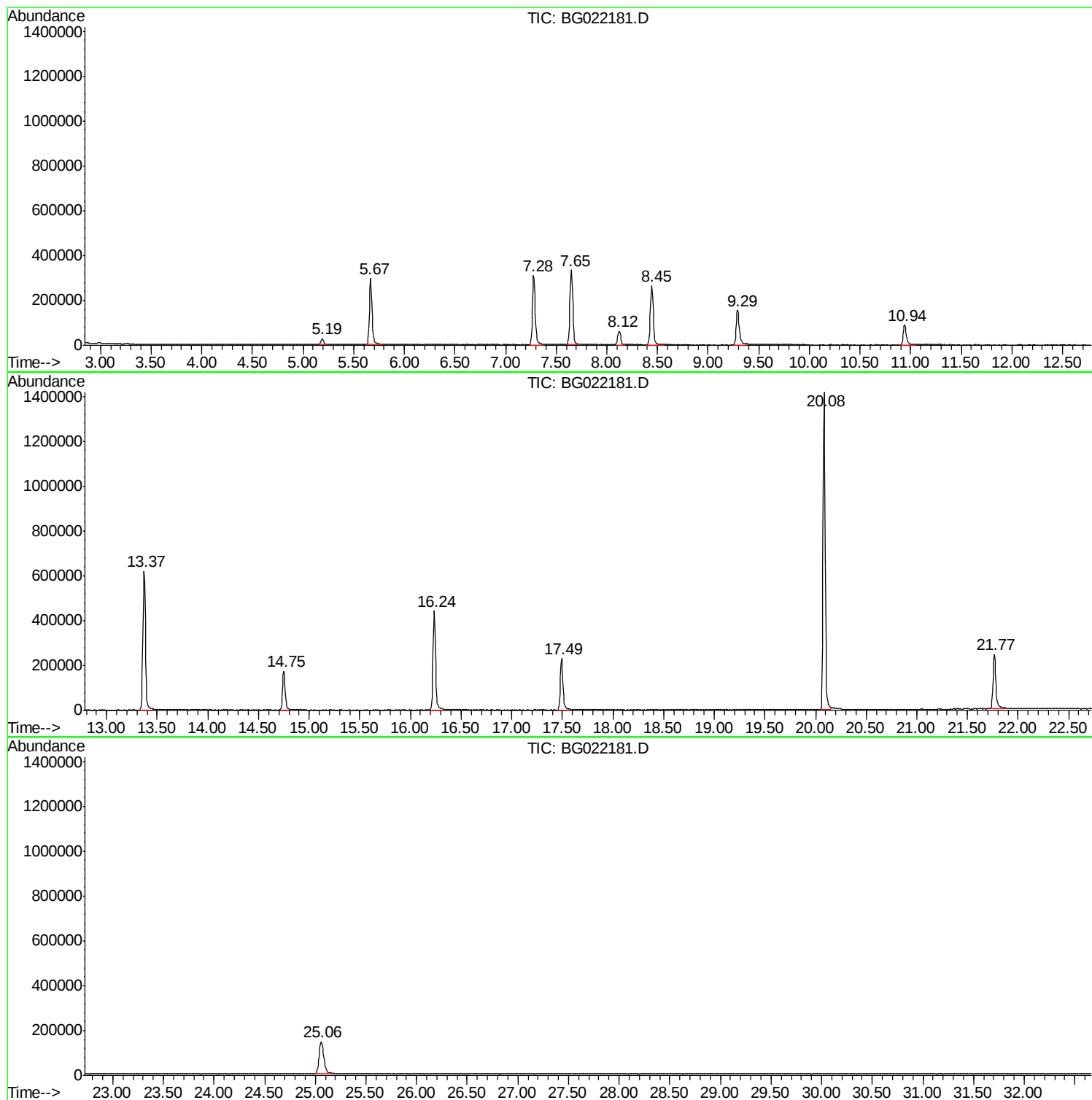
Sum of corrected areas: 8779762

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060616.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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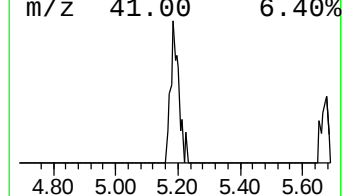
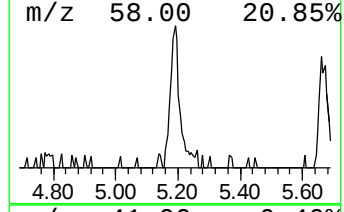
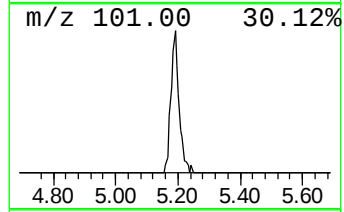
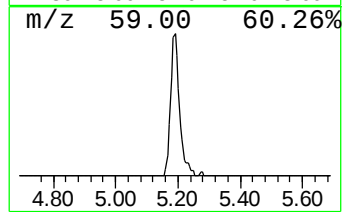
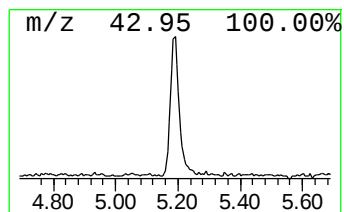
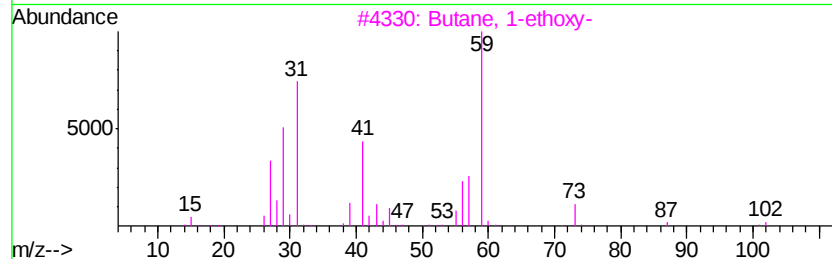
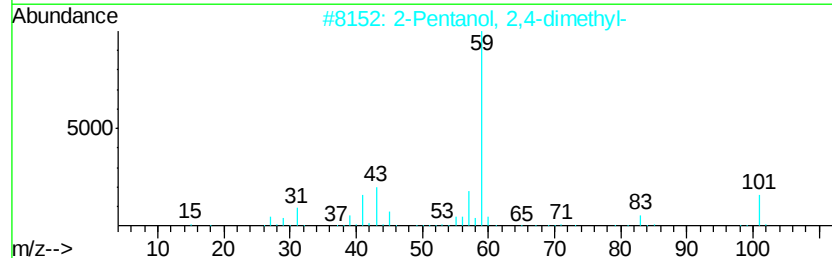
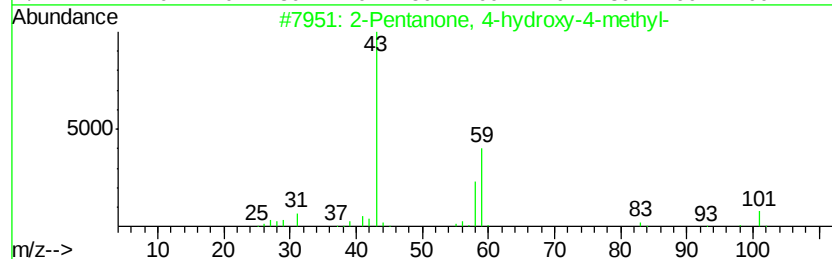
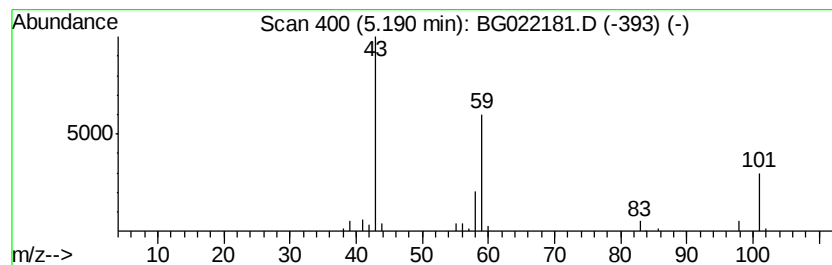
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	7.81 ng	50101	1,4-Dichlorobenzene-d4	8.12

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59	
2	2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	47	
3	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	10	
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9	
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9	



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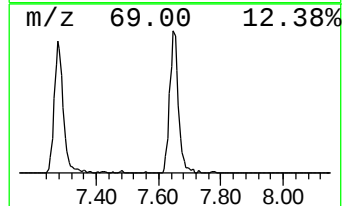
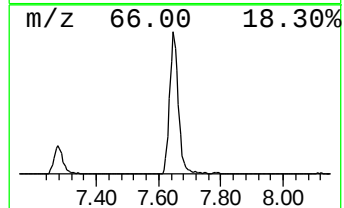
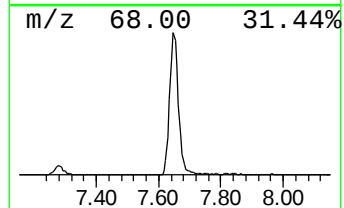
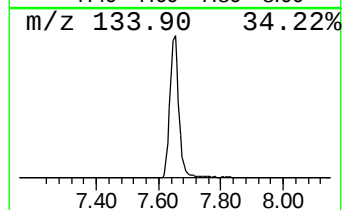
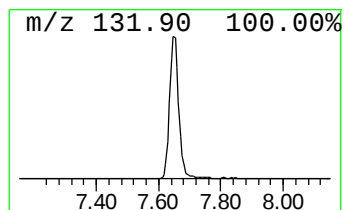
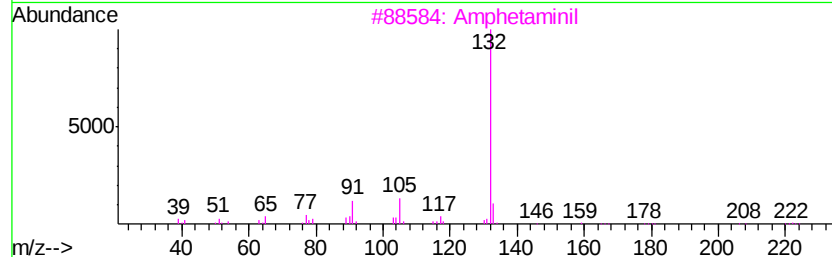
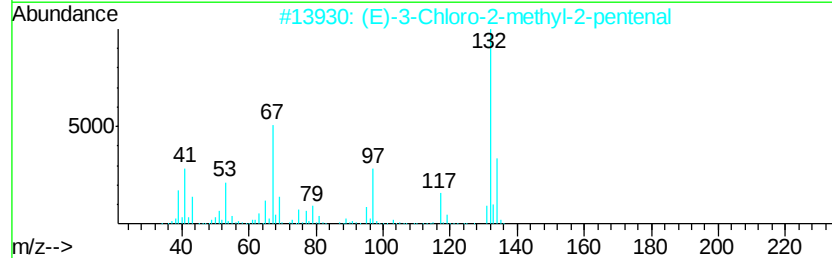
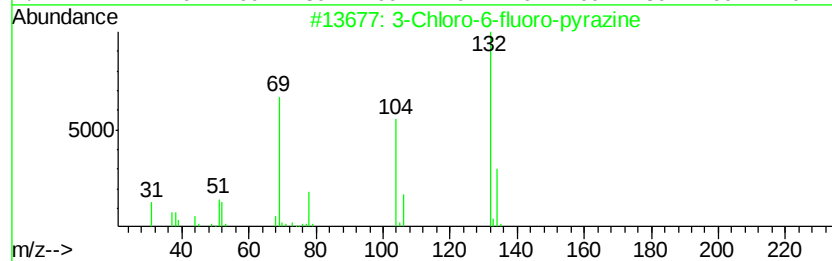
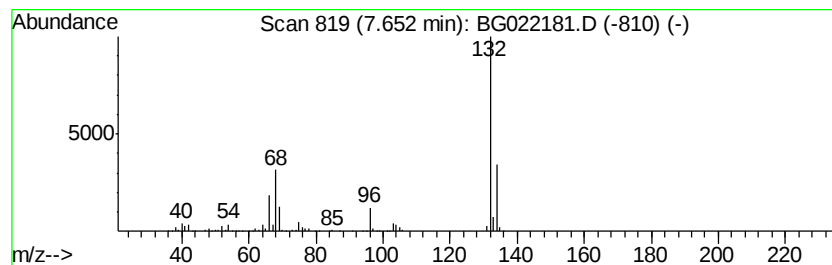
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Peak Number 2 unknown7.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.65	103.62 ng	664369	1,4-Dichlorobenzene-d4	8.12

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	25
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
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...	5.19	7.8	ng	50101	1	8.12	128233	20.0
unknown7.65	7.65	103.6	ng	664369	1	8.12	128233	20.0