

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061716\
 Data File : BG022418.D
 Acq On : 18 Jun 2016 11:20
 Operator : UM/SJ
 Sample : PB91233BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB91233BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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	2.868	3	5	16	rVB	28287	40080	2.18%	0.477%
2	5.101	378	385	394	rBV	14003	26693	1.45%	0.318%
3	5.600	463	470	486	rBV	278657	542504	29.51%	6.455%
4	7.228	740	747	765	rBV	283070	598276	32.54%	7.119%
5	7.580	798	807	821	rBV	330417	641091	34.87%	7.628%
6	8.038	878	885	899	rBV	64256	127476	6.93%	1.517%
7	8.362	932	940	952	rBV	251188	502657	27.34%	5.981%
8	9.219	1078	1086	1102	rBV	148414	346370	18.84%	4.121%
9	10.865	1358	1366	1384	rBV	88976	209701	11.41%	2.495%
10	13.303	1772	1781	1804	rBV	628819	1114146	60.60%	13.257%
11	14.684	2008	2016	2030	rBV2	193084	357638	19.45%	4.255%
12	16.176	2263	2270	2290	rBV	401327	722512	39.30%	8.597%
13	17.427	2475	2483	2499	rBV2	230584	439576	23.91%	5.230%
14	20.024	2918	2925	2938	rBV	1189485	1838501	100.00%	21.876%
15	21.693	3201	3209	3228	rBV	200466	466570	25.38%	5.552%
16	24.930	3743	3760	3779	rBV3	114189	430369	23.41%	5.121%

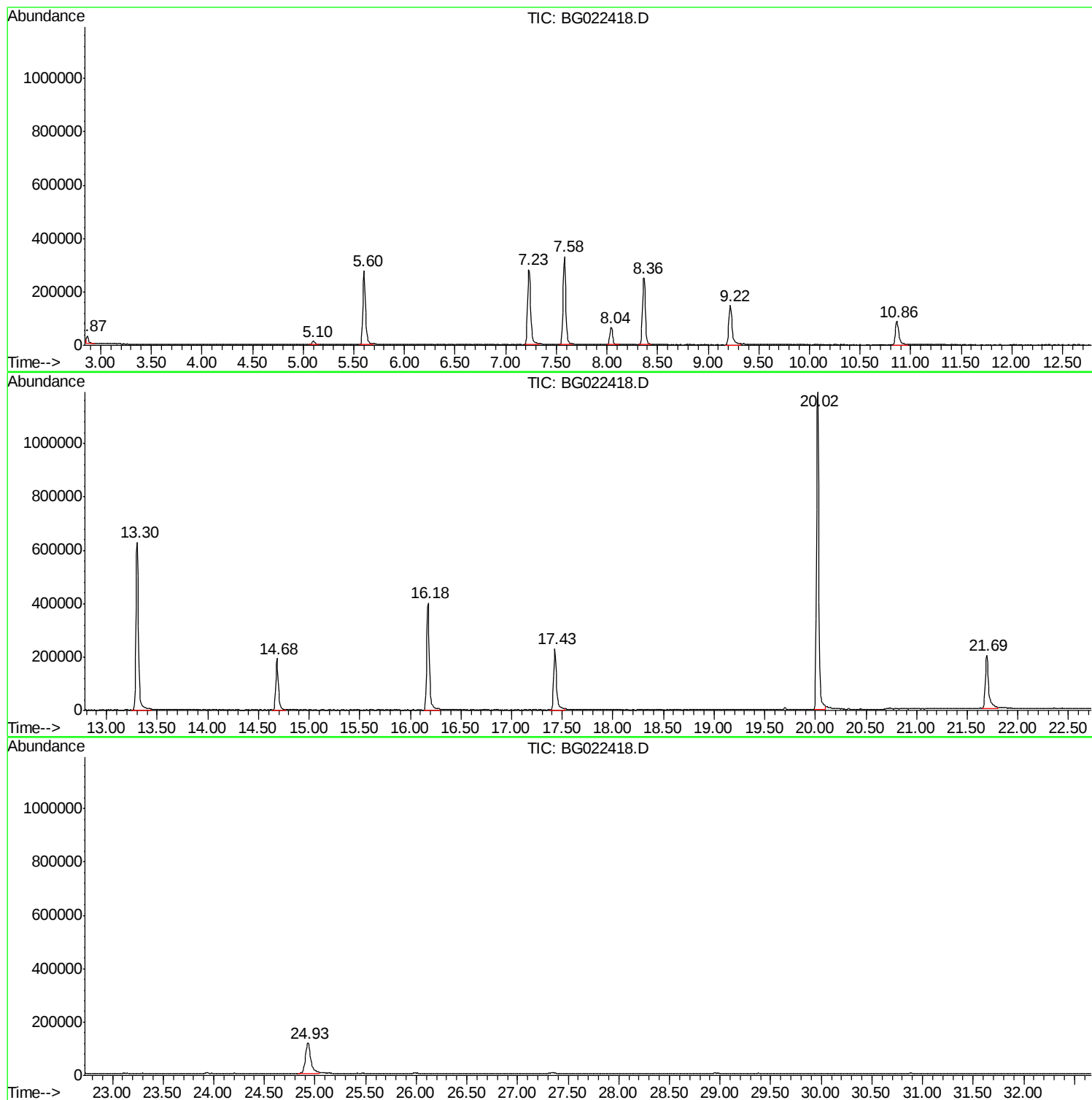
Sum of corrected areas: 8404160

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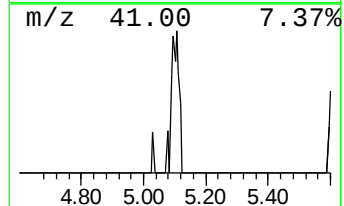
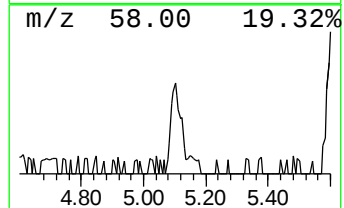
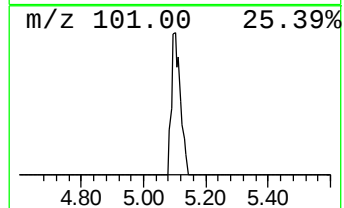
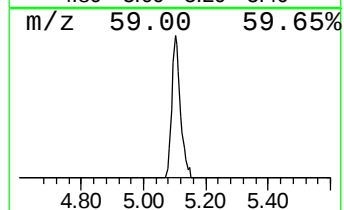
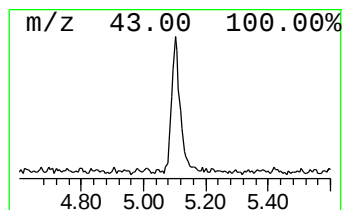
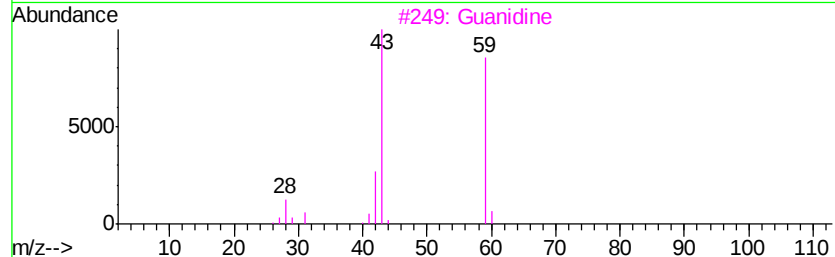
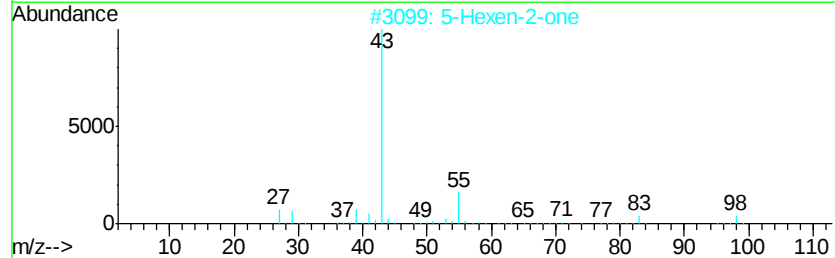
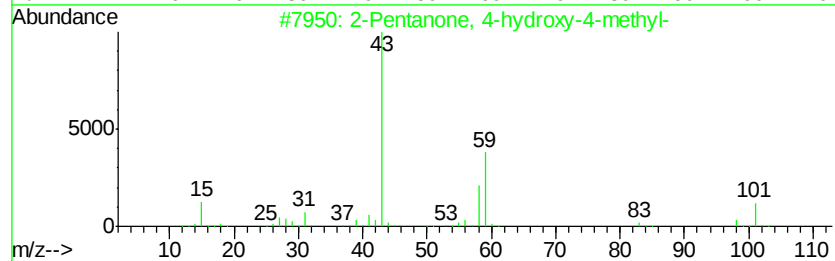
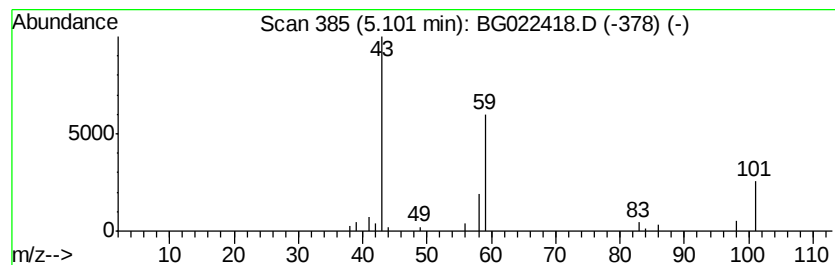
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.10	4.19 ng	26693	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	28
2		5-Hexen-2-one	98	C6H10O	000109-49-9	9
3		Guanidine	59	CH5N3	000113-00-8	9
4		2-Propanone, 1-cyclopropyl-	98	C6H10O	004160-75-2	9
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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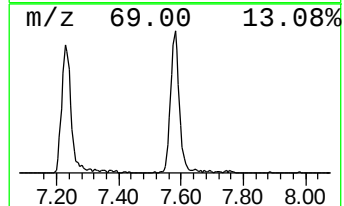
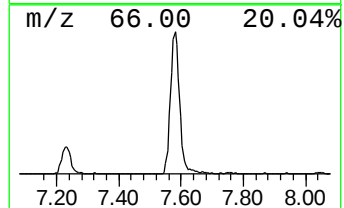
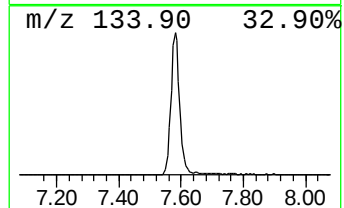
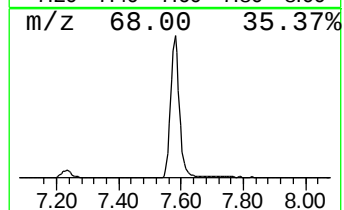
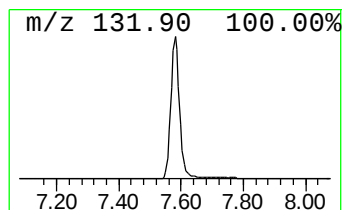
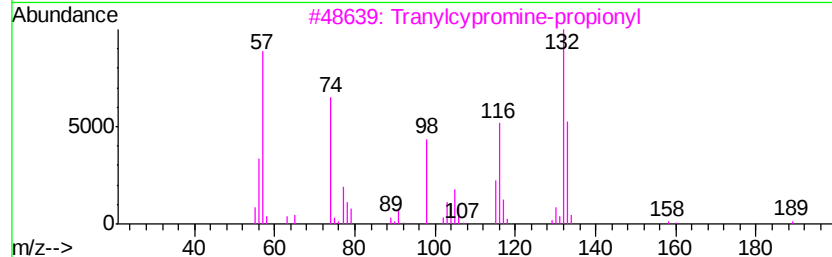
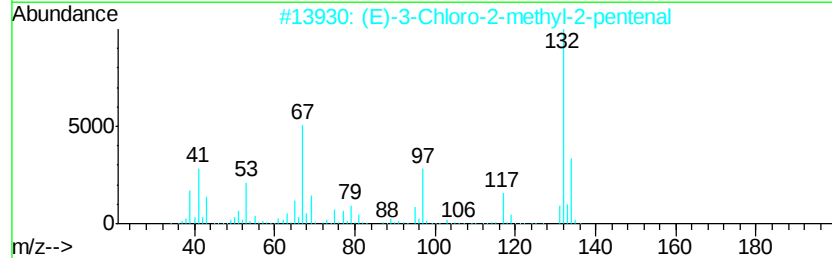
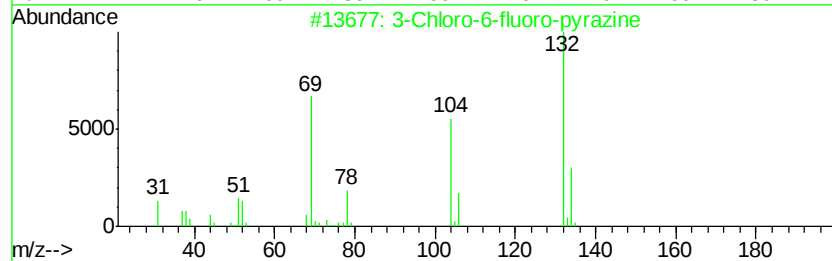
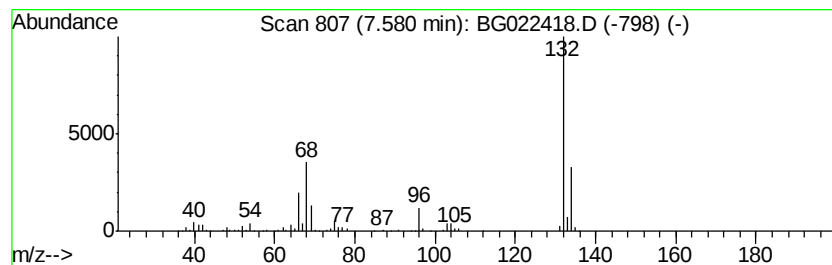
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Peak Number 3 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	100.58 ng	641091	1,4-Dichlorobenzene-d4	8.04

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		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		Amphetaminil	250	C17H18N2	017590-01-1	12
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.10	4.2	ng	26693	1	8.04	127476	20.0
unknown7.58	7.58	100.6	ng	641091	1	8.04	127476	20.0