

Data Path : Z:\HPCHEM1\BNA G\DATA\BG020317\
 Data File : BG025768.D
 Acq On : 2 Feb 2017 21:37
 Operator : SJ/MA
 Sample : PB96675BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB96675BL

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-BG013017.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	5.220	399	405	417	rBV	650764	1097133	13.02%	2.785%
2	5.702	477	487	502	rBV	1598846	2858717	33.93%	7.257%
3	7.324	755	763	778	rBV	1614650	3142461	37.30%	7.977%
4	7.694	818	826	840	rBV	1676036	2976670	35.33%	7.556%
5	8.164	899	906	916	rVB2	291287	530946	6.30%	1.348%
6	8.487	951	961	972	rBV	1238905	2198904	26.10%	5.582%
7	9.351	1100	1108	1123	rBV	1011395	1969534	23.38%	5.000%
8	10.996	1380	1388	1397	rBV	360600	766449	9.10%	1.946%
9	13.416	1792	1800	1809	rBV	2764268	4520839	53.66%	11.476%
10	14.797	2027	2035	2047	rBV	699088	1192702	14.16%	3.028%
11	16.284	2280	2288	2304	rBV	2435590	4208689	49.96%	10.684%
12	17.541	2495	2502	2513	rBV2	957031	1584972	18.81%	4.023%
13	20.126	2935	2942	2951	rBV	6189697	8424957	100.00%	21.386%
14	21.830	3225	3232	3242	rBV	1107376	2008176	23.84%	5.098%
15	25.209	3796	3807	3818	rBV2	606511	1913118	22.71%	4.856%

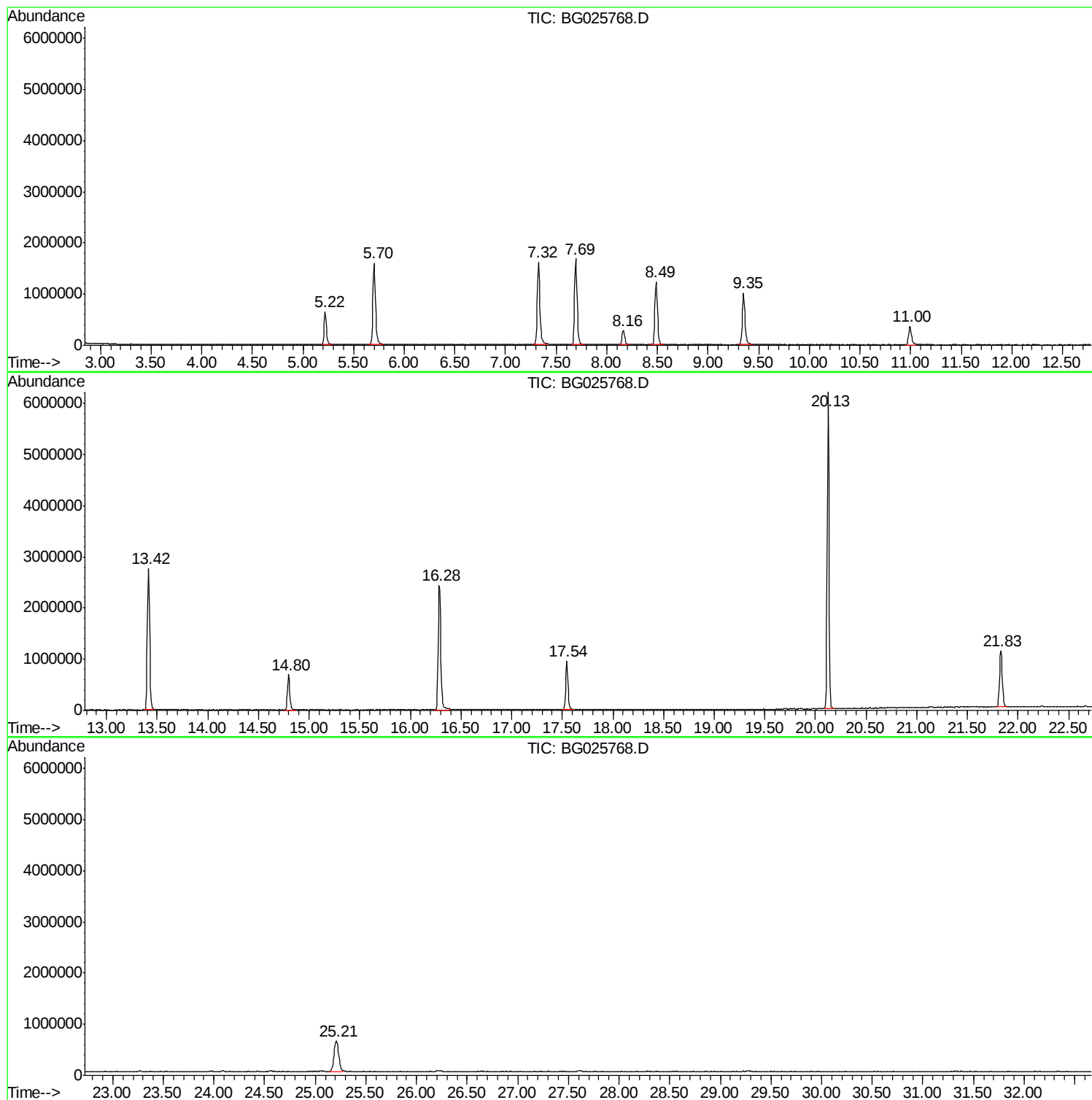
Sum of corrected areas: 39394267

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG013017.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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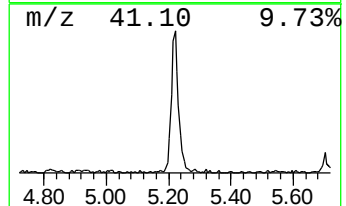
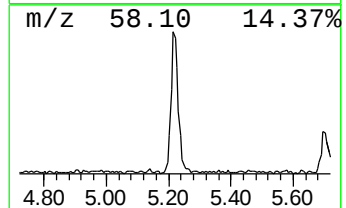
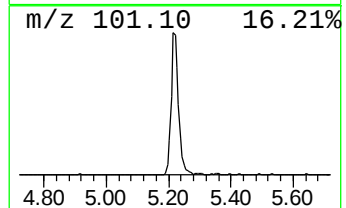
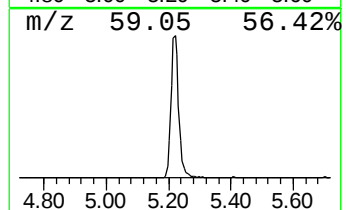
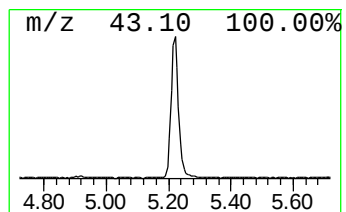
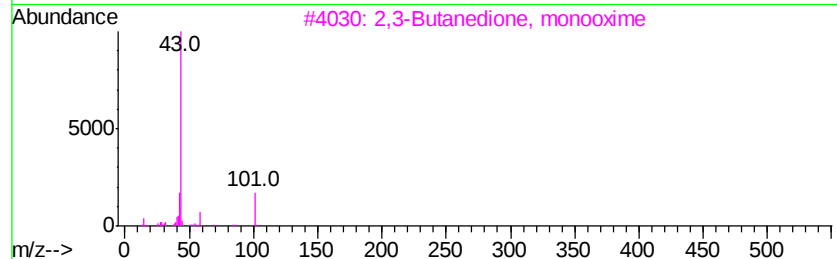
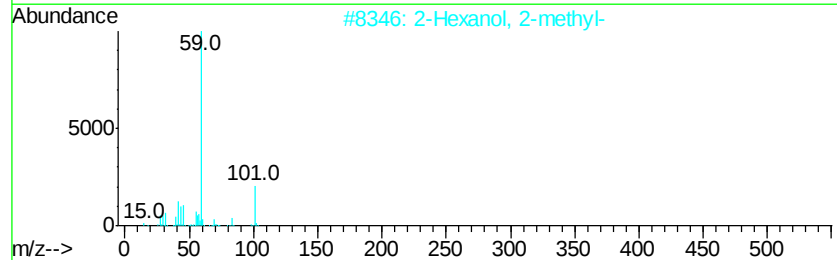
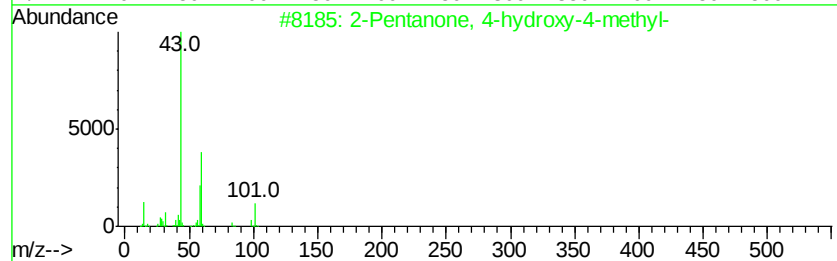
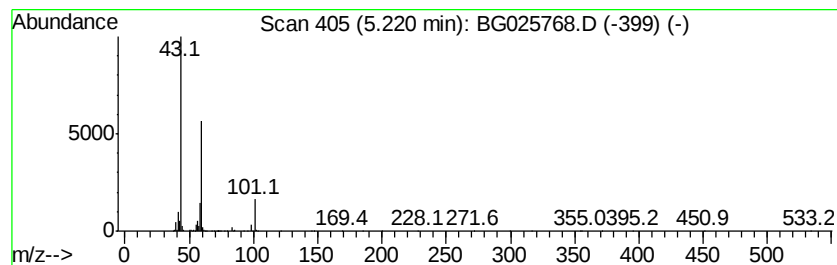
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.22	41.33 ng	1097130	1,4-Dichlorobenzene-d4	8.16

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	25
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
4			3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	9
5			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9



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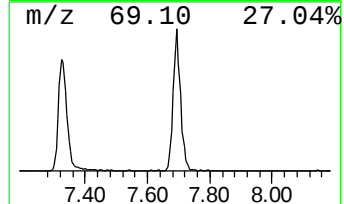
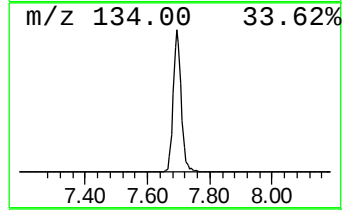
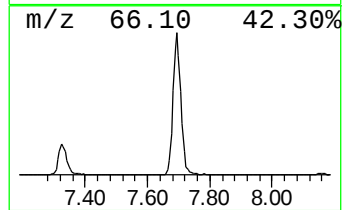
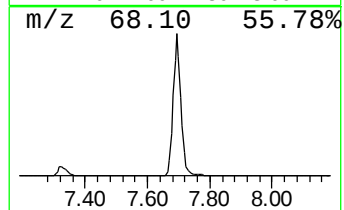
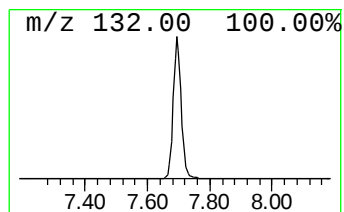
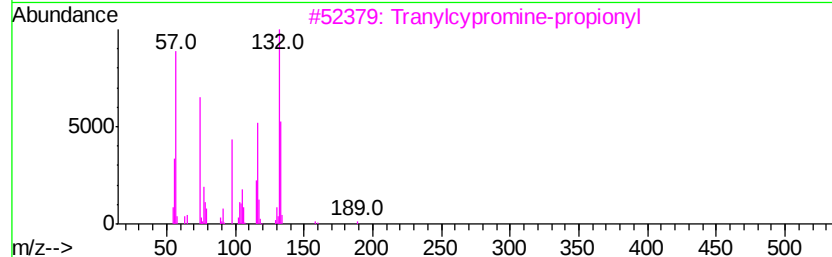
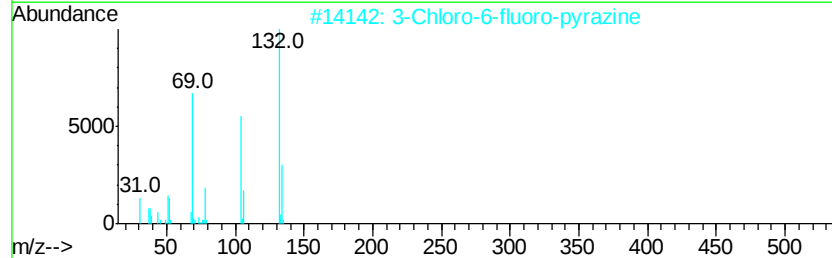
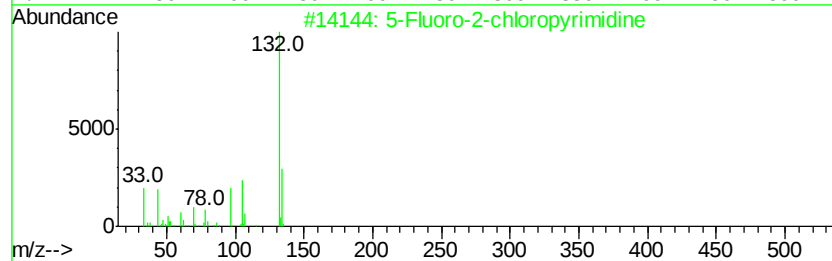
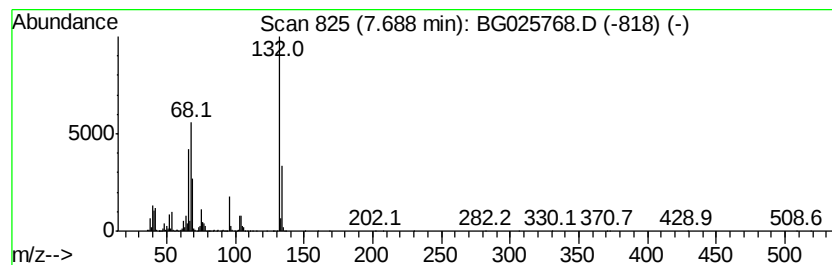
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Peak Number 2 unknown7.69 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.69	112.13 ng	2976670	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	14
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
5		1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	10



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.22	41.3	ng	1097130	1	8.16	530946	20.0
unknown7.69	7.69	112.1	ng	2976670	1	8.16	530946	20.0