

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060517\
 Data File : BG027225.D
 Acq On : 5 Jun 2017 23:11
 Operator : SJ/MA
 Sample : PB99436BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB99436BL

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-BG060517.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	5.715	338	345	357	rBV	88955	153996	3.08%	0.654%
2	6.150	411	419	430	rBV	1057147	1842111	36.80%	7.822%
3	7.701	674	683	697	rBV	1086904	1984931	39.66%	8.428%
4	8.101	743	751	761	rBV	1028023	1904799	38.06%	8.088%
5	8.571	822	831	840	rVB	168479	306878	6.13%	1.303%
6	8.900	877	887	897	rVB	689257	1321536	26.40%	5.611%
7	9.734	1019	1029	1039	rBV	559683	1044615	20.87%	4.436%
8	11.391	1303	1311	1321	rVB	257652	493906	9.87%	2.097%
9	13.782	1710	1718	1725	rBV	1838371	3065526	61.25%	13.016%
10	15.151	1944	1951	1959	rBV2	436222	739954	14.78%	3.142%
11	16.632	2196	2203	2218	rBV	1488108	2456198	49.07%	10.429%
12	17.901	2413	2419	2427	rBV2	599860	987273	19.73%	4.192%
13	20.486	2852	2859	2865	rBV	3602107	5005117	100.00%	21.252%
14	22.296	3160	3167	3178	rVB	605182	1159082	23.16%	4.922%
15	26.003	3789	3798	3813	rVB2	342467	1085268	21.68%	4.608%

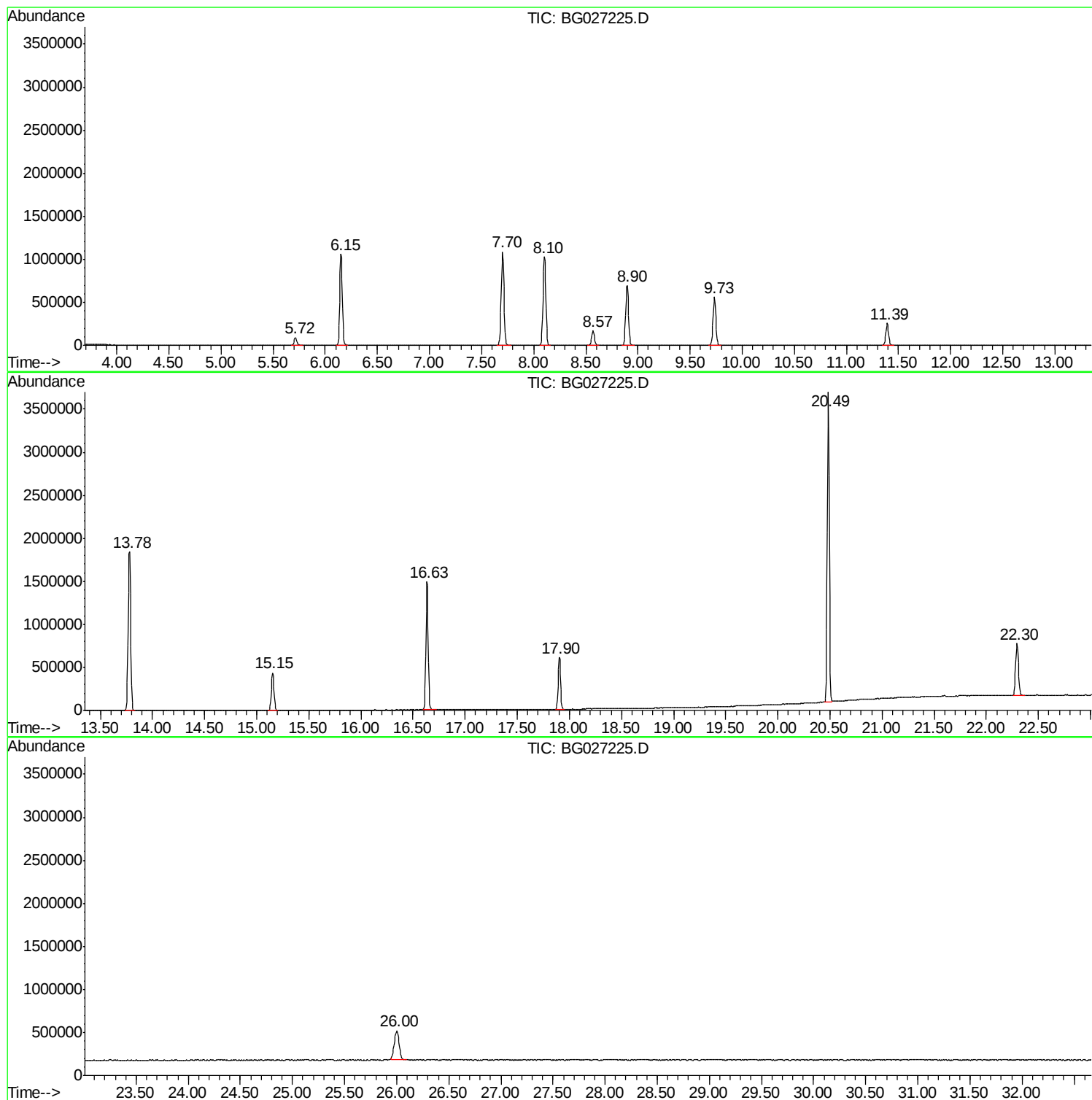
Sum of corrected areas: 23551190

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060517.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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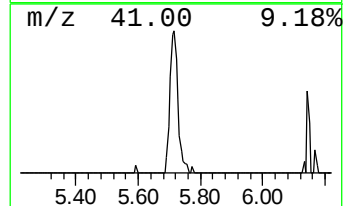
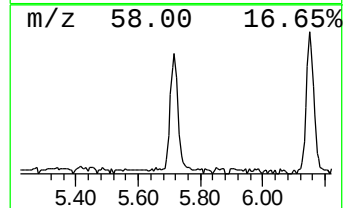
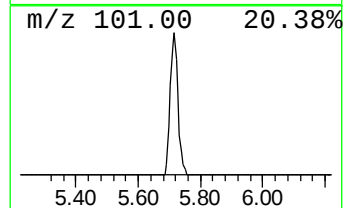
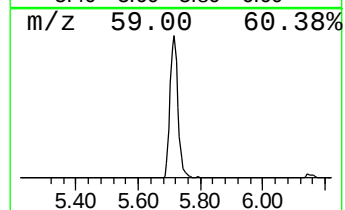
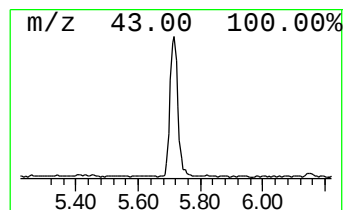
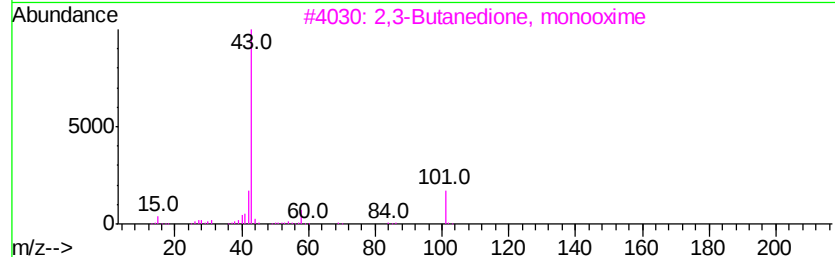
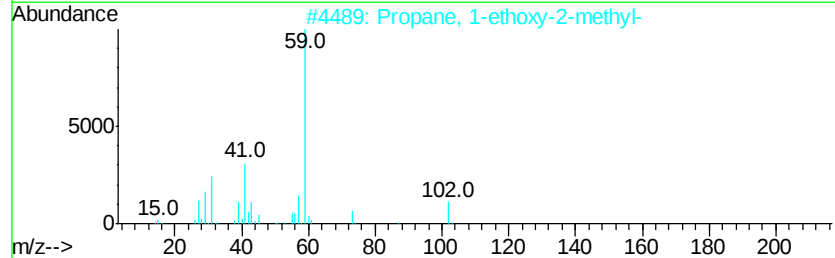
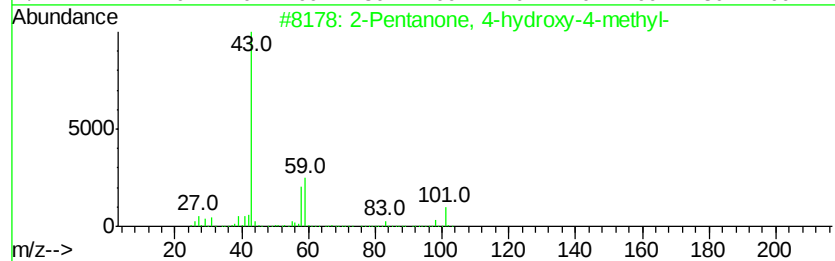
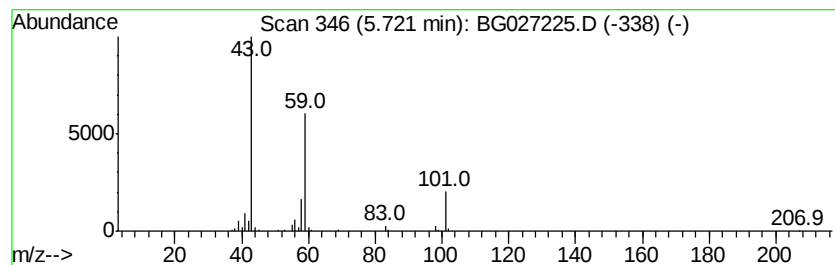
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.72	10.04 ng	153996	1,4-Dichlorobenzene-d4	8.57

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	38
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
5			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	17



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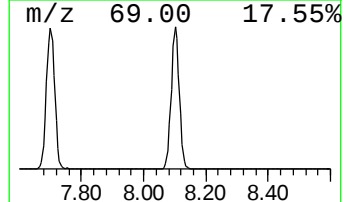
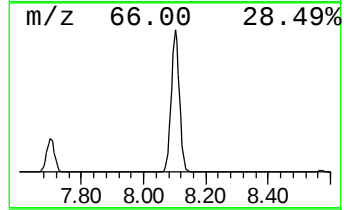
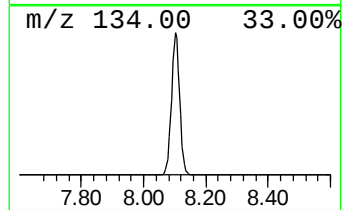
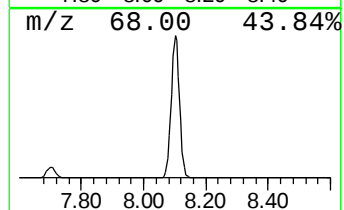
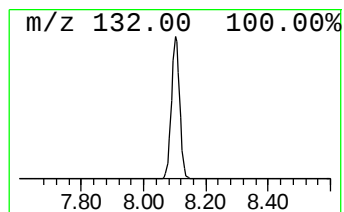
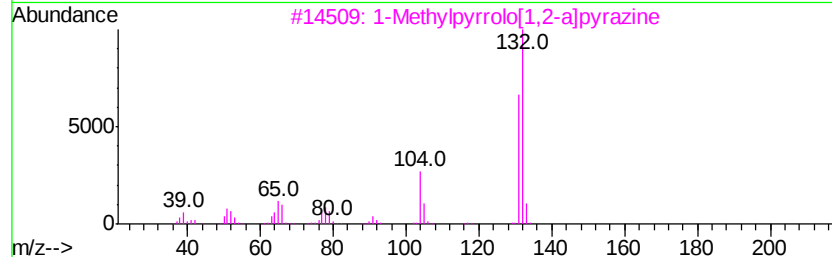
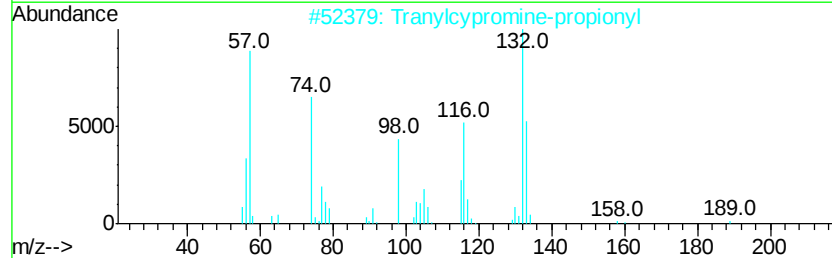
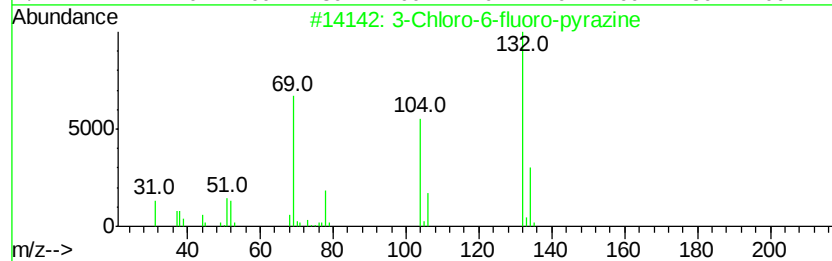
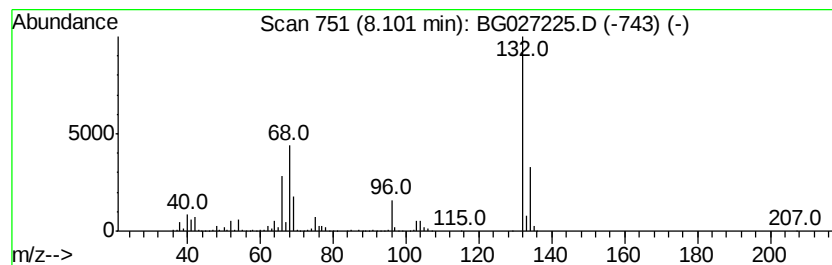
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Peak Number 2 unknown8.10 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.10	124.14 ng	1904800	1,4-Dichlorobenzene-d4	8.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
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...	5.72	10.0	ng	153996	1	8.57	306878	20.0
unknown8.10	8.10	124.1	ng	1904800	1	8.57	306878	20.0