

Data Path : Z:\HPCHEM1\BNA G\DATA\BG050115\
 Data File : BG016851.D
 Acq On : 1 May 2015 18:48
 Operator : TP/IZ
 Sample : PB83121BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB83121BL

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-BG042915.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	4.380	283	288	297	rBV	64882	88145	3.52%	0.650%
2	4.867	365	371	390	rBV	756849	1131017	45.20%	8.337%
3	6.483	640	646	670	rBV	688306	1200792	47.99%	8.852%
4	6.800	693	700	719	rBV	773382	1276639	51.02%	9.411%
5	7.259	773	778	797	rBV	137565	270218	10.80%	1.992%
6	7.576	825	832	859	rBV	585989	984455	39.34%	7.257%
7	8.422	969	976	1004	rBV	360787	690829	27.61%	5.092%
8	10.032	1243	1250	1266	rBV	197095	373779	14.94%	2.755%
9	12.552	1672	1679	1696	rBV	1065937	1652973	66.06%	12.185%
10	13.945	1910	1916	1933	rBV2	301508	489792	19.57%	3.610%
11	15.449	2166	2172	2188	rBV	656383	1028221	41.09%	7.579%
12	16.695	2378	2384	2398	rBV2	340805	565544	22.60%	4.169%
13	19.356	2831	2837	2850	rBV	2091560	2502303	100.00%	18.446%
14	20.907	3096	3101	3110	rBV	529589	658486	26.32%	4.854%
15	22.981	3447	3454	3466	rBV	357684	652678	26.08%	4.811%

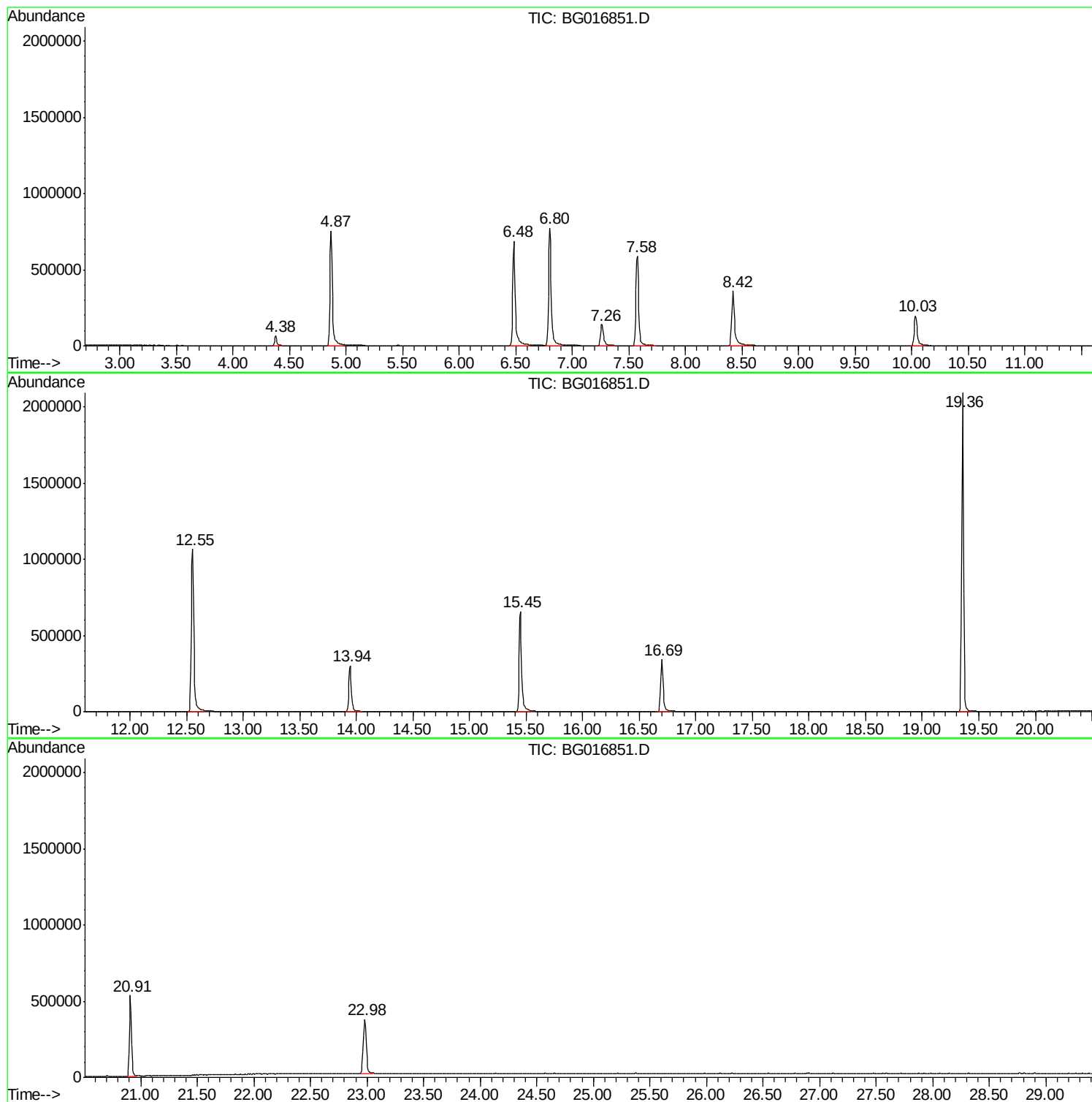
Sum of corrected areas: 13565871

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042915.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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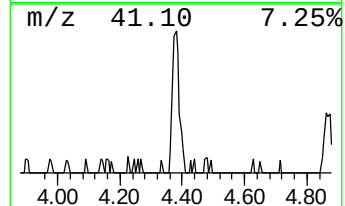
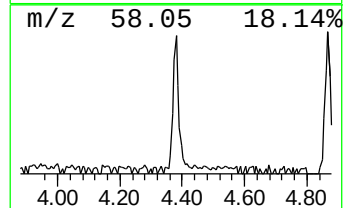
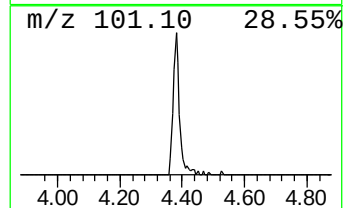
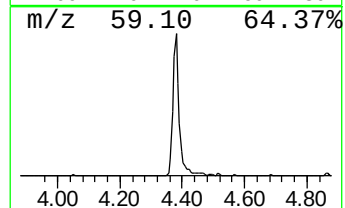
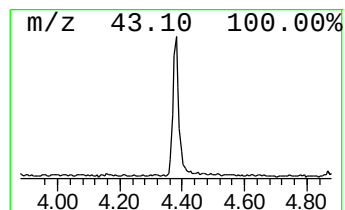
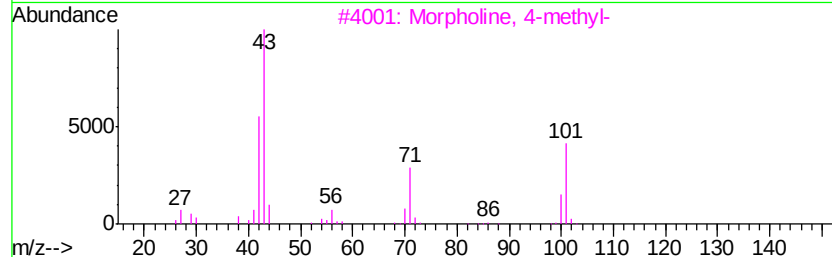
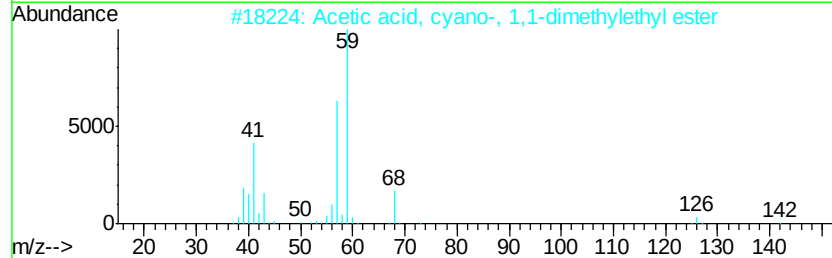
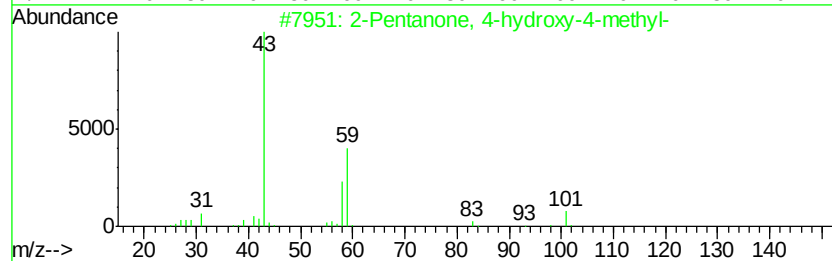
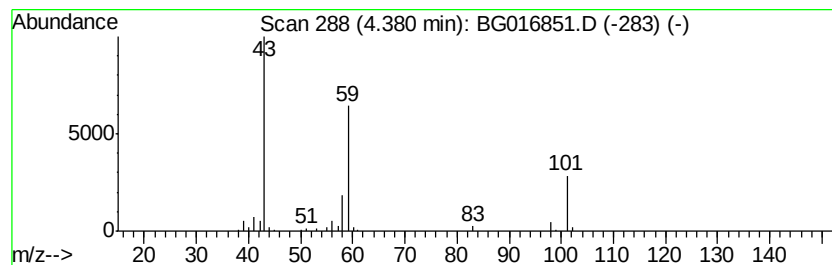
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.38	6.52 ng	88145	1,4-Dichlorobenzene-d4	7.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4	Guanidine	59	CH5N3	000113-00-8	9
5	Propanamide, N-ethyl-	101	C5H11NO	005129-72-6	9



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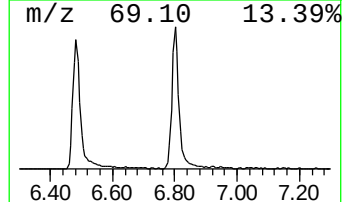
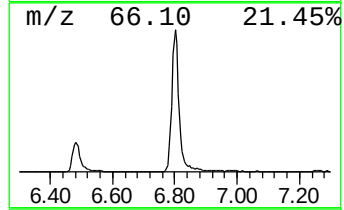
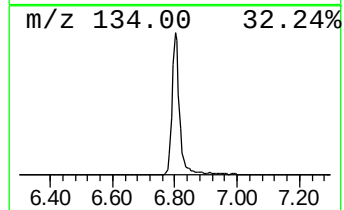
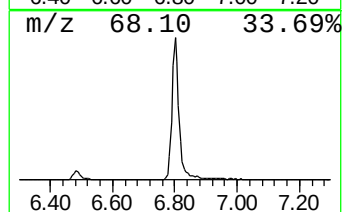
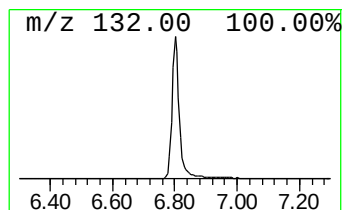
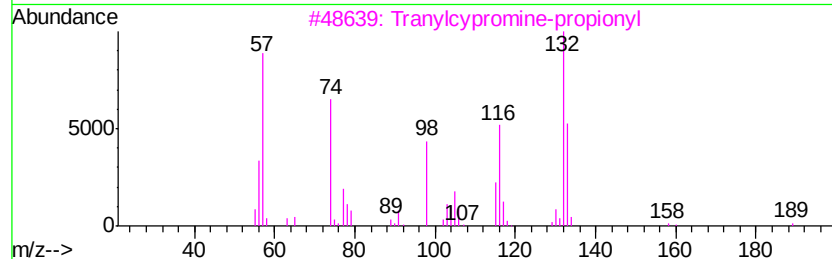
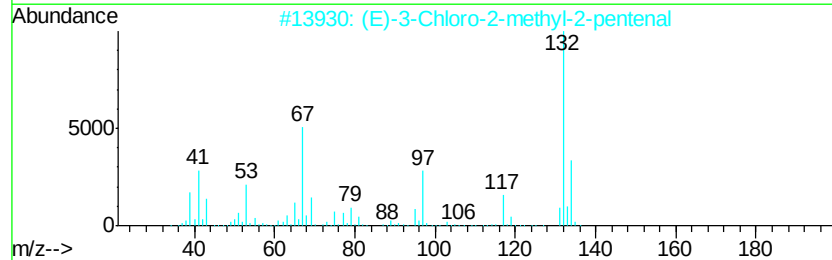
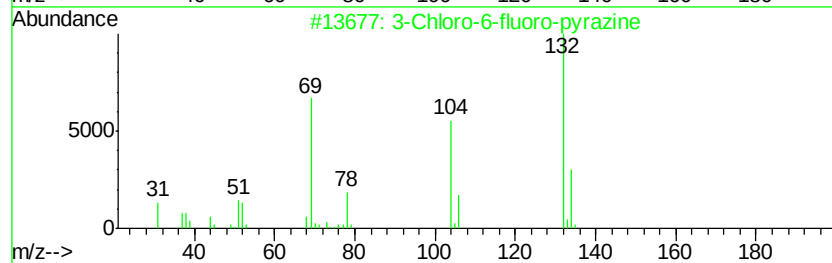
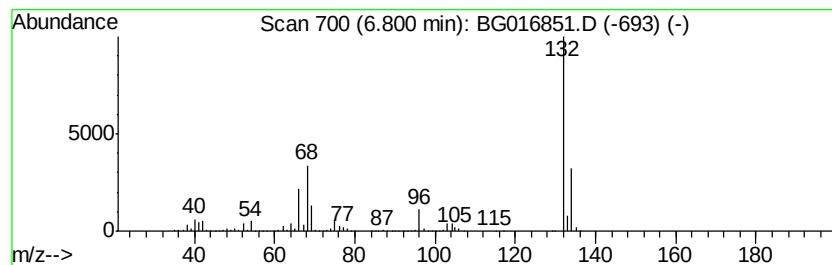
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Peak Number 2 unknown6.80 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.80	94.49 ng	1276640	1,4-Dichlorobenzene-d4	7.26

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	35
3		Tranvlcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.38	6.5	ng	88145	1	7.26	270218	20.0
unknown6.80	6.80	94.5	ng	1276640	1	7.26	270218	20.0