

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060315\
 Data File : BG017431.D
 Acq On : 4 Jun 2015 6:52
 Operator : TP/IZ
 Sample : PB83703BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB83703BL

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-BG060315.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	4.729	343	347	354	rVB2	18316	24185	1.49%	0.340%
2	5.217	424	430	439	rBV	360282	499075	30.66%	7.006%
3	6.827	697	704	716	rBV	357526	538911	33.11%	7.565%
4	7.168	754	762	771	rBV	351385	544283	33.44%	7.641%
5	7.626	833	840	846	rVB	77743	116376	7.15%	1.634%
6	7.943	888	894	901	rBV	255754	388384	23.86%	5.452%
7	8.789	1029	1038	1047	rBV	206770	334565	20.55%	4.697%
8	10.423	1308	1316	1324	rBV	93251	161445	9.92%	2.266%
9	12.908	1732	1739	1750	rBV	510136	738686	45.38%	10.370%
10	14.289	1968	1974	1982	rBV2	155401	232275	14.27%	3.261%
11	15.787	2223	2229	2241	rBV	508816	707935	43.49%	9.938%
12	17.038	2436	2442	2448	rBV2	233956	321467	19.75%	4.513%
13	19.377	2835	2840	2844	rBV2	22642	26518	1.63%	0.372%
14	19.676	2885	2891	2900	rVB	1341567	1627670	100.00%	22.850%
15	20.423	3013	3018	3026	rBV5	17902	27585	1.69%	0.387%
16	21.233	3150	3156	3160	rBV	347422	434366	26.69%	6.098%
17	23.466	3529	3536	3542	rBV2	224245	399688	24.56%	5.611%

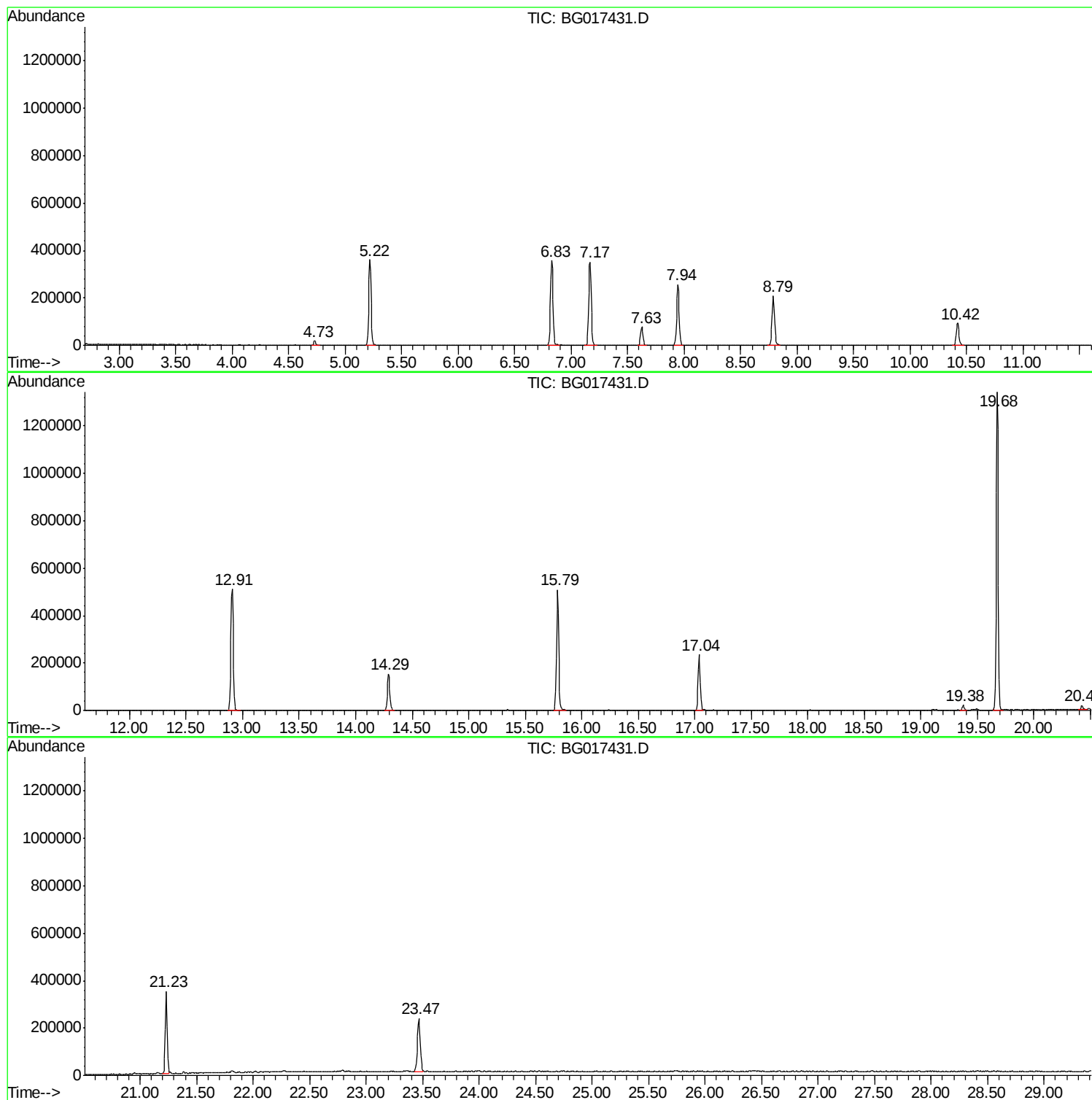
Sum of corrected areas: 7123414

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Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060315.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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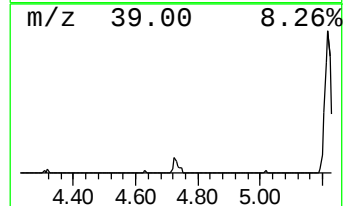
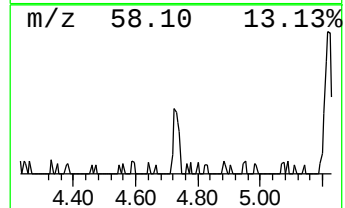
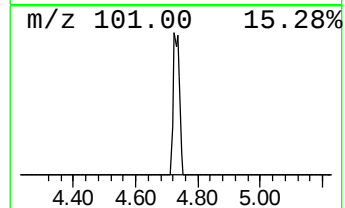
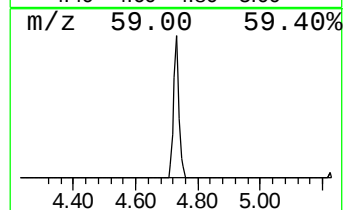
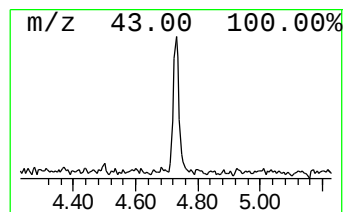
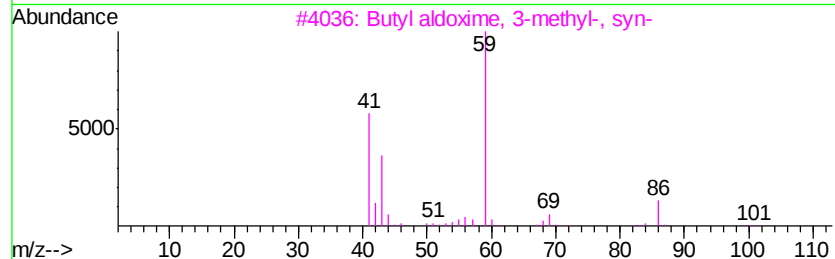
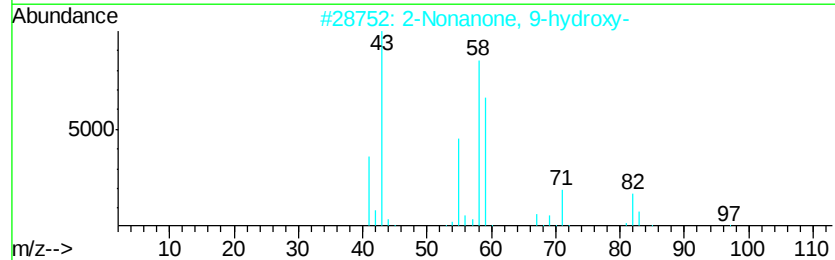
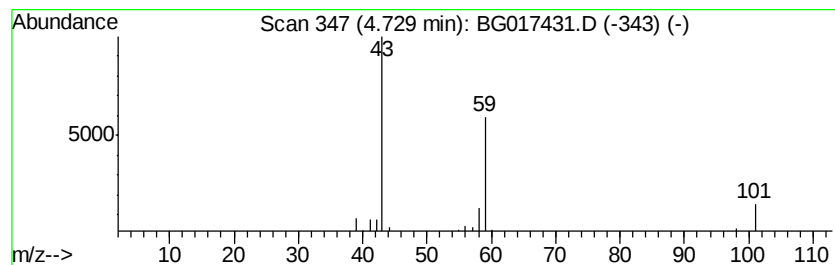
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.73	4.16 ng	24185	1,4-Dichlorobenzene-d4	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2	2-Nonanone, 9-hydroxy-	158	C9H18O2	025368-56-3	38
3	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	37
4	Butane, 1-(1-methylethoxy)-	116	C7H16O	001860-27-1	32
5	Diacetamide	101	C4H7NO2	000625-77-4	9



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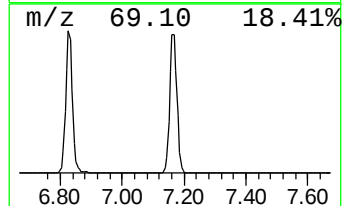
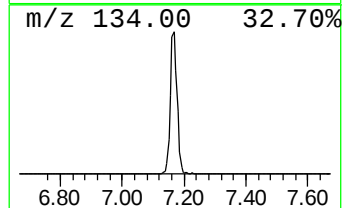
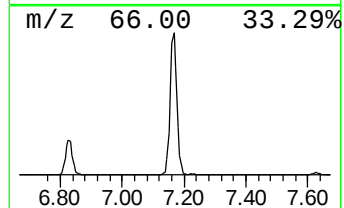
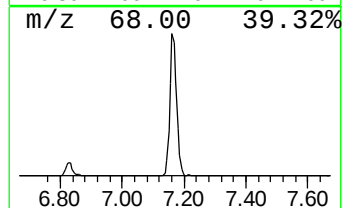
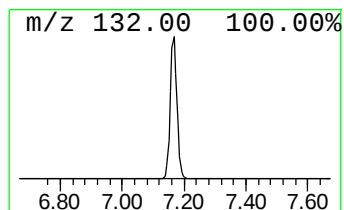
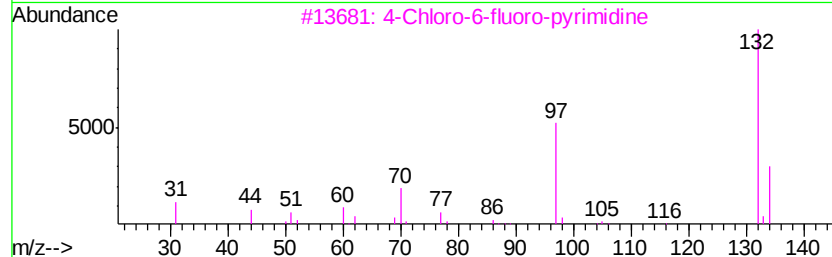
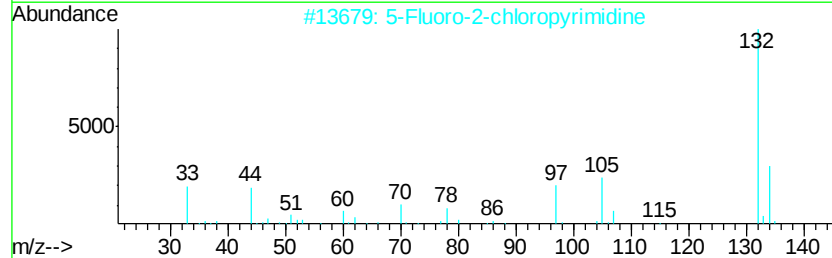
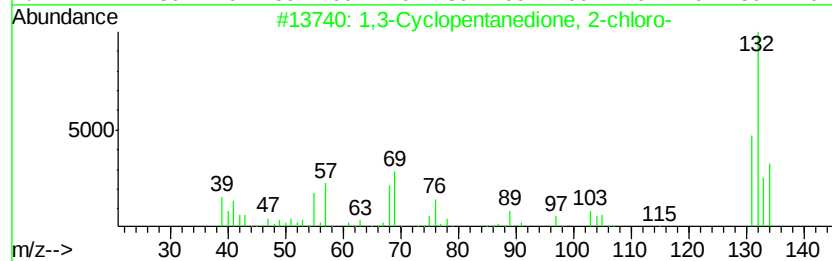
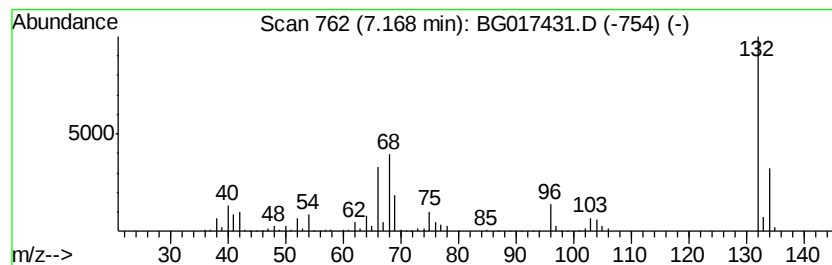
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Peak Number 2 unknown7.17 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.17	93.54 ng	544283	1,4-Dichlorobenzene-d4	7.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	38
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
3		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.73	4.2	ng	24185	1	7.63	116376	20.0
unknown7.17	7.17	93.5	ng	544283	1	7.63	116376	20.0