

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG062217\
 Data File : BG027626.D
 Acq On : 23 Jun 2017 11:14
 Operator : SJ/JU
 Sample : PB100013BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB100013BL

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-BG062117.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	5.639	292	298	308	rVB	46449	74115	3.36%	0.873%
2	6.085	367	374	386	rBV	341380	579218	26.25%	6.825%
3	7.642	631	639	649	rBV	335767	604151	27.38%	7.119%
4	8.036	698	706	718	rBV	335315	593562	26.90%	6.994%
5	8.500	779	785	793	rBB	57848	102767	4.66%	1.211%
6	8.829	834	841	850	rBV	229377	421511	19.10%	4.967%
7	9.669	976	984	995	rBV	185357	352337	15.97%	4.152%
8	11.320	1256	1265	1275	rBV	80399	151882	6.88%	1.790%
9	13.712	1665	1672	1680	rBV	584199	907925	41.14%	10.698%
10	15.086	1900	1906	1916	rBB2	131815	225046	10.20%	2.652%
11	16.573	2151	2159	2171	rBV	613580	983365	44.56%	11.587%
12	17.836	2367	2374	2383	rBV2	236470	369077	16.73%	4.349%
13	20.415	2806	2813	2823	rBV	1645751	2206700	100.00%	26.002%
14	22.207	3110	3118	3125	rBV	250287	472123	21.39%	5.563%
15	25.838	3722	3736	3749	rBV3	132393	442716	20.06%	5.217%

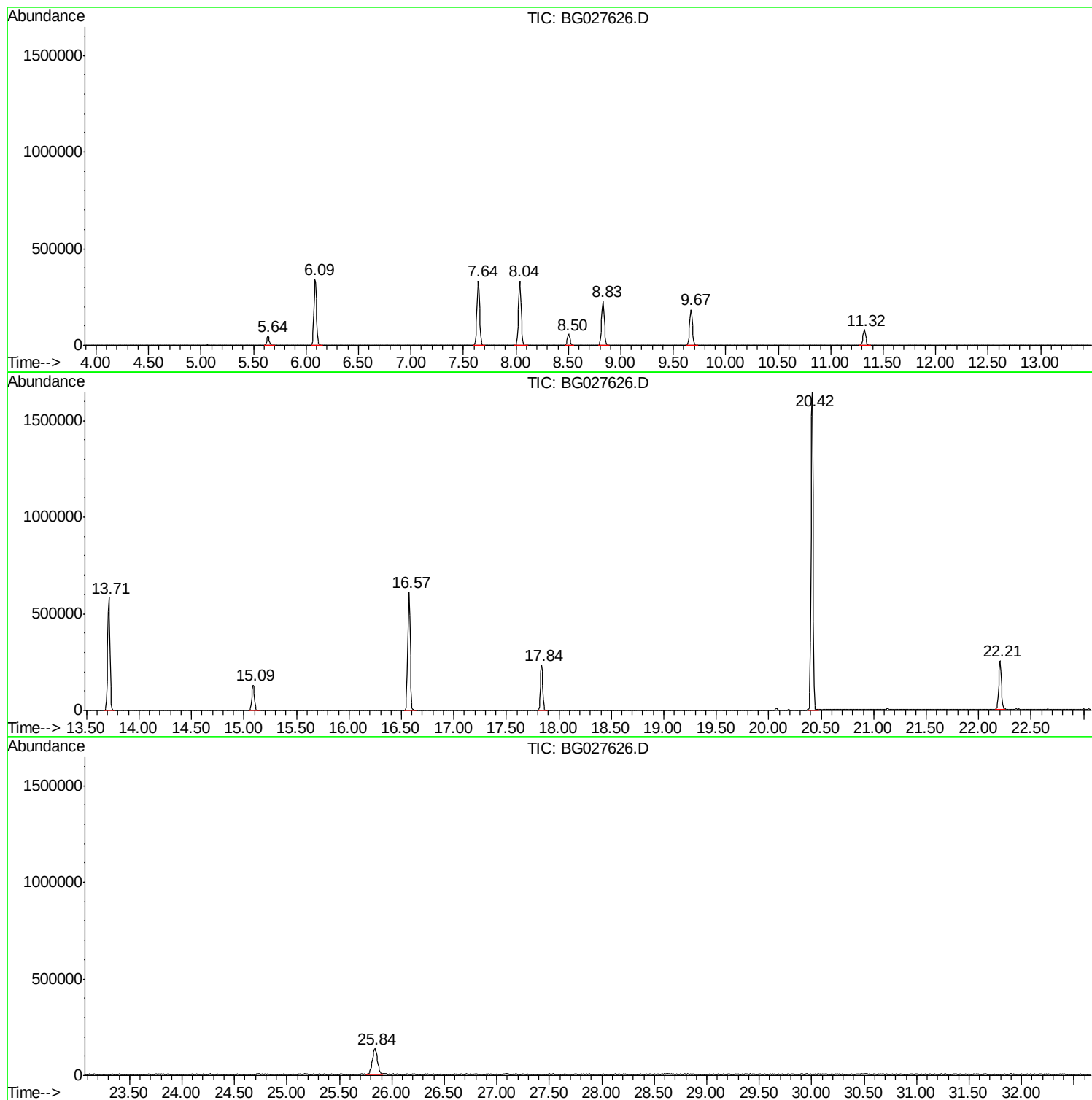
Sum of corrected areas: 8486495

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG062117.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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Instrument :
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ClientSampled :
PB100013BL

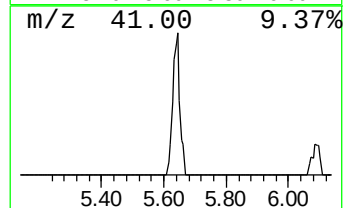
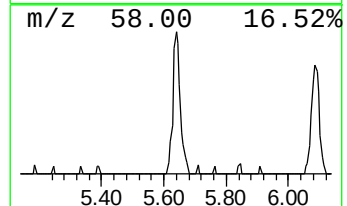
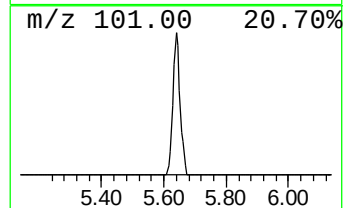
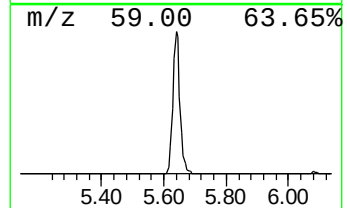
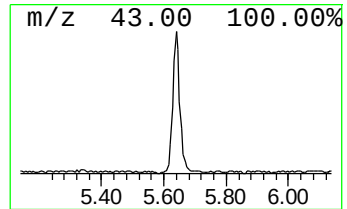
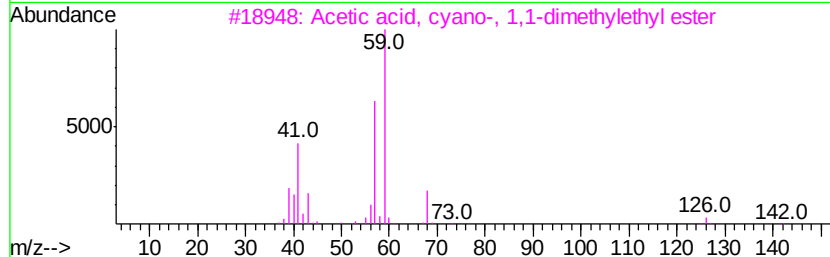
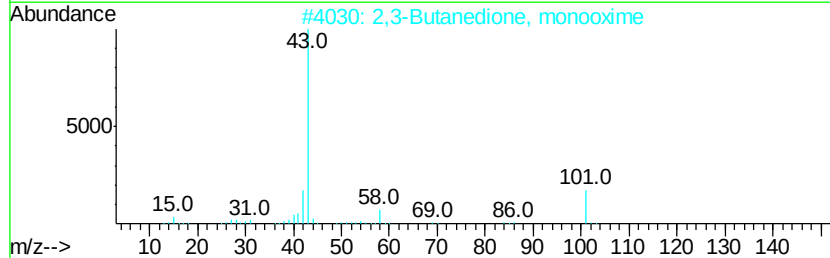
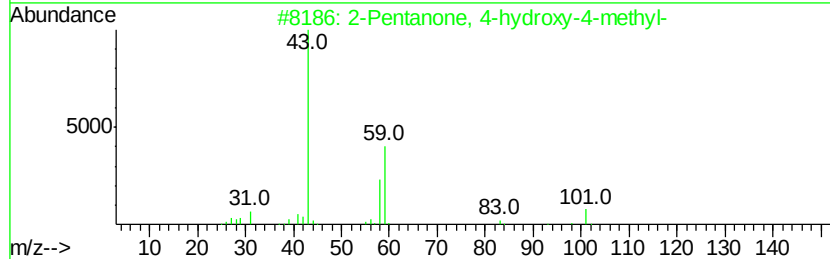
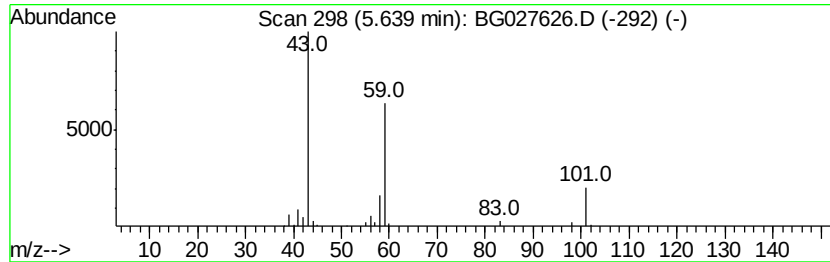
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.64	14.42 ng	74115	1,4-Dichlorobenzene-d4	8.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	32
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4		Acetone	58	C3H6O	000067-64-1	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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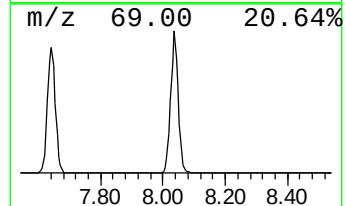
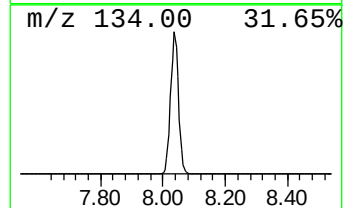
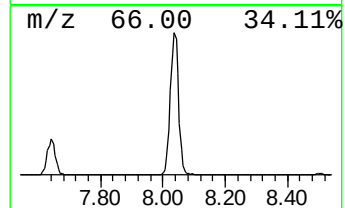
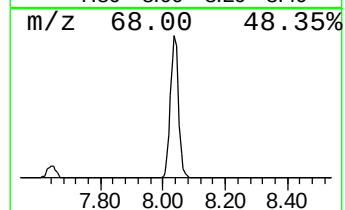
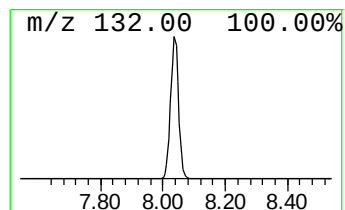
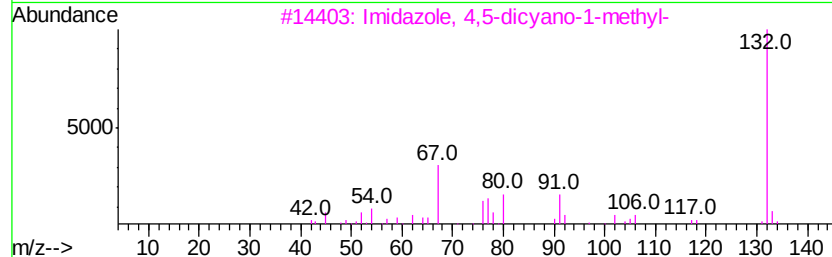
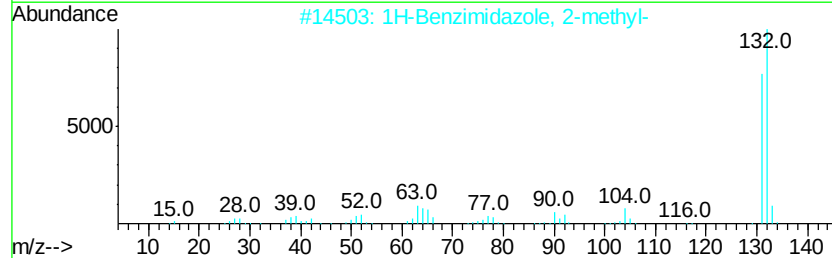
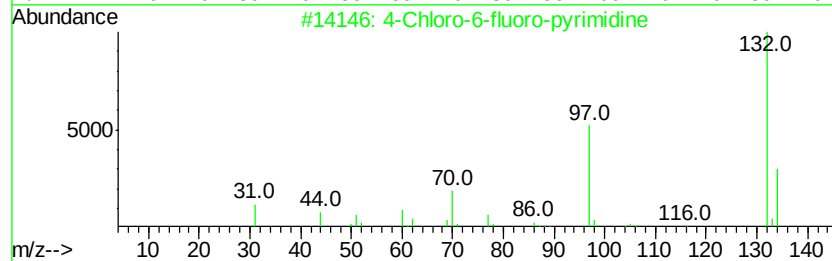
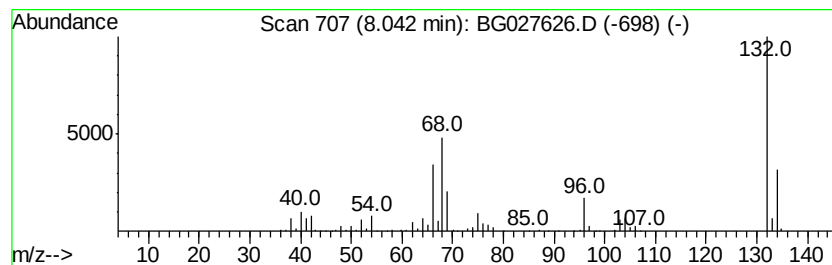
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Peak Number 2 unknown8.04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.04	115.52 ng	593562	1,4-Dichlorobenzene-d4	8.50

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	22
4		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
5		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.64	14.4	ng	74115	1	8.50	102767	20.0
unknown8.04	8.04	115.5	ng	593562	1	8.50	102767	20.0