

# BeadDataPackR

October 5, 2010

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`compressBeadData`     *Write raw bead level data to a compressed format.*

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## Description

Given raw bead level data, in the form of .txt and .locs file, this function combines the two producing a new file with the data stored in a compressed format.

## Usage

```
compressBeadData(txtFile, locsGrn, locsRed = NULL, outputFile = NULL,  
                 path = NULL, nBytes = 8, base2 = TRUE, fullLocsIndex = FALSE,  
                 nrow = NULL, ncol = NULL, progressBar = TRUE)
```

## Arguments

<code>txtFile</code>	The name of the .txt file to be read in.
<code>locsGrn</code>	The locs file for the green channel.
<code>locsRed</code>	The locs file for the red channel. Only needed for two channel data.
<code>outputFile</code>	Name of the file to be created.
<code>path</code>	Path to where the input files can be found. If NULL the current working directory is used. This is also the directory where the output files will be written.
<code>nBytes</code>	Gives the number of bytes that are used to store the fractional parts of the bead coordinates. For a single channel array the maximum value is 4, whilst it is 8 for a two channel array. Any number larger than this is automatically set the maximum value. If the maximum value is used the coordinates are stored in the .bab file as single precision floating point numbers, as they are in the .locs files. If a value smaller than the maximum is chosen then the integer parts of each coordinate are stored separately. The requested number of bytes are then used to store the fractional parts, with a corresponding loss of precision as the number of bytes decreases.
<code>base2</code>	If not using the full precision coordinates, the approximations can be stored as either a binary or decimal fraction. Using a binary fraction ( <code>base2 = TRUE</code> ) provides a greater accuracy, but can lead to a meandering number of decimal places in the reconstructed .txt files. If one wants a consistent number of decimal places, set <code>base2 = FALSE</code> .

fullLocsIndex	Default value of 0 uses a linear model fitted to each segment of the array to allow reconstruct the locs file when the file is decompressed. Using a value of 1 a simple index is used to record the locs file order, but requires more space.
ncol	This specifies the number of columns in each grid segment on the array and, if left blank, can normally be inferred from the grid coordinates. However, this can fail for particularly small grids. If one wants or needs to specify them explicitly, these values can be found in the .sdf which accompanies the bead level output from the scanner. The number of columns per segment can be found within the tag <SizeGridX>
nrow	See ncol. If needed can be found within the <SizeGridY> tag in the .sdf file.
progressBar	By default the function uses a <code>txtProgressBar</code> to indicate progress through the compression. Setting this argument to FALSE suppresses the drawing of this progress bar.

### Details

In the future the file names will be determined automatically, rather than requiring manual entry of each. The path argument may also be amended so there are separate options for the locations of the input and output files.

### Value

Primarily invoked for its side effect, which is to produce a compressed version of the input files. The function returns, invisibly, the total number of beads stored in the compressed file.

### Author(s)

Mike L. Smith

### Examples

```
dataPath <- system.file("data", package = "BeadDataPackR")
compressBeadData(txtFile = "example.txt", locsGrn = "example_Grn.locs", outputFile =
```

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`decompressBeadData` *Decompress a file in the beadarray binary format*

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### Description

Decompresses a file created by `BeadDataPackR`. The original files that were compressed will be restored as accurately as possible, depending upon the degree of precision specified during the compression.

### Usage

```
decompressBeadData(inputFile, inputPath = ".", outputMask = NULL, outputPath = "
outputNonDecoded = FALSE, roundValues = TRUE, progressBar = T
```

**Arguments**

<code>inputFile</code>	The name of the file to be read.
<code>inputPath</code>	Path where the compress file is located. The default is to use the current working directory.
<code>outputMask</code>	Text specify the names of the output files. The output files will have ".txt", "_Grn.locs" and (if appropriate "_Red.locs") appended to this mask. If left NULL the original names of the section will be used.
<code>outputPath</code>	Path to where the uncompressed version of the files should be written to. The default is to use the current working directory.
<code>outputNonDecoded</code>	If TRUE the undecoded beads will be included in the output .txt file. They will have ProbeID 0 and intensity 0, but the bead centre coordinates will be included.
<code>roundValues</code>	The original Illumina text files give the bead centre coordinates to 7 significant figures. When this argument is TRUE decompressed files are also truncated in this manner, whilst FALSE writes them to the full precision they are stored in the compressed file.
<code>progressBar</code>	By default the function uses a <code>txtProgressBar</code> to indicate progress through the compression. Setting this argument to FALSE suppresses the drawing of this progress bar.

**Value**

Called primarily for its side effect, in which two (or three) files are written to the disk. These files should be representative of the original files that were compressed. The function returns, invisibly, the number of lines written in the .txt file.

**Author(s)**

Mike L. Smith

**Examples**

```
dataPath <- system.file("data", package = "BeadDataPackR")
decompressBeadData(inputFile = "example.bab", inputPath = dataPath, outputPath = ".")
```

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