Package ‘biganalytics’

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Title Utilities for 'big.matrix' Objects from Package 'bigmemory'

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Depends stats, utils, bigmemory (>= 4.0.0), foreach, biglm

Suggests methods

LinkingTo Rcpp, bigmemory, BH

Description Extend the 'bigmemory' package with various analytics.
   Functions 'bigkmeans' and 'binit' may also be used with native R objects.
   For 'tapply'-like functions, the bigtabulate package may also be helpful.
   For linear algebra support, see 'bigalgebra'. For mutex (locking) support
   for advanced shared-memory usage, see 'synchronicity'.

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biganalytics-package

Utilities for big.matrix objects of package bigmemory

Description

Extend the bigmemory package with various analytics. In addition to the more obvious summary statistics (see colmean, etc...), biganalytics offers biglm.big.matrix, bigglm.big.matrix, bigkmeans, binit, and apply for big.matrix objects. Some of the functions may be used with native R objects, as well, providing gains in speed and memory-efficiency.

apply.big.matrix-method

apply() for big.matrix objects

Description

apply for big.matrix objects. Note that the performance may be degraded (compared to apply with regular R matrices) because of S4 overhead associated with extracting data from big.matrix objects. This sort of limitation is unavoidable and would be the case (or even worse) with other "custom" data structures. Of course, this would only be partially significant if you are applying over lengthy rows or columns.

Usage

## S4 method for signature 'big.matrix'
apply(X, MARGIN, FUN, ..., simplify)

Arguments

- X: a big.matrix object.
- MARGIN: the margin. May be may only be 1 or 2, but otherwise conforming to what you would expect from apply().
- FUN: the function to apply.
- ...: other parameters to pass to the FUN parameter.
- simplify: see the base::apply documentation.
**Examples**

```r
library(bigmemory)
options(bigmemory.typecast.warning=FALSE)
x <- big.matrix(5, 2, type="integer", init=0,
dimnames=list(NULL, c("alpha", "beta")))
x[,] <- round(rnorm(10))
biganalytics::apply(x, 1, mean)
```

**Description**

This is a wrapper to Thomas Lumley’s `biglm` package, allowing it to be used with massive data stored in `big.matrix` objects.

**Usage**

```r
bigglm.big.matrix(
  formula,
  data,
  chunksize = NULL,
  ...,
  fc = NULL,
  getNextChunkFunc = NULL
)
```

```r
biglm.big.matrix(
  formula,
  data,
  chunksize = NULL,
  ...,
  fc = NULL,
  getNextChunkFunc = NULL
)
```

**Arguments**

- `formula`: a model formula.
- `data`: a `big.matrix`.
- `chunksize`: an integer maximum size of chunks of data to process iteratively.
- `fc`: either column indices or names of variables that are factors.
- `...`: options associated with the `biglm`.
- `getNextChunkFunc`: a function which retrieves chunk data.
bigkmeans

Memory-efficient k-means cluster analysis

Description

k-means cluster analysis without the memory overhead, and possibly in parallel using shared memory.

Usage

bigkmeans(x, centers, iter.max = 10, nstart = 1, dist = "euclid")

Arguments

x a big.matrix object.

centers a scalar denoting the number of clusters, or for k clusters, a k by ncol(x) matrix.

iter.max the maximum number of iterations.

nstart number of random starts, to be done in parallel if there is a registered backend (see below).

dist the distance function. Can be "euclid" or "cosine".
Details

The real benefit is the lack of memory overhead compared to the standard \texttt{kmeans} function. Part of the overhead from \texttt{kmeans()} stems from the way it looks for unique starting centers, and could be improved upon. The \texttt{bigkmeans()} function works on either regular \texttt{R} matrix objects, or on \texttt{big.matrix} objects. In either case, it requires no extra memory (beyond the data, other than recording the cluster memberships), whereas \texttt{kmeans()} makes at least two extra copies of the data. And \texttt{kmeans()} is even worse if multiple starts (\texttt{nstart}>1) are used. If \texttt{nstart}>1 and you are using \texttt{bigkmeans()} in parallel, a vector of cluster memberships will need to be stored for each worker, which could be memory-intensive for large data. This isn’t a problem if you use are running the multiple starts sequentially.

Unless you have a really big data set (where a single run of \texttt{kmeans} not only burns memory but takes more than a few seconds), use of parallel computing for multiple random starts is unlikely to be much faster than running iteratively.

Only the algorithm by MacQueen is used here.

Value

An object of class \texttt{kmeans}, just as produced by \texttt{kmeans}.

Note

A comment should be made about the excellent package \texttt{foreach}. By default, it provides \texttt{foreach}, which is used much like a for loop, here over the \texttt{nstart} and doing a final comparison of all results).

When a parallel backend has been registered (see packages \texttt{doSNOW}, \texttt{doMC}, and \texttt{doMPI}, for example), \texttt{bigkmeans()} automatically distributes the \texttt{nstart} random starting points across the available workers. This is done in shared memory on an SMP, but is distributed on a cluster *IF* the \texttt{big.matrix} is file-backed. If used on a cluster with an in-RAM \texttt{big.matrix}, it will fail horribly. We’re considering an extra option as an alternative to the current behavior.

\begin{verbatim}
\textbf{binit}\phantom{10}
\textit{Count elements appearing in bins of one or two variables}
\end{verbatim}

Description

Provides preliminary counting functionality to eventually support graphical exploration or as an alternative to \texttt{table}. Note the availability of \texttt{bigtabulate}.

Usage

\begin{verbatim}
binit(x, cols, breaks = 10)
\end{verbatim}
Arguments

x  
a big.matrix or a matrix.

cols  
a vector of column indices or names of length 1 or 2.

breaks  
a number of bins to span the range from the maximum to the minimum, or a vector (1-variable case) or list of two vectors (2-variable case) where each vector is a triplet of min, max, and number of bins.

Details

The user may specify the number of bins to be used, of equal widths, spanning the range of the data (the default is 10 bins). The user may also specify the range to be spanned along with the number of bins, in case a summary of a subrange of the data is desired. Either univariate or bivariate counting is supported.

The function uses left-closed intervals [a,b) except in the right-most bin, where the interval is entirely closed.

Value

a list containing (a) a vector (1-variable case) or a matrix (2-variable case) of counts of the numbers of cases appearing in each of the bins, (b) description(s) of bin centers, and (c) description(s) of breaks between the bins.

Examples

```r
y <- matrix(rnorm(40), 20, 2)
y[,1] <- NA
x <- as.big.matrix(y, type="double")
x[,]
binit(y, 1:2, list(c(-1,1,5), c(-1,1,2)))
binit(x, 1:2, list(c(-1,1,5), c(-1,1,2)))
binit(y, 1:2)
binit(x, 1:2)

binit(y, 1:2, 5)
binit(x, 1:2, 5)

binit(y, 1)
binit(x, 1)

x <- as.big.matrix(matrix(rnorm(400), 200, 2), type="double")
x[,1] <- x[,1] + 3
x.binit <- binit(x, 1:2)
filled.contour(round(x.binit$rowcenters,2), round(x.binit$colcenters,2),
x.binit$counts, xlab="Variable 1",
ylab="Variable 2")
```
Description

Functions operate on columns of a big.matrix object

Usage

colmin(x, cols = NULL, na.rm = FALSE)

## S4 method for signature 'big.matrix'
colmin(x, cols = NULL, na.rm = FALSE)

## S4 method for signature 'big.matrix'
min(x, ..., na.rm = FALSE)

colmax(x, cols = NULL, na.rm = FALSE)

## S4 method for signature 'big.matrix'
colmax(x, cols = NULL, na.rm = FALSE)

## S4 method for signature 'big.matrix'
max(x, ..., na.rm = FALSE)

colprod(x, cols = NULL, na.rm = FALSE)

## S4 method for signature 'big.matrix'
colprod(x, cols = NULL, na.rm = FALSE)

## S4 method for signature 'big.matrix'
prod(x, ..., na.rm = FALSE)

colsum(x, cols = NULL, na.rm = FALSE)

## S4 method for signature 'big.matrix'
colsum(x, cols = NULL, na.rm = FALSE)

## S4 method for signature 'big.matrix'
sum(x, ..., na.rm = FALSE)

colrange(x, cols = NULL, na.rm = FALSE)

## S4 method for signature 'big.matrix'
colrange(x, cols = NULL, na.rm = FALSE)
range(x, ..., na.rm = FALSE)
colmean(x, cols = NULL, na.rm = FALSE)
## S4 method for signature 'big.matrix'
colmean(x, cols = NULL, na.rm = FALSE)
## S4 method for signature 'big.matrix'
mean(x, ...)
colvar(x, cols = NULL, na.rm = FALSE)
## S4 method for signature 'big.matrix'
colvar(x, cols = NULL, na.rm = FALSE)
colsd(x, cols = NULL, na.rm = FALSE)
## S4 method for signature 'big.matrix'
colsd(x, cols = NULL, na.rm = FALSE)
colna(x, cols = NULL)
## S4 method for signature 'big.matrix'
colna(x, cols = NULL)
## S4 method for signature 'big.matrix'
summary(object)

**Arguments**

- `x`  
  a `big.matrix` object.
- `cols`  
  a scalar or vector of column(s) to be summarized.
- `na.rm`  
  if `TRUE`, remove NA values before summarizing.
- `...`  
  options associated with the corresponding default R function.
- `object`  
  a `big.matrix` object.

**Details**

These functions essentially apply summary functions to each column (or each specified column) of the `big.matrix` in turn.

**Value**

For `colrange`, a matrix with two columns and `length(cols)` rows; column 1 contains the minimum, and column 2 contains the maximum for that column. The other functions return vectors of length `length(cols)`.
Examples

```r
x <- as.big.matrix(
  matrix( sample(1:10, 20, replace=TRUE), 5, 4,
     dimnames=list( NULL, c("a", "b", "c", "d")) ) )

x[,]
mean(x)
colmean(x)
colmin(x)
colmax(x)
colmax(x, "b")
colsd(x)
colrange(x)
range(x)
colsum(x)
colprod(x)
```
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