## Package: matrixStats (via r-universe)

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## Contents

matrixStats-package ..... 2
anyMissing ..... 3
binCounts ..... 4
binMeans ..... 5
indexByRow ..... 7
logSumExp ..... 8
product ..... 10
rowAlls ..... 11
rowCollapse ..... 13
rowCounts ..... 14
rowCumsums ..... 16
rowDiffs ..... 18
rowIQRs ..... 19
rowLogSumExps ..... 20
rowMads ..... 21
rowMeans2 ..... 22
rowMedians ..... 23
rowOrderStats ..... 24
rowQuantiles ..... 26
rowRanges ..... 27
rowRanks ..... 28
rowSums2 ..... 30
rowTabulates ..... 31
rowVars ..... 33
rowWeightedMeans ..... 35
rowWeightedMedians ..... 37
varDiff ..... 38
weightedMad ..... 40
weightedMean ..... 42
weightedMedian ..... 44
weightedVar ..... 46
Index ..... 49
matrixStats-package Package matrixStats

## Description

High-performing functions operating on rows and columns of matrices, e.g. col / rowMedians(), col / rowRanks(), and col / rowSds(). Functions optimized per data type and for subsetted calculations such that both memory usage and processing time is minimized. There are also optimized vectorbased methods, e.g. binMeans(), madDiff() and weightedMedian().

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## Author(s)

Henrik Bengtsson, Hector Corrada Bravo, Robert Gentleman, Ola Hossjer, Harris Jaffee, Dongcan Jiang, Peter Langfelder

## See Also

Useful links:

- https://github.com/HenrikBengtsson/matrixStats
- Report bugs at https://github.com/HenrikBengtsson/matrixStats/issues

```
anyMissing Checks if there are any missing values in an object or not
```


## Description

Checks if there are any missing values in an object or not. Please use base: : anyNA() instead of anyMissing(), colAnyNAs() instead of colAnyMissings(), and rowAnyNAs() instead of rowAnyMissings().

## Usage

```
anyMissing(x, idxs = NULL, ...)
colAnyMissings(x, rows = NULL, cols = NULL, ..., useNames = TRUE)
rowAnyMissings(x, rows = NULL, cols = NULL, ..., useNames = TRUE)
colAnyNAs(x, rows = NULL, cols = NULL, ..., useNames = TRUE)
rowAnyNAs(x, rows = NULL, cols = NULL, ..., useNames = TRUE)
```


## Arguments

x
... Not used.
useNames
idxs A vector indicating subset of elements to operate over. If NULL, no subsetting is done.
rows A vector indicating subset of rows to operate over. If NULL, no subsetting is done.
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
A vector, a list, a matrix, a data. frame, or NULL.

If TRUE (default), names attributes of the result are set, otherwise not.

## Details

The implementation of this method is optimized for both speed and memory. The method will return TRUE as soon as a missing value is detected.

## Value

Returns TRUE if a missing value was detected, otherwise FALSE.

## Author(s)

## Henrik Bengtsson

## See Also

Starting with R v3.1.0, there is anyNA() in the base, which provides the same functionality as anyMissing().

## Examples

```
x <- rnorm(n = 1000)
x[seq(300, length(x), by = 100)] <- NA
stopifnot(anyMissing(x) == any(is.na(x)))
```


## binCounts Fast element counting in non-overlapping bins

## Description

Counts the number of elements in non-overlapping bins

## Usage

binCounts(x, idxs = NULL, bx, right = FALSE, ...)

## Arguments

x
idxs
$b x \quad$ A numeric vector of $B+1$ ordered positions specifying the $B>0$ bins [bx[1], $b x[2]),[b x[2], b x[3]), \ldots,[b x[B], b x[B+1])$.
right If TRUE, the bins are right-closed (left open), otherwise left-closed (right open).
... Not used.

## Details

binCounts ( $x$, $b x$, right $=$ TRUE ) gives equivalent results as rev(binCounts $(-x, b x=r e v(-b x)$, right $=$ FALSE $)$ ), but is faster and more memory efficient.

## Value

Returns an integer vector of length B with non-negative integers.

## Missing and non-finite values

Missing values in x are ignored/dropped. Missing values in bx are not allowed and gives an error.

## Author(s)

Henrik Bengtsson

## See Also

An alternative for counting occurrences within bins is hist, e.g. hist ( x , breaks $=\mathrm{bx}, \mathrm{plot}=$ FALSE) $\$$ counts. That approach is $\sim 30-60 \%$ slower than binCounts (... , right $=$ TRUE).
To count occurrences of indices $x$ (positive integers) in [1, B], use tabulate ( x , nbins $=\mathrm{B}$ ), where $x$ does not have to be sorted first. For details, see tabulate().
To average values within bins, see binMeans().

## binMeans Fast mean calculations in non-overlapping bins

## Description

Computes the sample means in non-overlapping bins

## Usage

binMeans(y, x, idxs = NULL, bx, na.rm = TRUE, count = TRUE, right $=$ FALSE, $\ldots$ )

## Arguments

| $y$ | A numeric or logical vector of $K$ values to calculate means on. |
| :--- | :--- |
| $x$ | A numeric vector of $K$ positions for to be binned. |
| $i d x s$ | A vector indicating subset of elements to operate over. If NULL, no subsetting <br> is done. |
| $b x$ | A numeric vector of $B+1$ ordered positions specifying the $B>0$ bins $[b x[1]$, <br> $b x[2]),[b x[2], b x[3]), \ldots,[b x[B], b x[B+1])$. |
| na.rm | If TRUE, missing values in $y$ are dropped before calculating the mean, otherwise <br> not. |

$$
\begin{array}{ll}
\text { count } & \begin{array}{l}
\text { If TRUE, the number of data points in each bins is returned as attribute count, } \\
\text { which is an integer vector of length B. }
\end{array} \\
\text { right } & \text { If TRUE, the bins are right-closed (left open), otherwise left-closed (right open). } \\
\ldots & \text { Not used. }
\end{array}
$$

## Details

binMeans ( $x$, bx, right $=T R U E$ ) gives equivalent results as rev(binMeans $(-x, b x=\operatorname{sort}(-b x)$, right = FALSE) ), but is faster.

## Value

Returns a numeric vector of length $B$.

## Missing and non-finite values

Data points where either of y and x is missing are dropped (and therefore are also not counted). Non-finite values in $y$ are not allowed and gives an error. Missing values in bx are not allowed and gives an error.

## Author(s)

Henrik Bengtsson with initial code contributions by Martin Morgan [1].

## References

[1] R-devel thread Fastest non-overlapping binning mean function out there? on Oct 3, 2012

## See Also

binCounts(). aggregate and mean().

## Examples

```
x <- 1:200
mu <- double(length(x))
mu[1:50] <- 5
mu[101:150] <- -5
y <- mu + rnorm(length(x))
# Binning
bx <- c(0, 50, 100, 150, 200) + 0.5
y_s <- binMeans(y, x = x, bx = bx)
plot(x, y)
for (kk in seq_along(y_s)) {
    lines(bx[c(kk, kk + 1)], y_s[c(kk, kk)], col = "blue", lwd = 2)
}
```


## Description

Translates matrix indices by rows into indices by columns.

## Usage

indexByRow(dim, idxs = NULL, ...)

## Arguments

| dim | A numeric vector of length two specifying the length of the "template" matrix. |
| :--- | :--- |
| idxs | A vector indicating subset of elements to operate over. If NULL, no subsetting |
| is done. |  |
| $\ldots$ | Not used. |

## Value

Returns an integer vector of indices.

## Known limitations

The current implementation does not support long-vector indices, because both input and output indices are of type integers. This means that the indices in argument idxs can only be in range [1, $\left.2^{\wedge} 31-1\right]$. Using a greater value will be coerced to NA_integer_. Moreover, returned indices can only be in the same range $\left[1,2^{\wedge} 31-1\right]$.

## Author(s)

## Henrik Bengtsson

## Examples

```
dim <- c(5, 4)
X <- matrix(NA_integer_, nrow = dim[1], ncol = dim[2])
Y <- t(X)
idxs <- seq_along(X)
# Assign by columns
X[idxs] <- idxs
print(X)
# Assign by rows
Y[indexByRow(dim(Y), idxs)] <- idxs
print(Y)
stopifnot(X == t(Y))
```


## Description

Accurately computes the logarithm of the sum of exponentials, that is, $\log (\operatorname{sum}(\exp (l x)))$. If $l x=\log (x)$, then this is equivalently to calculating $\log (\operatorname{sum}(x))$.

## Usage

$\operatorname{logSumExp}(1 x, i d x s=$ NULL, na.rm = FALSE, $\ldots$ )

## Arguments

| $l x$ | A numeric vector. Typically $l x \operatorname{are} \log (x)$ values. |
| :--- | :--- |
| idxs | A vector indicating subset of elements to operate over. If NULL, no subsetting |
| is done. |  |
| na.rm | If TRUE, missing values are excluded. |
| $\ldots$ | Not used. |

## Details

This function, which avoid numerical underflow, is often used when computing the logarithm of the sum of small numbers $(|x| \ll 1)$ such as probabilities.
This is function is more accurate than $\log (\operatorname{sum}(\exp (l x)))$ when the values of $x=\exp (l x)$ are $|x| \ll 1$. The implementation of this function is based on the observation that
$\log (a+b)=[l a=\log (a), l b=\log (b)]=\log (\exp (l a)+\exp (l b))=l a+\log (1+\exp (l b-l a))$
Assuming $l a>l b$, then $|l b-l a|<|l b|$, and it is less likely that the computation of $1+\exp (l b-l a)$ will not underflow/overflow numerically. Because of this, the overall result from this function should be more accurate. Analogously to this, the implementation of this function finds the maximum value of $1 x$ and subtracts it from the remaining values in $1 x$.

## Value

Returns a numeric scalar.

## Benchmarking

This method is optimized for correctness, that avoiding underflowing. It is implemented in native code that is optimized for speed and memory.

## Author(s)

Henrik Bengtsson

## References

[1] R Core Team, Writing R Extensions, v3.0.0, April 2013.
[2] Laurent El Ghaoui, Hyper-Textbook: Optimization Models and Applications, University of California at Berkeley, August 2012. (Chapter 'Log-Sum-Exp (LSE) Function and Properties')
[3] R-help thread logsumexp function in R, 2011-02-17. https://stat.ethz.ch/pipermail/ r-help/2011-February/269205.html

## See Also

To compute this function on rows or columns of a matrix, see rowLogSumExps().
For adding two double values in native code, R provides the C function logspace_add() [1]. For properties of the log-sum-exponential function, see [2].

## Examples

```
## EXAMPLE #1
lx <- c(1000.01, 1000.02)
y0 <- log(sum(exp(lx)))
print(y0) ## Inf
y1 <- logSumExp(lx)
print(y1) ## 1000.708
## EXAMPLE #2
lx <- c(-1000.01, -1000.02)
y0 <- log(sum(exp(lx)))
print(y0) ## -Inf
y1 <- logSumExp(lx)
print(y1) ## -999.3218
## EXAMPLE #3
## R-help thread 'Beyond double-precision?' on May 9, 2009.
set.seed(1)
x <- runif(50)
## The logarithm of the harmonic mean
y0 <- log(1 / mean(1 / x))
print(y0) ## -1.600885
lx <- log(x)
y1 <- log(length(x)) - logSumExp(-lx)
print(y1) ## [1] -1.600885
# Sanity check
stopifnot(all.equal(y1, y0))
```


## Description

Calculates the product for each row (column) in a matrix.

## Usage

product $(x$, idxs $=$ NULL, na.rm $=$ FALSE, ...)
rowProds(x, rows $=$ NULL, cols $=$ NULL, na.rm $=$ FALSE, method = c("direct", "expSumLog"), ..., useNames = TRUE)
colProds $(x$, rows $=$ NULL, cols $=$ NULL, na.rm $=$ FALSE, method = c("direct", "expSumLog"), ..., useNames = TRUE)

## Arguments

| $x$ | An NxK matrix or, if dim. is specified, an $\mathrm{N} * \mathrm{~K}$ vector. |
| :--- | :--- |
| idxs | A vector indicating subset of elements to operate over. If NULL, no subsetting <br> is done. |
| na.rm | If TRUE, missing values are excluded. |
| $\ldots$ | Not used. |
| rows | A vector indicating subset of rows to operate over. If NULL, no subsetting is <br> done. |
| cols | A vector indicating subset of columns to operate over. If NULL, no subsetting <br> is done. |
| method | A character string specifying how each product is calculated. |
| useNames | If TRUE (default), names attributes of the result are set, otherwise not. |

## Details

If method = "expSumLog", then then product() function is used, which calculates the product via the logarithmic transform (treating negative values specially). This improves the precision and lowers the risk for numeric overflow. If method = "direct", the direct product is calculated via the $\operatorname{prod}()$ function.

## Value

Returns a numeric vector of length $\mathrm{N}(\mathrm{K})$.

## Missing values

Note, if method = "expSumLog", na.rm = FALSE, and $x$ contains missing values (NA or NaN), then the calculated value is also missing value. Note that it depends on platform whether NaN or NA is returned when an NaN exists, cf. is. nan().

## Author(s)

Henrik Bengtsson

```
rowAlls
```

Checks if a value exists / does not exist in each row (column) of a matrix

## Description

Checks if a value exists / does not exist in each row (column) of a matrix.

## Usage

```
    rowAlls (x, rows \(=\) NULL, cols \(=\) NULL, value \(=\) TRUE, na. \(\mathrm{rm}=\) FALSE,
        \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\), useNames \(=\) TRUE \()\)
    colAlls(x, rows \(=\) NULL, cols \(=\) NULL, value \(=\) TRUE, na.rm \(=\) FALSE,
        dim. \(=\operatorname{dim}(x), \ldots\) useNames \(=\) TRUE)
    allValue(x, idxs = NULL, value = TRUE, na.rm = FALSE, ...)
    rowAnys(x, rows \(=\) NULL, cols \(=\) NULL, value \(=\) TRUE, na.rm \(=\) FALSE,
        dim. \(=\operatorname{dim}(x), \ldots\), useNames \(=\) TRUE)
    colAnys \((x\), rows \(=\) NULL, cols \(=\) NULL, value \(=\) TRUE, na.rm \(=\) FALSE,
        \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\) useNames \(=\) TRUE)
    anyValue(x, idxs = NULL, value = TRUE, na.rm = FALSE, ...)
```


## Arguments

x
rows A vector indicating subset of rows to operate over. If NULL, no subsetting is done.
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
value A value to search for.
na.rm If TRUE, missing values are excluded.
dim. An integer vector of length two specifying the dimension of $x$, also when not a matrix. Comment: The reason for this argument being named with a period at the end is purely technical (we get a run-time error if we try to name it dim).
... Not used.
useNames If TRUE (default), names attributes of the result are set, otherwise not.
idxs A vector indicating subset of elements to operate over. If NULL, no subsetting is done.

## Details

These functions takes either a matrix or a vector as input. If a vector, then argument dim. must be specified and fulfill prod(dim.) $==$ length $(x)$. The result will be identical to the results obtained when passing matrix $(x$, nrow $=\operatorname{dim} .[1 \mathrm{~L}]$, ncol $=\operatorname{dim} .[2 \mathrm{~L}]$ ), but avoids having to temporarily create/allocate a matrix, if only such is needed only for these calculations.

## Value

rowAlls() (colAlls()) returns an logical vector of length $\mathrm{N}(\mathrm{K})$. Analogously for rowAnys() (rowAlls()).

## Logical value

When value is logical, the result is as if the function is applied on as.logical( $x$ ). More specifically, if $x$ is numeric, then all zeros are treated as FALSE, non-zero values as TRUE, and all missing values as NA.

## Author(s)

Henrik Bengtsson

## See Also

rowCounts

## Examples

```
x <- matrix(FALSE, nrow = 10, ncol = 5)
x[3:7, c(2, 4)] <- TRUE
x[2:4, ] <- TRUE
x[, 1] <- TRUE
x[5, ] <- FALSE
x[, 5] <- FALSE
print(x)
print(rowCounts(x)) # 1 4 4 4 0 3 3 1 1 1
print(colCounts(x)) # 9 5 3 5 0
print(rowAnys(x))
print(which(rowAnys(x))) # 1 1 2 < 3
print(colAnys(x))
print(which(colAnys(x))) # 1 2 3 4
```

```
rowCollapse Extracts one cell per row (column) from a matrix
```


## Description

Extracts one cell per row (column) from a matrix. The implementation is optimized for memory and speed.

## Usage

rowCollapse(x, idxs, rows $=$ NULL, $\operatorname{dim} .=\operatorname{dim}(x), \ldots$, useNames $=$ TRUE)
colCollapse(x, idxs, cols = NULL, dim. $=\operatorname{dim}(x), \ldots$, useNames $=$ TRUE)

## Arguments

x
An NxK matrix or, if dim. is specified, an $\mathrm{N} * \mathrm{~K}$ vector.
idxs An index vector of (maximum) length $N(K)$ specifying the columns (rows) to be extracted.
rows A vector indicating subset of rows to operate over. If NULL, no subsetting is done.
dim. An integer vector of length two specifying the dimension of $x$, also when not a matrix. Comment: The reason for this argument being named with a period at the end is purely technical (we get a run-time error if we try to name it dim).
... Not used.
useNames If TRUE (default), names attributes of the result are set, otherwise not.
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.

## Value

Returns a vector of length $\mathrm{N}(\mathrm{K})$.

## Author(s)

Henrik Bengtsson

## See Also

Matrix indexing to index elements in matrices and arrays, cf. [().

## Examples

```
x <- matrix(1:27, ncol = 3)
y <- rowCollapse(x, 1)
stopifnot(identical(y, x[, 1]))
y <- rowCollapse(x, 2)
stopifnot(identical(y, x[, 2]))
y <- rowCollapse(x, c(1, 1, 1, 1, 1, 3, 3, 3, 3))
stopifnot(identical(y, c(x[1:5, 1], x[6:9, 3])))
y <- rowCollapse(x, 1:3)
print(y)
y_truth <- c(x[1, 1], x[2, 2], x[3, 3], x[4, 1], x[5, 2],
    x[6, 3], x[7, 1], x[8, 2], x[9, 3])
stopifnot(identical(y, y_truth))
```

rowCounts

Counts the number of occurrences of a specific value

## Description

The row- and column-wise functions take either a matrix or a vector as input. If a vector, then argument dim. must be specified and fulfill prod (dim.) $==$ length $(x)$. The result will be identical to the results obtained when passing matrix (x, nrow = dim.[1L], ncol = dim.[2L]), but avoids having to temporarily create/allocate a matrix, if only such is needed only for these calculations.

## Usage

rowCounts (x, rows $=$ NULL, cols $=$ NULL, value $=$ TRUE, na.rm = FALSE, dim. $=\operatorname{dim}(x), \ldots$, useNames $=$ TRUE)
colCounts ( $x$, rows $=$ NULL, cols $=$ NULL, value $=$ TRUE, na.rm = FALSE, $\operatorname{dim} .=\operatorname{dim}(x), \ldots$, useNames $=$ TRUE $)$
count (x, idxs = NULL, value = TRUE, na.rm = FALSE, ...)

## Arguments

x
rows A vector indicating subset of rows to operate over. If NULL, no subsetting is done.
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
value A value to search for.
na.rm
An NxK matrix or, if dim. is specified, an $\mathrm{N}^{*} \mathrm{~K}$ vector.
.
If TRUE, missing values are excluded.

| dim. | An integer vector of length two specifying the dimension of $x$, also when not <br> a matrix. Comment: The reason for this argument being named with a period <br> at the end is purely technical (we get a run-time error if we try to name it dim). |
| :--- | :--- |
| $\ldots$ | Not used. |
| useNames | If TRUE (default), names attributes of the result are set, otherwise not. |
| idxs | A vector indicating subset of elements to operate over. If NULL, no subsetting <br> is done. |

## Value

rowCounts() (colCounts()) returns an integer vector of length $N(K)$. count () returns a scalar of type integer if the count is less than 2^31-1 (= . Machine\$integer.max) otherwise a scalar of type double.

## Author(s)

Henrik Bengtsson

## See Also

rowAlls

## Examples

```
x <- matrix(0:11, nrow = 4, ncol = 3)
x[2:3, 2:3] <- 2:5
x[3, 3] <- NA_integer_
print(x)
print(rowCounts(x, value = 2))
## [1] 0 1 NA 0
print(colCounts(x, value = 2))
## [1] 1 1 NA
print(colCounts(x, value = NA_integer_))
## [1] 0 0 1
print(rowCounts(x, value = 2, na.rm = TRUE))
## [1] 0 1 1 0
print(colCounts(x, value = 2, na.rm = TRUE))
## [1] 1 1 0
print(rowAnys(x, value = 2))
## [1] FALSE TRUE TRUE FALSE
print(rowAnys(x, value = NA_integer_))
## [1] FALSE FALSE TRUE FALSE
print(colAnys(x, value = 2))
## [1] TRUE TRUE NA
print(colAnys(x, value = 2, na.rm = TRUE))
## [1] TRUE TRUE FALSE
```

```
print(colAlls(x, value = 2))
## [1] FALSE FALSE FALSE
```

```
rowCumsums
```

Cumulative sums, products, minima and maxima for each row (column) in a matrix

## Description

Cumulative sums, products, minima and maxima for each row (column) in a matrix.

## Usage

```
rowCumsums (x, rows \(=\) NULL, cols \(=\) NULL, \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\),
    useNames = TRUE)
    colCumsums(x, rows \(=\) NULL, cols \(=\) NULL, \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\),
        useNames = TRUE)
    rowCumprods(x, rows \(=\) NULL, cols \(=\) NULL, \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\),
        useNames = TRUE)
    colCumprods \((x\), rows \(=\) NULL, cols \(=\) NULL, \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\),
        useNames = TRUE)
    rowCummins(x, rows \(=\) NULL, cols \(=\) NULL, \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\),
    useNames = TRUE)
    colCummins(x, rows \(=\) NULL, cols \(=\) NULL, \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\),
        useNames = TRUE)
    rowCummaxs(x, rows \(=\) NULL, cols \(=\) NULL, \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\),
    useNames \(=\) TRUE)
    colCummaxs(x, rows \(=\) NULL, cols \(=\) NULL, \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\),
    useNames = TRUE)
```


## Arguments

| $x$ | An NxK matrix or, if dim. is specified, an $\mathrm{N} * \mathrm{~K}$ vector. |
| :--- | :--- |
| rows | A vector indicating subset of rows to operate over. If NULL, no subsetting is <br> done. |
| cols | A vector indicating subset of columns to operate over. If NULL, no subsetting <br> is done. |
| dim. | An integer vector of length two specifying the dimension of $x$, also when not <br> a matrix. Comment: The reason for this argument being named with a period <br> at the end is purely technical (we get a run-time error if we try to name it dim). |


| $\ldots$ | Not used. |
| :--- | :--- |
| useNames | If TRUE (default), names attributes of the result are set, otherwise not. |

## Value

Returns a numeric NxK matrix of the same mode as $x$, except when $x$ is of mode logical, then the return type is integer.

## Author(s)

Henrik Bengtsson

## See Also

See cumsum(), cumprod(), cummin(), and cummax().

## Examples

```
x <- matrix(1:12, nrow = 4, ncol = 3)
print(x)
yr <- rowCumsums(x)
print(yr)
yc <- colCumsums(x)
print(yc)
yr <- rowCumprods(x)
print(yr)
yc <- colCumprods(x)
print(yc)
yr <- rowCummaxs(x)
print(yr)
yc <- colCummaxs(x)
print(yc)
yr <- rowCummins(x)
print(yr)
yc <- colCummins(x)
print(yc)
```

```
rowDiffs Calculates difference for each row (column) in a matrix
```


## Description

Calculates difference for each row (column) in a matrix.

## Usage

```
rowDiffs(x, rows \(=\) NULL, cols \(=\) NULL, lag = 1L, differences = 1L,
    dim. = dim(x), ..., useNames = TRUE)
    colDiffs(x, rows \(=\) NULL, cols \(=\) NULL, lag = 1L, differences = 1L,
    \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\), useNames \(=\) TRUE \()\)
```


## Arguments

x
rows A vector indicating subset of rows to operate over. If NULL, no subsetting is done.
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
lag An integer specifying the lag.
differences
dim. An integer vector of length two specifying the dimension of $x$, also when not a matrix. Comment: The reason for this argument being named with a period at the end is purely technical (we get a run-time error if we try to name it dim).
... Not used.
useNames If TRUE (default), names attributes of the result are set, otherwise not.

## Value

Returns a numeric $\mathrm{Nx}(\mathrm{K}-1)$ or (N-1)xK matrix.

## Author(s)

Henrik Bengtsson

## See Also

See also diff2().

## Examples

```
x <- matrix(1:27, ncol = 3)
d1 <- rowDiffs(x)
print(d1)
d2 <- t(colDiffs(t(x)))
stopifnot(all.equal(d2, d1))
```

```
rowIQRs
```

Estimates of the interquartile range for each row (column) in a matrix

## Description

Estimates of the interquartile range for each row (column) in a matrix.

## Usage

```
rowIQRs(x, rows = NULL, cols = NULL, na.rm = FALSE, ...,
    useNames = TRUE)
colIQRs(x, rows = NULL, cols = NULL, na.rm = FALSE, ...,
    useNames = TRUE)
iqr(x, idxs = NULL, na.rm = FALSE, ...)
```


## Arguments

x
rows A vector indicating subset of rows to operate over. If NULL, no subsetting is done.
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
na.rm If TRUE, missing values are excluded.
Additional arguments passed to rowQuantiles() (colQuantiles()).
useNames
idxs A vector indicating subset of elements to operate over. If NULL, no subsetting is done.

## Value

Returns a numeric vector of length $\mathrm{N}(\mathrm{K})$.

## Missing values

Contrary to IQR, which gives an error if there are missing values and na. rm = FALSE, iqr() and its corresponding row and column-specific functions return NA_real_.

## Author(s)

Henrik Bengtsson

## See Also

See IQR. See rowSds().

## Examples

```
set.seed(1)
x <- matrix(rnorm(50 * 40), nrow = 50, ncol = 40)
str(x)
# Row IQRs
q <- rowIQRs(x)
print(q)
q0 <- apply(x, MARGIN = 1, FUN = IQR)
stopifnot(all.equal(q0, q))
# Column IQRs
q <- colIQRs(x)
print(q)
q0 <- apply(x, MARGIN = 2, FUN = IQR)
stopifnot(all.equal(q0,q))
```

```
    rowLogSumExps
```

Accurately computes the logarithm of the sum of exponentials across rows or columns

## Description

Accurately computes the logarithm of the sum of exponentials across rows or columns.

## Usage

```
rowLogSumExps(lx, rows = NULL, cols = NULL, na.rm = FALSE,
        dim. = dim(lx), ..., useNames = TRUE)
    colLogSumExps(lx, rows = NULL, cols = NULL, na.rm = FALSE,
        dim. = dim(lx), ..., useNames = TRUE)
```


## Arguments

$1 x$
rows, cols
na.rm

A numeric NxK matrix. Typically 1 x are $\log (x)$ values.
A vector indicating subset of rows (and/or columns) to operate over. If NULL, no subsetting is done.
If TRUE, any missing values are ignored, otherwise not.
dim. An integer vector of length two specifying the dimension of $x$, also when not a matrix.
... Not used.
useNames If TRUE (default), names attributes of the result are set, otherwise not.

## Value

A numeric vector of length $N(K)$.

## Benchmarking

These methods are implemented in native code and have been optimized for speed and memory.

## Author(s)

Native implementation by Henrik Bengtsson. Original R code by Nakayama ??? (Japan).

## See Also

To calculate the same on vectors, logSumExp().

```
rowMads
Standard deviation estimates for each row (column) in a matrix
```


## Description

Standard deviation estimates for each row (column) in a matrix.

## Usage

```
    rowMads(x, rows = NULL, cols = NULL, center = NULL, constant = 1.4826,
        na.rm = FALSE, dim. = dim(x), ..., useNames = TRUE)
    colMads(x, rows = NULL, cols = NULL, center = NULL, constant = 1.4826,
        na.rm = FALSE, dim. = dim(x), ..., useNames = TRUE)
    rowSds(x, rows = NULL, cols = NULL, na.rm = FALSE, refine = TRUE,
        center = NULL, dim. = dim(x), ..., useNames = TRUE)
    colSds(x, rows = NULL, cols = NULL, na.rm = FALSE, refine = TRUE,
        center = NULL, dim. = dim(x), ..., useNames = TRUE)
```


## Arguments

| x |  |
| :--- | :--- |
| rows | An NxK matrix or, if dim. is specified, an $\mathrm{N} * \mathrm{~K}$ vector. <br> A vector indicating subset of rows to operate over. If NULL, no subsetting is <br> done. |
| cols | A vector indicating subset of columns to operate over. If NULL, no subsetting <br> is done. <br> (optional) The center, defaults to the row means for the SD estimators and row <br> medians for the MAD estimators. |
| center | A scale factor. See mad for details. |
| na.rm | If TRUE, missing values are excluded. |
| dim. | An integer vector of length two specifying the dimension of x, also when not <br> a matrix. Comment: The reason for this argument being named with a period <br> at the end is purely technical (we get a run-time error if we try to name it dim). |
| $\ldots$ | Additional arguments passed to rowMeans() and rowSums(). <br> useNames |
| If TRUE (default), names attributes of the result are set, otherwise not. |  |

## Value

Returns a numeric vector of length $\mathrm{N}(\mathrm{K})$.

## Author(s)

Henrik Bengtsson

## See Also

sd, mad and var. rowIQRs().

## rowMeans2

Calculates the mean for each row (column) in a matrix

## Description

Calculates the mean for each row (column) in a matrix.

## Usage

rowMeans2(x, rows $=$ NULL, cols $=$ NULL, na.rm = FALSE, refine $=$ TRUE, dim. = dim(x), ..., useNames = TRUE)
colMeans2(x, rows = NULL, cols = NULL, na.rm = FALSE, refine = TRUE, dim. = dim(x), ..., useNames = TRUE)
rowMedians

## Arguments

x
rows
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
na.rm If TRUE, missing values are excluded.
refine If TRUE and $x$ is numeric, then extra effort is used to calculate the average with greater numerical precision, otherwise not.
dim. An integer vector of length two specifying the dimension of $x$, also when not a matrix. Comment: The reason for this argument being named with a period at the end is purely technical (we get a run-time error if we try to name it dim).
... Not used.
useNames If TRUE (default), names attributes of the result are set, otherwise not.

## Details

The implementation of rowMeans2() and colMeans2() is optimized for both speed and memory.

## Value

Returns a numeric vector of length $\mathrm{N}(\mathrm{K})$.

## Author(s)

Henrik Bengtsson

## Description

Calculates the median for each row (column) in a matrix.

## Usage

rowMedians(x, rows $=$ NULL, cols $=$ NULL, na.rm = FALSE, dim. $=\operatorname{dim}(x)$, ..., useNames = TRUE)
colMedians(x, rows $=$ NULL, cols $=$ NULL, na.rm $=$ FALSE, $\operatorname{dim} .=\operatorname{dim}(x)$, ..., useNames = TRUE)

## Arguments

na.rm If TRUE, NAs are excluded first, otherwise not.
dim. An integer vector of length two specifying the dimension of $x$, also when not

X
rows, cols
. . Not used.
useNames a matrix.

An NxK matrix or, if dim. is specified, an $\mathrm{N}^{*} \mathrm{~K}$ vector.
A vector indicating subset of rows (and/or columns) to operate over. If NULL, no subsetting is done.

## Details

The implementation of rowMedians() and colMedians() is optimized for both speed and memory. To avoid coercing to doubles (and hence memory allocation), there is a special implementation for integer matrices. That is, if $x$ is an integer matrix, then rowMedians(as.double(x)) (rowMedians(as.double(x))) would require three times the memory of rowMedians (x) (colMedians(x)), but all this is avoided.

## Value

Returns a numeric vector of length $\mathrm{N}(\mathrm{K})$.

## Author(s)

Henrik Bengtsson, Harris Jaffee

## See Also

See rowWeightedMedians() and colWeightedMedians() for weighted medians. For mean estimates, see rowMeans2() and rowMeans().

```
rowOrderStats

\section*{Description}

Gets an order statistic for each row (column) in a matrix.

\section*{Usage}
rowOrderStats(x, rows \(=\) NULL, cols \(=\) NULL, which, \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\), useNames = TRUE)
colOrderStats(x, rows \(=\) NULL, cols \(=\) NULL, which, \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\), useNames = TRUE)

\section*{Arguments}

X
rows A vector indicating subset of rows to operate over. If NULL, no subsetting is done.
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
which An integer index in \([1, \mathrm{~K}]([1, \mathrm{~N}])\) indicating which order statistic to be returned.
dim. An integer vector of length two specifying the dimension of \(x\), also when not a matrix. Comment: The reason for this argument being named with a period at the end is purely technical (we get a run-time error if we try to name it dim).
... Not used.
useNames If TRUE (default), names attributes of the result are set, otherwise not.

\section*{Details}

The implementation of rowOrderStats() is optimized for both speed and memory. To avoid coercing to doubles (and hence memory allocation), there is a unique implementation for integer matrices.

\section*{Value}

Returns a numeric vector of length \(\mathrm{N}(\mathrm{K})\).

\section*{Missing values}

This method does not handle missing values, that is, the result corresponds to having na. \(\mathrm{rm}=\) FALSE (if such an argument would be available).

\section*{Author(s)}

The native implementation of rowOrderStats() was adopted by Henrik Bengtsson from Robert Gentleman's rowQ() in the Biobase package.

\section*{See Also}

See rowMeans() in colSums().

\section*{Description}

Estimates quantiles for each row (column) in a matrix.

\section*{Usage}
rowQuantiles \((x\), rows \(=\) NULL, cols \(=\) NULL, probs \(=\operatorname{seq}(\) from \(=0\), to \(=1\), by \(=0.25\) ), na.rm \(=\) FALSE, type \(=7 \mathrm{~L}\), digits \(=7 \mathrm{~L}, \ldots\), useNames = TRUE, drop = TRUE)
colQuantiles(x, rows \(=\) NULL, cols \(=\) NULL, probs \(=\operatorname{seq}(f r o m=0\), to \(=1\), by \(=0.25\) ), na.rm = FALSE, type \(=7 \mathrm{~L}\), digits \(=7 \mathrm{~L}, \ldots\), useNames = TRUE, drop = TRUE)

\section*{Arguments}
x
rows
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
probs A numeric vector of J probabilities in \([0,1]\).
na.rm If TRUE, missing values are excluded.
type An integer specifying the type of estimator. See quantile for more details.
digits An integer specifying the precision of the formatted percentages. Not used
when 'useNames \(=\) FALSE'. In \(* *\) matrixStats \(* *(<0.63 .0)\), the default used to be 'max(2L, getOption("digits"))' inline with R (<4.1.0).
... Additional arguments passed to quantile.
useNames If TRUE (default), names attributes of the result are set, otherwise not.
drop
An integer, numeric or logical NxK matrix with \(\mathrm{N}>=0\).
A vector indicating subset of rows to operate over. If NULL, no subsetting is done.

If TRUE, singleton dimensions in the result are dropped, otherwise not.

\section*{Value}

Returns a \(\mathrm{NxJ}(\mathrm{KxJ})\) matrix, where \(\mathrm{N}(\mathrm{K})\) is the number of rows (columns) for which the J quantiles are calculated. The return type is either integer or numeric depending on type.

\section*{Author(s)}

Henrik Bengtsson

\section*{See Also}
quantile.
rowRanges

\section*{Examples}
```

set.seed(1)
x <- matrix(rnorm(50 * 40), nrow = 50, ncol = 40)
str(x)
probs <- c(0.25, 0.5, 0.75)

# Row quantiles

q <- rowQuantiles(x, probs = probs)
print(q)
q_0 <- apply(x, MARGIN = 1, FUN = quantile, probs = probs)
stopifnot(all.equal(q_0, t(q)))

# Column IQRs

q <- colQuantiles(x, probs = probs)
print(q)
q_0 <- apply(x, MARGIN = 2, FUN = quantile, probs = probs)
stopifnot(all.equal(q_0, t(q)))

```
rowRanges

\section*{Description}

Gets the range of values in each row (column) of a matrix.

\section*{Usage}
```

rowRanges(x, rows $=$ NULL, cols $=$ NULL, na.rm = FALSE, dim. $=\operatorname{dim}(x)$,
..., useNames = TRUE)
rowMins(x, rows $=$ NULL, cols $=$ NULL, na.rm $=$ FALSE, $\operatorname{dim} .=\operatorname{dim}(x), \ldots$,
useNames = TRUE)
rowMaxs(x, rows $=$ NULL, cols $=$ NULL, na. $r m=$ FALSE, $\operatorname{dim} .=\operatorname{dim}(x), \ldots$,
useNames = TRUE)
colRanges $(x$, rows $=$ NULL, cols $=$ NULL, na.rm $=$ FALSE, dim. $=\operatorname{dim}(x)$,
..., useNames = TRUE)
colMins(x, rows $=$ NULL, cols $=$ NULL, na.rm $=$ FALSE, $\operatorname{dim} .=\operatorname{dim}(x), \ldots$,
useNames = TRUE)
colMaxs(x, rows $=$ NULL, cols $=$ NULL, na.rm $=$ FALSE, $\operatorname{dim} .=\operatorname{dim}(x), \ldots$,
useNames = TRUE)

```

\section*{Arguments}
x
rows
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
na.rm If TRUE, missing values are excluded.
dim. An integer vector of length two specifying the dimension of \(x\), also when not a matrix. Comment: The reason for this argument being named with a period at the end is purely technical (we get a run-time error if we try to name it dim).
... Not used.
useNames If TRUE (default), names attributes of the result are set, otherwise not.

\section*{Value}
rowRanges() (colRanges()) returns a numeric \(\mathrm{Nx} 2(\mathrm{Kx} 2)\) matrix, where \(\mathrm{N}(\mathrm{K})\) is the number of rows (columns) for which the ranges are calculated.
rowMins()/rowMaxs() (colMins()/colMaxs()) returns a numeric vector of length \(N(K)\).

\section*{Author(s)}

Henrik Bengtsson

\section*{See Also}
rowOrderStats() and pmin.int().
```

rowRanks
Gets the rank of the elements in each row (column) of a matrix

```

\section*{Description}

Gets the rank of the elements in each row (column) of a matrix.

\section*{Usage}
```

    rowRanks(x, rows = NULL, cols = NULL, ties.method = c("max", "average",
        "first", "last", "random", "max", "min", "dense"), dim. = dim(x), ...,
        useNames = TRUE)
    colRanks(x, rows = NULL, cols = NULL, ties.method = c("max", "average",
        "first", "last", "random", "max", "min", "dense"), dim. = dim(x),
        preserveShape = FALSE, ..., useNames = TRUE)
    ```

\section*{Arguments}
x
rows A vector indicating subset of rows to operate over. If NULL, no subsetting is done.
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
ties.method A character string specifying how ties are treated. For details, see below.
dim. An integer vector of length two specifying the dimension of \(x\), also when not a matrix. Comment: The reason for this argument being named with a period at the end is purely technical (we get a run-time error if we try to name it dim).

Not used.
useNames
If TRUE (default), names attributes of the result are set, otherwise not.
preserveShape A logical specifying whether the matrix returned should preserve the input shape of \(x\), or not.

\section*{Details}

These functions rank values and treats missing values the same way as rank(). For equal values ("ties"), argument ties.method determines how these are ranked among each other. More precisely, for the following values of ties.method, each index set of ties consists of:
- "first" - increasing values that are all unique
- "last" - decreasing values that are all unique
- "min" - identical values equaling the minimum of their original ranks
- "max" - identical values equaling the maximum of their original ranks
- "average" - identical values that equal the sample mean of their original ranks. Because the average is calculated, the returned ranks may be non-integer values
- "random" - randomly shuffled values of their original ranks.
- "dense" - increasing values that are all unique and, contrary to "first", never contain any gaps

For more information on ties.method \(=\) "dense", see frank() of the data.table package. For more information on the other alternatives, see rank().

Note that, due to different randomization strategies, the shuffling order produced by these functions when using ties.method = "random" does not reproduce that of rank().

WARNING: For backward-compatibility reasons, the default is ties.method = "max", which differs from rank() which uses ties.method = "average" by default. Since we plan to change the default behavior in a future version, we recommend to explicitly specify the intended value of argument ties.method.

\section*{Value}

A matrix of type integer is returned, unless ties.method = "average" when it is of type numeric.
The rowRanks() function always returns an NxK matrix, where \(N(K)\) is the number of rows (columns) whose ranks are calculated.
The colRanks() function returns an NxK matrix, if preserveShape = TRUE, otherwise a KxN matrix.
Any names of \(x\) are ignored and absent in the result.

\section*{Missing values}

Missing values are ranked as NA_integer_, as with na.last = "keep" in the rank() function.

\section*{Performance}

The implementation is optimized for both speed and memory. To avoid coercing to doubles (and hence memory allocation), there is a unique implementation for integer matrices. Furthermore, it is more memory efficient to do colRanks ( \(x\), preserveShape \(=\) TRUE) than \(t\) (colRanks \((x\), preserveShape \(=\) FALSE) ).

\section*{Author(s)}

Hector Corrada Bravo and Harris Jaffee. Peter Langfelder for adding 'ties.method' support. Brian Montgomery for adding more 'ties.method's. Henrik Bengtsson adapted the original native implementation of rowRanks() from Robert Gentleman's rowQ() in the Biobase package.

\section*{See Also}

For developers, see also Section Utility functions' in 'Writing R Extensions manual', particularly the native functions R_qsort_I () and R_qsort_int_I ().
```

rowSums2 Calculates the sum for each row (column) in a matrix

```

\section*{Description}

Calculates the sum for each row (column) in a matrix.

\section*{Usage}
```

rowSums2(x, rows = NULL, cols = NULL, na.rm = FALSE, dim. = dim(x),
..., useNames = TRUE)
colSums2(x, rows $=$ NULL, cols $=$ NULL, na.rm $=$ FALSE, $\operatorname{dim} .=\operatorname{dim}(x)$,
..., useNames = TRUE)

```

\section*{Arguments}

X
rows
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
na.rm If TRUE, missing values are excluded.
dim.
. . Not used.
useNames If TRUE (default), names attributes of the result are set, otherwise not.

\section*{Details}

The implementation of rowSums2() and colSums2() is optimized for both speed and memory.

\section*{Value}

Returns a numeric vector of length \(\mathrm{N}(\mathrm{K})\).

\section*{Author(s)}

Henrik Bengtsson
```

rowTabulates

```

Tabulates the values in a matrix by row (column).

\section*{Description}

Tabulates the values in a matrix by row (column).

\section*{Usage}
```

rowTabulates(x, rows $=$ NULL, cols $=$ NULL, values = NULL, ...,
useNames = TRUE)
colTabulates(x, rows $=$ NULL, cols $=$ NULL, values $=$ NULL,... ,
useNames = TRUE)

```

\section*{Arguments}

X
rows A vector indicating subset of rows to operate over. If NULL, no subsetting is done.
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
values An vector of \(\mathbf{J}\) values of count. If NULL, all (unique) values are counted.
...
useNames If TRUE (default), names attributes of the result are set, otherwise not.

\section*{Details}

An alternative to these functions, is to use table \((x, \operatorname{row}(x))\) and \(\operatorname{table}(x, \operatorname{col}(x))\), with the exception that the latter do not support the raw data type. When there are no missing values in \(x\), we have that all (rowTabulates \((x)==t(\operatorname{table}(x\), row \((x)))\) ) and all (colTabulates \((x)==\) \(t(\operatorname{table}(x, \operatorname{col}(x))))\). When there are missing values, we have that all(rowTabulates \((x)==\) \(t(\operatorname{table}(x, \operatorname{row}(x)\), useNA = "always") [, seq_len(nrow(x))])) and all(colTabulates(x) \(==t(\operatorname{table}(x, \operatorname{col}(x)\), useNA = "always")\([\), seq_len \((n c o l(x))])\) ).

\section*{Value}

Returns a NxJ (KxJ) matrix where \(\mathrm{N}(\mathrm{K})\) is the number of row (column) vectors tabulated and \(\mathbf{J}\) is the number of values counted.

\section*{Author(s)}

Henrik Bengtsson

\section*{Examples}
```

x <- matrix(1:5, nrow = 10, ncol = 5)
print(x)
print(rowTabulates(x))
print(colTabulates(x))

# Count only certain values

print(rowTabulates(x, values = 1:3))
y <- as.raw(x)
dim(y) <- dim(x)
print(y)
print(rowTabulates(y))
print(colTabulates(y))

```

\section*{rowVars Variance estimates for each row (column) in a matrix}

\section*{Description}

Variance estimates for each row (column) in a matrix.

\section*{Usage}
```

    rowVars(x, rows \(=\) NULL, cols \(=\) NULL, na. \(\mathrm{rm}=\) FALSE, refine \(=\) TRUE,
        center \(=\) NULL, \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\), useNames \(=\) TRUE)
    colVars(x, rows \(=\) NULL, cols \(=\) NULL, na.rm \(=\) FALSE, refine \(=\) TRUE,
        center \(=\) NULL, \(\operatorname{dim} .=\operatorname{dim}(x), \ldots\), useNames \(=\) TRUE)
    ```

\section*{Arguments}
\(\left.\begin{array}{ll}\mathrm{x} & \begin{array}{l}\text { An NxK matrix or, if dim. is specified, an } \mathrm{N} * \mathrm{~K} \text { vector. } \\ \text { rows } \\ \text { A vector indicating subset of rows to operate over. If NULL, no subsetting is } \\ \text { done. }\end{array} \\ \text { na.rm } & \begin{array}{l}\text { A vector indicating subset of columns to operate over. If NULL, no subsetting } \\ \text { is done. }\end{array} \\ \text { refine } & \begin{array}{l}\text { If TRUE, missing values are excluded. }\end{array} \\ \text { If TRUE, 'center' is NULL, and x is numeric, then extra effort is used to calculate } \\ \text { the average with greater numerical precision, otherwise not. } \\ \text { (optional; a vector or length N (K)) If the row (column) means are already esti- } \\ \text { mated, they can be pre-specified using this argument. This avoid re-estimating } \\ \text { them again. _Warning: It is important that a non-biased sample mean estimate } \\ \text { is passed. If not, then the variance estimate of the spread will also be biased._ If }\end{array}\right\}\)

\section*{Value}

Returns a numeric vector of length \(\mathrm{N}(\mathrm{K})\).

\section*{Providing center estimates}

The sample variance is estimated as
\(n /(n-1) * \operatorname{mean}\left((x-\text { center })^{2}\right)\),
where center is estimated as the sample mean, by default. In matrixStats \((<0.58 .0)\),
\(n /(n-1) *\left(\operatorname{mean}\left(x^{2}\right)-\right.\) center \(\left.^{2}\right)\)
was used. Both formulas give the same result _when_ 'center' is the sample mean estimate.
Argument 'center' can be used to provide an already existing estimate. It is important that the sample mean estimate is passed. If not, then the variance estimate of the spread will be biased.
For the time being, in order to lower the risk for such mistakes, argument 'center' is occasionally validated against the sample-mean estimate. If a discrepancy is detected, an informative error is provided to prevent incorrect variance estimates from being used. For performance reasons, this check is only performed once every 50 times. The frequency can be controlled by R option 'matrixStats.vars.formula.freq', whose default can be set by environment variable 'R_MATRIXSTATS_VARS_FORMULA_FREQ'.

\section*{Author(s)}

Henrik Bengtsson

\section*{See Also}

See rowMeans() and rowSums() in colSums().

\section*{Examples}
```

set.seed(1)
x <- matrix(rnorm(20), nrow = 5, ncol = 4)
print(x)

# Row averages

print(rowMeans(x))
print(rowMedians(x))

# Column averages

print(colMeans(x))
print(colMedians(x))

# Row variabilities

print(rowVars(x))
print(rowSds(x))
print(rowMads(x))
print(rowIQRs(x))

# Column variabilities

print(rowVars(x))
print(colSds(x))
print(colMads(x))
print(colIQRs(x))

```
```


# Row ranges

print(rowRanges(x))
print(cbind(rowMins(x), rowMaxs(x)))
print(cbind(rowOrderStats(x, which = 1), rowOrderStats(x, which = ncol(x))))

# Column ranges

print(colRanges(x))
print(cbind(colMins(x), colMaxs(x)))
print(cbind(colOrderStats(x, which = 1), colOrderStats(x, which = nrow(x))))
x <- matrix(rnorm(2000), nrow = 50, ncol = 40)

# Row standard deviations

d <- rowDiffs(x)
s1 <- rowSds(d) / sqrt(2)
s2 <- rowSds(x)
print(summary(s1 - s2))

# Column standard deviations

d <- colDiffs(x)
s1 <- colSds(d) / sqrt(2)
s2 <- colSds(x)
print(summary(s1 - s2))

```
rowWeightedMeans Calculates the weighted means for each row (column) in a matrix

\section*{Description}

Calculates the weighted means for each row (column) in a matrix.

\section*{Usage}
```

rowWeightedMeans(x, w = NULL, rows = NULL, cols = NULL, na.rm = FALSE,
..., useNames = TRUE)
colWeightedMeans(x, w = NULL, rows = NULL, cols = NULL, na.rm = FALSE,
..., useNames = TRUE)

```

\section*{Arguments}
x
w

\section*{rows}
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
\begin{tabular}{ll} 
na.rm & If TRUE, missing values are excluded. \\
\(\ldots\) & Not used. \\
useNames & If TRUE (default), names attributes of the result are set, otherwise not.
\end{tabular}

\section*{Details}

The implementations of these methods are optimized for both speed and memory. If no weights are given, the corresponding rowMeans()/colMeans() is used.

\section*{Value}

Returns a numeric vector of length \(\mathrm{N}(\mathrm{K})\).

\section*{Author(s)}

\section*{Henrik Bengtsson}

\section*{See Also}

See rowMeans() and colMeans() in colSums() for non-weighted means. See also weighted.mean.

\section*{Examples}
```

x <- matrix(rnorm(20), nrow = 5, ncol = 4)
print(x)

# Non-weighted row averages

mu_0 <- rowMeans(x)
mu <- rowWeightedMeans(x)
stopifnot(all.equal(mu, mu_0))

# Weighted row averages (uniform weights)

w <- rep(2.5, times = ncol(x))
mu <- rowWeightedMeans(x, w = w)
stopifnot(all.equal(mu, mu_0))

# Weighted row averages (excluding some columns)

w <- c(1, 1, 0, 1)
mu_0 <- rowMeans(x[, (w == 1), drop = FALSE])
mu <- rowWeightedMeans(x, w = w)
stopifnot(all.equal(mu, mu_0))

# Weighted row averages (excluding some columns)

w <- c(0, 1, 0, 0)
mu_0 <- rowMeans(x[, (w == 1), drop = FALSE])
mu <- rowWeightedMeans(x, w = w)
stopifnot(all.equal(mu, mu_0))

# Weighted averages by rows and columns

w <- 1:4
mu_1 <- rowWeightedMeans(x, w = w)

```
```

mu_2 <- colWeightedMeans(t(x), w = w)
stopifnot(all.equal(mu_2, mu_1))

```

\section*{Description}

Calculates the weighted medians for each row (column) in a matrix.

\section*{Usage}
```

rowWeightedMedians(x, w = NULL, rows = NULL, cols = NULL,
na.rm = FALSE, ..., useNames = TRUE)
colWeightedMedians(x, w = NULL, rows = NULL, cols = NULL,
na.rm = FALSE, ..., useNames = TRUE)

```

\section*{Arguments}

X
W
rows A vector indicating subset of rows to operate over. If NULL, no subsetting is done.
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
na.rm If TRUE, missing values are excluded.
... Additional arguments passed to weightedMedian().
useNames If TRUE (default), names attributes of the result are set, otherwise not.

\section*{Details}

The implementations of these methods are optimized for both speed and memory. If no weights are given, the corresponding rowMedians()/colMedians() is used.

\section*{Value}

Returns a numeric vector of length \(\mathrm{N}(\mathrm{K})\).

\section*{Author(s)}

\section*{Henrik Bengtsson}

\section*{See Also}

Internally, weightedMedian() is used. See rowMedians() and colMedians() for non-weighted medians.

\section*{Examples}
```

x <- matrix(rnorm(20), nrow = 5, ncol = 4)
print(x)

# Non-weighted row averages

mu_0 <- rowMedians(x)
mu <- rowWeightedMedians(x)
stopifnot(all.equal(mu, mu_0))

# Weighted row averages (uniform weights)

w <- rep(2.5, times = ncol(x))
mu <- rowWeightedMedians(x, w = w)
stopifnot(all.equal(mu, mu_0))

# Weighted row averages (excluding some columns)

w <- c(1, 1, 0, 1)
mu_0 <- rowMedians(x[, (w == 1), drop = FALSE])
mu <- rowWeightedMedians(x, w = w)
stopifnot(all.equal(mu, mu_0))

# Weighted row averages (excluding some columns)

w <- c(0, 1, 0, 0)
mu_0 <- rowMedians(x[, (w == 1), drop = FALSE])
mu <- rowWeightedMedians(x, w = w)
stopifnot(all.equal(mu, mu_0))

# Weighted averages by rows and columns

w <- 1:4
mu_1 <- rowWeightedMedians(x, w = w)
mu_2 <- colWeightedMedians(t(x), w = w)
stopifnot(all.equal(mu_2, mu_1))

```
varDiff

\section*{Description}

Estimation of scale based on sequential-order differences, corresponding to the scale estimates provided by var, sd, mad and IQR.

\section*{Usage}
varDiff(x, idxs \(=\) NULL, na.rm \(=\) FALSE, \(\operatorname{diff}=1 \mathrm{~L}, \operatorname{trim}=0, \ldots\) )
sdDiff(x, idxs \(=\) NULL, na.rm \(=\) FALSE, \(\operatorname{diff}=1 \mathrm{~L}, \operatorname{trim}=0, \ldots\) )
madDiff(x, idxs = NULL, na.rm = FALSE, diff = 1L, trim = 0, constant \(=1.4826, \ldots\) )
```

iqrDiff(x, idxs = NULL, na.rm = FALSE, diff = 1L, trim = 0, ...)
rowVarDiffs(x, rows = NULL, cols = NULL, na.rm = FALSE, diff = 1L,
trim $=0, \ldots$ useNames $=$ TRUE)
colVarDiffs $(x$, rows $=$ NULL, cols $=$ NULL, na.rm $=$ FALSE, diff $=1 \mathrm{~L}$,
trim $=0, \ldots$, useNames $=$ TRUE)
rowSdDiffs(x, rows = NULL, cols = NULL, na. $\mathrm{rm}=$ FALSE, $\operatorname{diff}=1 \mathrm{~L}$,
trim $=0, \ldots$ useNames $=$ TRUE)
colSdDiffs(x, rows $=$ NULL, cols $=$ NULL, na.rm $=$ FALSE, diff $=1 \mathrm{~L}$,
trim $=0, \ldots$ useNames $=$ TRUE)
rowMadDiffs(x, rows $=$ NULL, cols $=$ NULL, na.rm $=$ FALSE, $\operatorname{diff}=1 \mathrm{~L}$,
trim $=0, \ldots$, useNames $=$ TRUE)
colMadDiffs(x, rows $=$ NULL, cols $=$ NULL, na.rm $=$ FALSE, diff $=1 \mathrm{~L}$,
trim $=0, \ldots$, useNames $=$ TRUE)
rowIQRDiffs(x, rows $=$ NULL, cols $=$ NULL, na.rm $=$ FALSE, diff $=1 \mathrm{~L}$,
trim $=0, \ldots$, useNames $=$ TRUE)
colIQRDiffs(x, rows = NULL, cols = NULL, na.rm = FALSE, diff = 1L,
trim $=0, \ldots$ useNames $=$ TRUE)

```

\section*{Arguments}
\begin{tabular}{ll}
x \\
idxs & A numeric vector of length N or a numeric NxK matrix. \\
na.rm & \begin{tabular}{l} 
A vector indicating subset of elements to operate over. If NULL, no subsetting \\
is done.
\end{tabular} \\
diff & \begin{tabular}{l} 
If TRUE, missing values are excluded.
\end{tabular} \\
trim & \begin{tabular}{l} 
The positional distance of elements for which the difference should be calcu- \\
lated.
\end{tabular} \\
\(\ldots\) & \begin{tabular}{l} 
A double in \([0,1 / 2]\) specifying the fraction of observations to be trimmed from \\
each
\end{tabular} \\
constant & \begin{tabular}{l} 
A scale factor adjusting for asymptotically normal consistency.
\end{tabular} \\
rows & \begin{tabular}{l} 
A vector indicating subset of rows to operate over. If NULL, no subsetting is \\
done.
\end{tabular} \\
cols & \begin{tabular}{l} 
A vector indicating subset of columns to operate over. If NULL, no subsetting \\
is done.
\end{tabular} \\
useNames & \begin{tabular}{l} 
If TRUE (default), names attributes of the result are set, otherwise not.
\end{tabular}
\end{tabular}

\section*{Details}

Note that n-order difference MAD estimates, just like the ordinary MAD estimate by mad, apply a correction factor such that the estimates are consistent with the standard deviation under Gaussian distributions.
The interquartile range (IQR) estimates does not apply such a correction factor. If asymptotically normal consistency is wanted, the correction factor for IQR estimate is \(1 /(2 *\) qnorm(3/4)), which is half of that used for MAD estimates, which is \(1 /\) qnorm(3/4). This correction factor needs to be applied manually, i.e. there is no constant argument for the IQR functions.

\section*{Value}

Returns a numeric vector of length 1, length N , or length K.

\section*{Author(s)}

Henrik Bengtsson

\section*{References}
[1] J. von Neumann et al., The mean square successive difference. Annals of Mathematical Statistics, 1941, 12, 153-162.

\section*{See Also}

For the corresponding non-differentiated estimates, see var, sd, mad and IQR. Internally, diff2() is used which is a faster version of \(\operatorname{diff}()\).
```

weightedMad Weighted Median Absolute Deviation (MAD)

```

\section*{Description}

Computes a weighted MAD of a numeric vector.

\section*{Usage}
weightedMad(x, w = NULL, idxs = NULL, na.rm = FALSE, constant = 1.4826, center \(=\) NULL, ...)
rowWeightedMads(x, w = NULL, rows = NULL, cols = NULL, na.rm = FALSE, constant \(=1.4826\), center \(=\) NULL,... , useNames \(=\) TRUE)
colWeightedMads(x, w = NULL, rows = NULL, cols = NULL, na.rm = FALSE, constant \(=1.4826\), center \(=\) NULL,... , useNames \(=\) TRUE)

\section*{Arguments}

X
w
idxs A vector indicating subset of elements to operate over. If NULL, no subsetting is done.
na.rm If TRUE, missing values are excluded.
constant A numeric scale factor, cf. mad.
center Optional numeric scalar specifying the center location of the data. If NULL, it is estimated from data.
... Not used.
rows A vector indicating subset of rows to operate over. If NULL, no subsetting is done.
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
useNames If TRUE (default), names attributes of the result are set, otherwise not.

\section*{Value}

Returns a numeric scalar.

\section*{Missing values}

Missing values are dropped at the very beginning, if argument na. rm is TRUE, otherwise not.

\section*{Author(s)}

\section*{Henrik Bengtsson}

\section*{See Also}

For the non-weighted MAD, see mad. Internally weightedMedian() is used to calculate the weighted median.

\section*{Examples}
```

x <- 1:10
n <- length(x)
m1 <- mad(x)
m2 <- weightedMad(x)
stopifnot(identical(m1, m2))
w <- rep(1, times = n)
m1 <- weightedMad(x, w)
stopifnot(identical(m1, m2))

```
```

\# All weight on the first value
w[1] <- Inf
m <- weightedMad(x, w)
stopifnot (m == 0)
\# All weight on the first two values
w[1:2] <- Inf
$m 1<-\operatorname{mad}(x[1: 2])$
m2 <- weightedMad(x, w)
stopifnot(identical(m1, m2))
\# All weights set to zero
$\mathrm{w}<-\operatorname{rep}(0$, times $=\mathrm{n})$
m <- weightedMad(x, w)
stopifnot(is.na(m))

```
weightedMean
Weighted Arithmetic Mean

\section*{Description}

Computes the weighted sample mean of a numeric vector.

\section*{Usage}
```

weightedMean(x, w = NULL, idxs = NULL, na.rm = FALSE, refine = FALSE,

```
    ...)

\section*{Arguments}
x
w a vector of weights the same length as \(x\) giving the weights to use for each element of \(x\). Negative weights are treated as zero weights. Default value is equal weight to all values. If a missing-value weight exists, the result is always a missing value.
idxs A vector indicating subset of elements to operate over. If NULL, no subsetting is done.
na.rm If TRUE, missing values are excluded.
refine If TRUE and \(x\) is numeric, then extra effort is used to calculate the average with greater numerical precision, otherwise not.
... Not used.

\section*{Value}

Returns a numeric scalar. If \(x\) is of zero length, then NaN is returned, which is consistent with mean().

\section*{Missing values}

This function handles missing values consistently with weighted.mean. More precisely, if na.rm \(=\) FALSE, then any missing values in either \(x\) or \(w\) will give result NA_real_. If na. rm \(=\) TRUE, then all \((x, w)\) data points for which \(x\) is missing are skipped. Note that if both \(x\) and \(w\) are missing for a data points, then it is also skipped (by the same rule). However, if only w is missing, then the final results will always be NA_real_ regardless of na.rm.

\section*{Author(s)}

\section*{Henrik Bengtsson}

\section*{See Also}
mean() and weighted. mean.

\section*{Examples}
```

x <- 1:10
n <- length(x)
w <- rep(1, times = n)
m0 <- weighted.mean(x, w)
m1 <- weightedMean(x, w)
stopifnot(identical(m1, m0))

# Pull the mean towards zero

w[1] <- 5
m0 <- weighted.mean(x, w)
m1 <- weightedMean(x, w)
stopifnot(identical(m1, m0))

# Put even more weight on the zero

w[1] <- 8.5
m0 <- weighted.mean(x, w)
m1 <- weightedMean(x, w)
stopifnot(identical(m1, m0))

# All weight on the first value

w[1] <- Inf
m0 <- weighted.mean(x, w)
m1 <- weightedMean(x, w)
stopifnot(identical(m1, m0))

# All weight on the last value

w[1] <- 1
w[n] <- Inf
m0 <- weighted.mean(x, w)
m1 <- weightedMean(x, w)
stopifnot(identical(m1, m0))

# All weights set to zero

```
```

w <- rep(0, times = n)
m0 <- weighted.mean(x, w)
m1 <- weightedMean(x, w)
stopifnot(identical(m1, m0))

```
weightedMedian Weighted Median Value

\section*{Description}

Computes a weighted median of a numeric vector.

\section*{Usage}
weightedMedian(x, w = NULL, idxs = NULL, na.rm = FALSE, interpolate \(=\) is.null(ties), ties = NULL, ...)

\section*{Arguments}
x
w a vector of weights the same length as \(x\) giving the weights to use for each element of \(x\). Negative weights are treated as zero weights. Default value is equal weight to all values.
idxs A vector indicating subset of elements to operate over. If NULL, no subsetting is done.
na.rm a logical value indicating whether NA values in \(x\) should be stripped before the computation proceeds, or not. If NA, no check at all for NAs is done.
interpolate If TRUE, linear interpolation is used to get a consistent estimate of the weighted median.
ties If interpolate ==FALSE, a character string specifying how to solve ties between two x's that are satisfying the weighted median criteria. Note that at most two values can satisfy the criteria. When ties is "min" ("lower weighted median"), the smaller value of the two is returned and when it is "max" ("upper weighted median"), the larger value is returned. If ties is "mean", the mean of the two values is returned. Finally, if ties is "weighted" (or NULL) a weighted average of the two are returned, where the weights are weights of all values \(x[i]\) \(<=x[k]\) and \(x[i]>=x[k]\), respectively.
... Not used.

\section*{Value}

Returns a numeric scalar.
For the \(n\) elements \(x=c(x[1], x[2], \ldots, x[n])\) with positive weights \(w=c(w[1], w[2], \ldots\), \(w[n])\) such that \(\operatorname{sum}(w)=S\), the weighted median is defined as the element \(x[k]\) for which the total weight of all elements \(x[i]<x[k]\) is less or equal to \(S / 2\) and for which the total weight of all elements \(x[i]>x[k]\) is less or equal to \(S / 2\) (c.f. [1]).

When using linear interpolation, the weighted mean of \(x[k-1]\) and \(x[k]\) with weights \(S[k-1]\) and \(S[k]\) corresponding to the cumulative weights of those two elements is used as an estimate.
If \(w\) is missing then all elements of \(x\) are given the same positive weight. If all weights are zero, NA_real_ is returned.

If one or more weights are Inf, it is the same as these weights have the same weight and the others have zero. This makes things easier for cases where the weights are result of a division with zero.
If there are missing values in \(w\) that are part of the calculation (after subsetting and dropping missing values in \(x\) ), then the final result is always NA of the same type as \(x\).

The weighted median solves the following optimization problem:
\[
\alpha^{*}=\arg _{\alpha} \min \sum_{i=1}^{n} w_{i}\left|x_{i}-\alpha\right|
\]
where \(x=\left(x_{1}, x_{2}, \ldots, x_{n}\right)\) are scalars and \(w=\left(w_{1}, w_{2}, \ldots, w_{n}\right)\) are the corresponding "weights" for each individual \(x\) value.

\section*{Author(s)}

Henrik Bengtsson and Ola Hossjer, Centre for Mathematical Sciences, Lund University. Thanks to Roger Koenker, Econometrics, University of Illinois, for the initial ideas.

\section*{References}
[1] T.H. Cormen, C.E. Leiserson, R.L. Rivest, Introduction to Algorithms, The MIT Press, Massachusetts Institute of Technology, 1989.

\section*{See Also}
median, mean() and weightedMean().

\section*{Examples}
```

x <- 1:10
n <- length(x)
m1 <- median(x) \# 5.5
m2 <- weightedMedian(x) \# 5.5
stopifnot(identical(m1, m2))
w <- rep(1, times = n)
m1 <- weightedMedian(x, w) \# 5.5 (default)
m2 <- weightedMedian(x, ties = "weighted") \# 5.5 (default)
m3 <- weightedMedian(x, ties = "min") \# 5
m4 <- weightedMedian(x, ties = "max") \# 6
stopifnot(identical(m1, m2))

# Pull the median towards zero

w[1] <- 5
m1 <- weightedMedian(x, w) \# 3.5
y <- c(rep(0, times = w[1]), x[-1]) \# Only possible for integer weights

```
```

    m2 <- median(y) # 3.5
    stopifnot(identical(m1, m2))
    # Put even more weight on the zero
    w[1] <- 8.5
    weightedMedian(x, w) # 2
    # All weight on the first value
    w[1] <- Inf
    weightedMedian(x, w) # 1
    # All weight on the last value
    w[1] <- 1
    w[n] <- Inf
    weightedMedian(x, w) # 10
    # All weights set to zero
    w <- rep(0, times = n)
    weightedMedian(x, w) # NA
    # Simple benchmarking
    bench <- function(N = 1e5, K = 10) {
        x <- rnorm(N)
        gc()
        t <- c()
        t[1] <- system.time(for (k in 1:K) median(x))[3]
        t[2] <- system.time(for (k in 1:K) weightedMedian(x))[3]
        t <- t / t[1]
        names(t) <- c("median", "weightedMedian")
        t
    }
print(bench(N = 5, K = 100))
print(bench(N = 50, K = 100))
print(bench(N = 200, K = 100))
print(bench(N = 1000, K = 100))
print(bench(N = 10e3, K = 20))
print(bench(N = 100e3, K = 20))

```
weightedVar Weighted variance and weighted standard deviation

\section*{Description}

Computes a weighted variance / standard deviation of a numeric vector or across rows or columns of a matrix.

\section*{Usage}
weightedVar(x, w = NULL, idxs = NULL, na.rm = FALSE, center = NULL,
weightedVar
```

    ...)
    weightedSd(...)
rowWeightedVars(x, w = NULL, rows = NULL, cols = NULL, na.rm = FALSE,
..., useNames = TRUE)
colWeightedVars(x, w = NULL, rows = NULL, cols = NULL, na.rm = FALSE,
..., useNames = TRUE)
rowWeightedSds(x, w = NULL, rows = NULL, cols = NULL, na.rm = FALSE,
..., useNames = TRUE)
colWeightedSds(x, w = NULL, rows = NULL, cols = NULL, na.rm = FALSE,
..., useNames = TRUE)

```

\section*{Arguments}
\(x \quad\) vector of type integer, numeric, or logical.
w a vector of weights the same length as \(x\) giving the weights to use for each element of \(x\). Negative weights are treated as zero weights. Default value is equal weight to all values.
idxs A vector indicating subset of elements to operate over. If NULL, no subsetting is done.
na.rm If TRUE, missing values are excluded.
center Optional numeric scalar specifying the center location of the data. If NULL, it is estimated from data.
... Not used.
rows A vector indicating subset of rows to operate over. If NULL, no subsetting is done.
cols A vector indicating subset of columns to operate over. If NULL, no subsetting is done.
useNames If TRUE (default), names attributes of the result are set, otherwise not.

\section*{Details}

The estimator used here is the same as the one used by the "unbiased" estimator of the Hmisc package. More specifically, weightedVar \((x, w=w)==\) Hmisc: \(: w t d . \operatorname{var}(x\), weights \(=w)\),

\section*{Value}

Returns a numeric scalar.

\section*{Missing values}

This function handles missing values consistently with weightedMean(). More precisely, if na.rm \(=\) FALSE, then any missing values in either \(x\) or \(w\) will give result NA_real_. If na.rm \(=\) TRUE, then
all ( \(\mathrm{x}, \mathrm{w}\) ) data points for which x is missing are skipped. Note that if both x and w are missing for a data points, then it is also skipped (by the same rule). However, if only \(w\) is missing, then the final results will always be NA_real_ regardless of na.rm.

\section*{Author(s)}

Henrik Bengtsson

\section*{See Also}

For the non-weighted variance, see var.

\section*{Index}

\section*{* array}
product, 10
rowAlls, 11
rowCounts, 14
rowCumsums, 16
rowDiffs, 18
rowIQRs, 19
rowLogSumExps, 20
rowMads, 21
rowMeans2, 22
rowMedians, 23
rowOrderStats, 24
rowQuantiles, 26
rowRanges, 27
rowRanks, 28
rowSums2, 30
rowVars, 33
rowWeightedMeans, 35
rowWeightedMedians, 37
* iteration
anyMissing, 3
indexByRow, 7
product, 10
rowAlls, 11
rowCounts, 14
rowCumsums, 16
rowDiffs, 18
rowIQRs, 19
rowMads, 21
rowMeans2, 22
rowMedians, 23
rowOrderStats, 24
rowQuantiles, 26
rowRanges, 27
rowRanks, 28
rowSums2, 30
rowVars, 33
rowWeightedMeans, 35
rowWeightedMedians, 37
varDiff, 38
* logic
anyMissing, 3
indexByRow, 7
rowAlls, 11
rowCounts, 14
* package
matrixStats-package, 2
* robust
product, 10
rowDiffs, 18
rowIQRs, 19
rowMads, 21
rowMeans2, 22
rowMedians, 23
rowOrderStats, 24
rowQuantiles, 26
rowRanges, 27
rowRanks, 28
rowSums2, 30
rowVars, 33
rowWeightedMeans, 35
rowWeightedMedians, 37
varDiff, 38
weightedMad, 40
weightedMean, 42
weightedMedian, 44
weightedVar, 46
* univar
binCounts, 4
binMeans, 5
product, 10
rowAlls, 11
rowCounts, 14
rowCumsums, 16
rowDiffs, 18
rowIQRs, 19
rowMads, 21
rowMeans2, 22
rowMedians, 23
rowOrderStats, 24
rowQuantiles, 26
rowRanges, 27
rowRanks, 28
rowSums2, 30
rowVars, 33
rowWeightedMeans, 35
rowWeightedMedians, 37
varDiff, 38
weightedMad, 40
weightedMean, 42
weightedMedian, 44
weightedVar, 46
* utilities
rowCollapse, 13
rowTabulates, 31
[, 13
aggregate, 6
allValue (rowAlls), 11
anyMissing, 3
anyValue (rowAlls), 11
binCounts, 4, 6
binMeans, 5, 5
character, 10,29
colAlls (rowAlls), 11
colAnyMissings (anyMissing), 3
colAnyNAs (anyMissing), 3
colAnys (rowAlls), 11
colCollapse (rowCollapse), 13
colCounts (rowCounts), 14
colCummaxs (rowCumsums), 16
colCummins (rowCumsums), 16
colCumprods (rowCumsums), 16
colCumsums (rowCumsums), 16
colDiffs (rowDiffs), 18
colIQRDiffs (varDiff), 38
colIQRs (rowIQRs), 19
colLogSumExps (rowLogSumExps), 20
colMadDiffs (varDiff), 38
colMads (rowMads), 21
colMaxs (rowRanges), 27
colMeans2 (rowMeans2), 22
colMedians (rowMedians), 23
colMins (rowRanges), 27
colOrderStats (rowOrderStats), 24
colProds (product), 10
colQuantiles (rowQuantiles), 26
colRanges (rowRanges), 27
colRanks (rowRanks), 28
colSdDiffs (varDiff), 38
colSds (rowMads), 21
colSums, 25, 34, 36
colSums2 (rowSums2), 30
colTabulates (rowTabulates), 31
colVarDiffs (varDiff), 38
colVars (rowVars), 33
colWeightedMads (weightedMad), 40
colWeightedMeans (rowWeightedMeans), 35
colWeightedMedians
(rowWeightedMedians), 37
colWeightedSds (weightedVar), 46
colWeightedVars (weightedVar), 46
count (rowCounts), 14
cummax, 17
cummin, 17
cumprod, 17
cumsum, 17
data.frame, 3
diff, 40
diff2, 18,40
double, 15, 24, 25, 30, 39
FALSE, 4
hist, 5
indexByRow, 7
integer, 5-7, 11, 13, 15-18, 21-26, 28-33, 41, 44, 47
IQR, 19, 20, 38, 40
iqr (rowIQRs), 19
iqrDiff (varDiff), 38
is.nan, 10
list, 3
logical, 5, 12, 17, 26, 29, 32, 41, 44, 47
logSumExp, 8, 21
mad, 22, 38, 40, 41
madDiff (varDiff), 38
matrix, 3, 10, 11, 13-26, 28-33, 35, 37, 39, 42
matrixStats (matrixStats-package), 2
matrixStats-package, 2
mean, \(6,42,43,45\)
median, 45
NA, 10, 19, 24, 44
NA_real_, 45
names, 30
\(\mathrm{NaN}, 10\)
NULL, 3-5, 7, 8, 10, 11, 13-16, 18-20, 22-26,
\(28,29,31-33,35,37,39,41,42,44\), 47
numeric, 4-8, 10, 17-26, 28, 30, 31, 33, 35-37, 39-42, 44, 47
pmin.int, 28
prod, 10
product, 10, 10
quantile, 26
rank, 29, 30
raw, 32
rowAlls, 11
rowAnyMissings (anyMissing), 3
rowAnyNAs (anyMissing), 3
rowAnys (rowAlls), 11
rowCollapse, 13
rowCounts, 14
rowCummaxs (rowCumsums), 16
rowCummins (rowCumsums), 16
rowCumprods (rowCumsums), 16
rowCumsums, 16
rowDiffs, 18
rowIQRDiffs (varDiff), 38
rowIQRs, 19, 22
rowLogSumExps, 9, 20
rowMadDiffs (varDiff), 38
rowMads, 21
rowMaxs (rowRanges), 27
rowMeans, 24
rowMeans2, 22, 24
rowMedians, 23, 37
rowMins (rowRanges), 27
rowOrderStats, 24, 28
rowProds (product), 10
rowQuantiles, 19, 26
rowRanges, 27
rowRanks, 28
rowSdDiffs (varDiff), 38
rowSds, 20
rowSds (rowMads), 21
rowSums2, 30
rowTabulates, 31
rowVarDiffs (varDiff), 38
rowVars, 33
rowWeightedMads (weightedMad), 40
rowWeightedMeans, 35
rowWeightedMedians, 24, 37
rowWeightedSds (weightedVar), 46
rowWeightedVars (weightedVar), 46
sd, 22, 38, 40
sdDiff(varDiff), 38
tabulate, 5
TRUE, 3-6, 8, 10, 11, 13-15, 17-26, 28, 29, 31-33, 36, 37, 39, 41, 42, 44, 47
var, 22, 38, 40, 48
varDiff, 38
vector, \(3-8,10-16,18-26,28,29,31-33\), 35-37, 39-42, 44, 47
weighted.mean, 36, 43
weightedMad, 40
weightedMean, \(42,45,47\)
weightedMedian, \(37,41,44\)
weightedSd (weightedVar), 46
weightedVar, 46```

