

# Package ‘EnrichedHeatmap’

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**Type** Package

**Title** Making Enriched Heatmaps

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**Author** Zuguang Gu

**Maintainer** Zuguang Gu <z.gu@dkfz.de>

**Depends** R (>= 3.1.2), methods, grid, ComplexHeatmap (>= 1.99.0),  
GenomicRanges

**Imports** matrixStats, stats, GetoptLong, Rcpp, utils, locfit, circlize  
(>= 0.4.5), IRanges

**Suggests** testthat (>= 0.3), knitr, markdown, genefilter, RColorBrewer

**VignetteBuilder** knitr

**Description** Enriched heatmap is a special type of heatmap which visualizes the enrichment of genomic signals on specific target regions. Here we implement enriched heatmap by ComplexHeatmap package. Since this type of heatmap is just a normal heatmap but with some special settings, with the functionality of ComplexHeatmap, it would be much easier to customize the heatmap as well as concatenating to a list of heatmaps to show correspondance between different data sources.

**biocViews** Software, Visualization, Sequencing, GenomeAnnotation,  
Coverage

**URL** <https://github.com/jokergoo/EnrichedHeatmap>

**License** MIT + file LICENSE

**LinkingTo** Rcpp

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## R topics documented:

|                                  |           |
|----------------------------------|-----------|
| anno_enriched . . . . .          | 2         |
| as.normalizedMatrix . . . . .    | 3         |
| copyAttr . . . . .               | 5         |
| default_smooth_fun . . . . .     | 5         |
| discretize . . . . .             | 6         |
| dist_by_closeness . . . . .      | 7         |
| EnrichedHeatmap . . . . .        | 8         |
| enriched_score . . . . .         | 10        |
| extract_anno_enriched . . . . .  | 11        |
| failed_rows . . . . .            | 11        |
| getSignalsFromList . . . . .     | 12        |
| makeWindows . . . . .            | 13        |
| normalizeToMatrix . . . . .      | 14        |
| print.normalizedMatrix . . . . . | 17        |
| rbind.normalizedMatrix . . . . . | 17        |
| [.normalizedMatrix . . . . .     | 18        |
| <b>Index</b>                     | <b>19</b> |

---

|               |   |
|---------------|---|
| anno_enriched | <i>Annotation Function to Show the Enrichment</i> |
|---------------|---|

---

### Description

Annotation Function to Show the Enrichment

### Usage

```
anno_enriched(gp = gpar(col = "red"), pos_line = NULL, pos_line_gp = NULL,
  ylim = NULL, value = c("mean", "sum", "abs_mean", "abs_sum"),
  yaxis = TRUE, axis = yaxis, axis_param = list(side = "right"),
  show_error = FALSE, height = unit(2, "cm"), ...)
```

### Arguments

|             |   |
|-------------|---|
| gp          | Graphic parameters. There are two non-standard parameters: neg_col and pos_col. If these two parameters are defined, the positive signals and negative signals are visualized separately. The graphic parameters can be set as vectors when the heatmap or heatmap list is split into several row clusters. |
| pos_line    | Whether draw vertical lines which represent positions of target?  |
| pos_line_gp | Graphic parameters for the position lines.  |
| ylim        | Ranges on y-axis. By default it is inferred from the data.  |
| value       | The method to summarize signals from columns of the normalized matrix.  |
| yaxis       | Deprecated, use axis instead.   |
| axis        | Whether show axis?  |
| axis_param  | parameters for controlling axis. See <a href="#">default_axis_param</a> for all possible settings and default parameters.   |

|            |   |
|------------|---|
| show_error | Whether show error regions which are one standard error to the mean value? Color of error area is same as the corresponding lines with 75 percent transparency. |
| height     | Height of the annotation.   |
| ...        | Other arguments.  |

### Details

This annotation functions shows mean values (or depends on the method set in value argument) of columns in the normalized matrix which summarises the enrichment of the signals to the targets.

If rows are splitted, the enriched lines are calculated for each row cluster and there will also be multiple lines in this annotation viewport.

It should only be placed as column annotation of the enriched heatmap.

### Value

A column annotation function which should be set to top\_annotation argument in [EnrichedHeatmap](#).

### Author(s)

Zuguang Gu <z.gu@dkfz.de>

### Examples

```
load(system.file("extdata", "chr21_test_data.RData", package = "EnrichedHeatmap"))
tss = promoters(genes, upstream = 0, downstream = 1)
mat1 = normalizeToMatrix(H3K4me3, tss, value_column = "coverage",
  extend = 5000, mean_mode = "w0", w = 50, keep = c(0, 0.99))
EnrichedHeatmap(mat1, col = c("white", "red"), name = "H3K4me3",
  top_annotation = HeatmapAnnotation(lines = anno_enriched(gp = gpar(col = 2:4))),
  km = 3, row_title_rot = 0)
```

---

as.normalizedMatrix     *Convert a Normal Matrix to a normalizedMatrix Object*

---

### Description

Convert a Normal Matrix to a normalizedMatrix Object

### Usage

```
as.normalizedMatrix(mat, k_upstream = 0, k_downstream = 0, k_target = 0,
  extend, signal_name = "signals", target_name = "targets",
  background = NA, smooth = FALSE, smooth_fun = default_smooth_fun,
  keep = c(0, 1), trim = NULL)
```

**Arguments**

|              |   |
|--------------|---|
| mat          | A matrix generated by other software.   |
| k_upstream   | Number of windows in the upstream.  |
| k_downstream | Number of windows in the downstream.  |
| k_target     | Number of windows in the target.  |
| extend       | Extension to the target. The length should be 1 (if one of k_upstream or k_downstream is zero). or 2 (if both of k_upstream and k_downstream are non-zero).   |
| signal_name  | The name of signal regions. It is only used for printing the object.  |
| target_name  | The name of the target names. It is only used for printing the object.  |
| background   | The background value in the matrix.   |
| smooth       | Whether apply smoothing on rows in the matrix.  |
| smooth_fun   | The smoothing function that is applied to each row in the matrix. This self-defined function accepts a numeric vector (may contain NA values) and returns a vector with same length. If the smoothing is failed, the function should call <a href="#">stop</a> to throw errors so that <a href="#">normalizeToMatrix</a> can catch how many rows are failed in smoothing. See the default <a href="#">default_smooth_fun</a> for example. |
| keep         | Percentiles in the normalized matrix to keep. The value is a vector of two percent values. Values less than the first percentile is replaces with the first percentile and values larger than the second percentile is replaced with the second percentile.   |
| trim         | Deprecated, please use keep instead.  |

**Details**

If users use the matrix from other software, they can use this function to convert it to the `normalizedMatrix` object and visualize it afterwards.

**Value**

A `normalizedMatrix` object.

**Author(s)**

`z.gu@dkfz.de`

**Examples**

```
# There is no example
NULL
```

---

`copyAttr`*Copy Attributes to Another Object*

---

**Description**

Copy Attributes to Another Object

**Usage**

```
copyAttr(x, y)
```

**Arguments**

|   |           |
|---|-----------|
| x | Object 1. |
| y | Object 2. |

**Details**

The `normalizeToMatrix` object is actually a matrix but with more additional attributes attached. When manipulating such matrix, there are some circumstances that the attributes are lost. This function is used to copy these specific attributes when dealing with the matrix.

**Author(s)**

Zuguang Gu <z.gu@dkfz.de>

**Examples**

```
gr = GRanges(seqnames = c("chr5", "chr5"),
  ranges = IRanges(start = c(98, 98),
    end = c(104, 104)))
target = GRanges(seqnames = "chr5",
  ranges = IRanges(start = 100,
    end = 100))
mat1 = normalizeToMatrix(gr, target, extend = 6, w = 1)
# attributes removed and you cannot use it for EnrichedHeatmap()
mat2 = mat1[]
# copy attributes to mat2 and now mat3 can be used for EnrichedHeatmap()
mat3 = copyAttr(mat1, mat2)
```

---

`default_smooth_fun`*Default Smoothing function*

---

**Description**

Default Smoothing function

**Usage**

```
default_smooth_fun(x)
```

**Arguments**

x                    Input numeric vector.

**Details**

The smoothing function is applied to every row in the normalized matrix. For this default smoothing function, `locfit` is first tried on the vector. If there is error, `loess` smoothing is tried afterwards. If both smoothing are failed, there will be an error.

**Author(s)**

Zuguang Gu <z.gu@dkfz.de>

**Examples**

```
# There is no example
NULL
```

---

discretize

*Discretize a Continuous Matrix to a Discrete Matrix*

---

**Description**

Discretize a Continuous Matrix to a Discrete Matrix

**Usage**

```
discretize(mat, rule, right_closed = FALSE)
```

**Arguments**

mat                    A normalized matrix from `normalizeToMatrix`.

rule                    A list of intervals which provide mapping between continuous values to discrete values. Note the order of intervals determines the order of corresponding discrete levels.

right\_closed          Is the interval right closed?

**Details**

Assuming we have a normalized matrix with both positive values and negative values, we only want to see the enrichment of the windows/regions showing significant positive values and negative values and we are only interested in the direction of the values while not the value itself, then we can define the rule as:

```
rule = list(
  "positive" = c(0.5, Inf),
  "negative" = c(-Inf, -0.5)
)
```

And we can convert the continuous matrix to a discrete matrix and visualize it:

```
mat2 = discretize(mat, rule)
EnrichedHeatmap(mat2, col = c("positive" = "red", "negative" = "green"))
```

Another example is to discretize the signals to discrete levels according to the intensities:

```
rule = list(
  "very_high" = c(100, Inf),
  "high" = c(50, 100),
  "intermediate" = c(25, 50),
  "low" = c(1e-6, 25)
)
```

### Author(s)

Zuguang Gu <z.gu@dkfz.de>

### Examples

```
# There is no example
NULL
```

---

|                   |                              |
|-------------------|------------------------------|
| dist_by_closeness | <i>Distance by Closeness</i> |
|-------------------|------------------------------|

---

### Description

Distance by Closeness

### Usage

```
dist_by_closeness(mat)
```

### Arguments

mat                    A numeric matrix where the distance is calculated by rows.

### Details

For two rows in the matrix, assume  $x_1, x_2, \dots, x_{n1}$  are the column index of non-zero values in row 1 and  $y_1, y_2, \dots, y_{n2}$  are the column index for non-zero values in row 2, the distance between the two rows based on the closeness is calculated as:

$$d_{\text{closeness}} = \frac{\sum_i \sum_j (|x_i - y_j|)}{(n_1 * n_2)}$$

### Value

A `dist` object.

### Author(s)

Zuguang Gu <z.gu@dkfz.de>

**Examples**

```
x1 = c(0, 0, 0, 0, 1, 1, 1, 0, 0, 0)
x2 = c(0, 0, 0, 1, 1, 1, 0, 0, 0, 0)
x3 = c(1, 0, 0, 0, 1, 1, 0, 0, 0, 0)
m = rbind(x1, x2, x3)
dist(m)
dist_by_closeness(m)
```

---

EnrichedHeatmap

*Constructor Method for the Enriched Heatmap*


---

**Description**

Constructor Method for the Enriched Heatmap

**Usage**

```
EnrichedHeatmap(mat,
  col,
  top_annotation = HeatmapAnnotation(enriched = anno_enriched()),
  row_order = order(enriched_score(mat), decreasing = TRUE),
  pos_line = TRUE,
  pos_line_gp = gpar(lty = 2),
  axis_name = NULL,
  axis_name_rot = 0,
  axis_name_gp = gpar(fontsize = 10),
  border = TRUE,
  cluster_rows = FALSE,
  row_dend_reorder = -enriched_score(mat),
  show_row_dend = FALSE,
  show_row_names = FALSE,
  heatmap_legend_param = list(),
  ...)
```

**Arguments**

|                |  |
|----------------|--|
| mat            | A matrix which is returned by <a href="#">normalizeToMatrix</a> .  |
| col            | Color settings. If the signals are categorical, color should be a vector with category levels as names.  |
| top_annotation | A special annotation which is always put on top of the enriched heatmap and is constructed by <a href="#">anno_enriched</a> .  |
| row_order      | Row order. Default rows are ordered by enriched scores calculated from <a href="#">enriched_score</a> .  |
| pos_line       | Whether draw vertical lines which represent the positions of target?   |
| pos_line_gp    | Graphic parameters for the position lines.   |
| axis_name      | Names for axis which is below the heatmap. If the targets are single points, axis_name is a vector of length three which corresponds to upstream, target itself and downstream. If the targets are regions with width larger than 1, axis_name should be a vector of length four which corresponds to upstream, start of targets, end of targets and downstream. |



|                                   |  |
|-----------------------------------|--|
| <code>axis_name_rot</code>        | Rotation for axis names.   |
| <code>axis_name_gp</code>         | Graphic parameters for axis names.   |
| <code>border</code>               | Whether show the border of the heatmap?  |
| <code>cluster_rows</code>         | Clustering on rows are turned off by default.  |
| <code>show_row_dend</code>        | Whether show dendrograms on rows if hierarchical clustering is applied on rows?                      |
| <code>row_dend_reorder</code>     | Weight for reordering the row dendrogram. It is reordered by enriched scores by default.             |
| <code>show_row_names</code>       | Whether show row names?  |
| <code>heatmap_legend_param</code> | A list of settings for heatmap legends. <code>at</code> and <code>labels</code> can not be set here. |
| <code>...</code>                  | Other arguments passed to <a href="#">Heatmap</a> .  |

## Details

The enriched heatmap is essentially a normal heatmap but with several special settings. Following parameters are set with pre-defined values:

```
cluster_columns enforced to be FALSE
show_column_names enforced to be FALSE
bottom_annotation enforced to be NULL
```

[EnrichedHeatmap](#) calls [Heatmap](#), thus, most of the arguments in [Heatmap](#) are usable in [EnrichedHeatmap](#) such as to apply clustering on rows, or to split rows by a data frame or k-means clustering. Users can also add more than one heatmaps by `+` operator. Enriched heatmaps and normal heatmaps can be concatenated mixed.

For detailed demonstration, please go to the vignette.

## Value

A [Heatmap-class](#) object.

## Author(s)

Zuguang Gu <z.gu@dkfz.de>

## Examples

```
load(system.file("extdata", "chr21_test_data.RData", package = "EnrichedHeatmap"))
mat3 = normalizeToMatrix(meth, cgi, value_column = "meth", mean_mode = "absolute",
  extend = 5000, w = 50, smooth = TRUE)
EnrichedHeatmap(mat3, name = "methylation", column_title = "methylation near CGI")
EnrichedHeatmap(mat3, name = "meth1") + EnrichedHeatmap(mat3, name = "meth2")
# for more examples, please go to the vignette
```

---

|                |                        |
|----------------|------------------------|
| enriched_score | <i>Enriched Scores</i> |
|----------------|------------------------|

---

### Description

Enriched Scores

### Usage

```
enriched_score(mat)
```

### Arguments

mat                    A normalized matrix from [normalizeToMatrix](#).

### Details

The function calculates how the signal is enriched in the target by weighting the distance to the target.

For a numeric vector, assume the vector is denoted as combination of three sub-vectors  $c(x_1, x_2, x_3)$  with length  $n_1, n_2$  and  $n_3$ , where  $x_1$  are data points in upstream windows,  $x_2$  are data points in target windows and  $x_3$  are data points in downstream windows, the enriched score is calculated as

$$\text{sum}(x_{1i} * i/n_1) + \text{sum}(x_{3j} * (n_3 - j + 1)/n_3) + \text{sum}(x_{2k} * \text{abs}(n_2/2 - \text{abs}(k - n_2/2)))$$

where the first two terms are the distance to the start or end position of the target by weighting the distance to the position that if it is closer to the start or end position of the target, it has higher weight. The second term weight the distance to the center point of the target and similar, if it is closer to the center position, it has higher weight.

### Value

A numeric vector.

### Author(s)

Zuguang Gu <z.gu@dkfz.de>

### Examples

```
# There is no example  
NULL
```

---

extract\_anno\_enriched *Extarct Enrichment Annotation Graphics as a Separated Plot*

---

### Description

Extarct Enrichment Annotation Graphics as a Separated Plot

### Usage

```
extract_anno_enriched(ht_list, which = NULL, newpage = TRUE, padding = NULL)
```

### Arguments

|         |  |
|---------|--|
| ht_list | The heatmap list returned by <a href="#">draw, HeatmapList-method</a> .  |
| which   | The index of enriched heatmap in the heatmap list. The value can be an integer index or a character index (the name of the heatmap). |
| newpage | Whether call <a href="#">grid.newpage</a> to create a new page?  |
| padding | Padding of the plot.   |

### Details

The extracted plot is exactly the same as that on the enriched heatmap.

### Author(s)

Zuguang Gu <z.gu@dkfz.de>

### Examples

```
# There is no example
NULL
```

---

failed\_rows *Indices of Rows Failed from Smoothing*

---

### Description

Indices of Rows Failed from Smoothing

### Usage

```
failed_rows(m)
```

### Arguments

|   |   |
|---|---|
| m | Matrix from <a href="#">normalizeToMatrix</a> . |
|---|---|

**Value**

A numeric vector or NULL.

**Examples**

```
# There is no example
NULL
```

---

getSignalsFromList      *Get Signals from a List*

---

**Description**

Get Signals from a List

**Usage**

```
getSignalsFromList(lt, fun = function(x) mean(x, na.rm = TRUE))
```

**Arguments**

|     |   |
|-----|---|
| lt  | A list of normalized matrices which are returned by <a href="#">normalizeToMatrix</a> . Matrices in the list should be generated with same settings (e.g. they should use same target regions, same extension to targets and same number of windows). |
| fun | A user-defined function to summarize signals.   |

**Details**

Let's assume you have a list of histone modification signals for different samples and you want to visualize the mean pattern across samples. You can first normalize histone mark signals for each sample and then calculate means values across all samples. In following example code, `hm_gr_list` is a list of GRanges objects which contain positions of histone modifications, `tss` is a GRanges object containing positions of gene TSS.

```
mat_list = NULL
for(i in seq_along(hm_gr_list)) {
  mat_list[[i]] = normalizeToMatrix(hm_gr_list[[i]], tss, value_column = "density")
}
```

If we compress the list of matrices as a three-dimension array where the first dimension corresponds to genes, the second dimension corresponds to windows and the third dimension corresponds to samples, the mean signal across all sample can be calculated on the third dimension. Here [getSignalsFromList](#) simplifies this job.

Applying `getSignalsFromList()` to `mat_list`, it gives a new normalized matrix which contains mean signals across all samples and can be directly used in `EnrichedHeatmap()`.

```
mat_mean = getSignalsFromList(mat_list)
EnrichedHeatmap(mat_mean)
```

The correlation between histone modification and gene expression can also be calculated on the third dimension of the array. In the user-defined function fun, x is the vector for gene i and window j in the array, and i is the index of current gene.

```
mat_corr = getSignalsFromList(mat_list,
  fun = function(x, i) cor(x, expr[i, ], method = "spearman"))
```

Then mat\_corr here can be used to visualize how gene expression is correlated to histone modification around TSS.

```
EnrichedHeatmap(mat_corr)
```

### Value

A [normalizeToMatrix](#) object which can be directly used for [EnrichedHeatmap](#).

### Author(s)

Zuguang Gu <z.gu@dkfz.de>

### Examples

```
NULL
```

---

makeWindows

*Split Regions into Windows*

---

### Description

Split Regions into Windows

### Usage

```
makeWindows(query, w = NULL, k = NULL, direction = c("normal", "reverse"),
  short.keep = FALSE)
```

### Arguments

|            |   |
|------------|---|
| query      | A <a href="#">GRanges-class</a> object.   |
| w          | Window size. A value larger than 1 means the number of base pairs and a value between 0 and 1 is the percent to the current region.   |
| k          | Number of partitions for each region. If it is set, all other arguments are ignored.  |
| direction  | Where to start the splitting? See 'Details' section.  |
| short.keep | If the the region can not be split equally under the window size, the argument controls whether to keep the windows that are smaller than the window size. See 'Details' section. |

**Details**

Following illustrates the meaning of `direction` and `short.keep`:

```
-->-->-->- one region, split by 3bp window (">" represents the direction of the sequence)
aaabbbccc  direction = "normal",  short.keep = FALSE
aaabbbcccd direction = "normal",  short.keep = TRUE
  aaabbbccc direction = "reverse", short.keep = FALSE
  abbbccddd direction = "reverse", short.keep = TRUE
```

**Value**

A `GRanges-class` object with two additional columns attached:

- `.i_query` which contains the correspondance between small windows and original regions in query
- `.i_window` which contains the index of the small window on the current region.

**Author(s)**

Zuguang gu <z.gu@dkfz.de>

**Examples**

```
query = GRanges(seqnames = "chr1", ranges = IRanges(start = c(1, 11, 21), end = c(10, 20, 30)))
makeWindows(query, w = 2)
makeWindows(query, w = 0.5)
makeWindows(query, w = 3)
makeWindows(query, w = 3, direction = "reverse")
makeWindows(query, w = 3, short.keep = TRUE)
makeWindows(query, w = 3, direction = "reverse", short.keep = TRUE)
makeWindows(query, w = 12)
makeWindows(query, w = 12, short.keep = TRUE)
makeWindows(query, k = 2)
makeWindows(query, k = 3)
query = GRanges(seqnames = "chr1", ranges = IRanges(start = c(1, 11, 31), end = c(10, 30, 70)))
makeWindows(query, w = 2)
makeWindows(query, w = 0.2)
```

---

|                   |  |
|-------------------|--|
| normalizeToMatrix | <i>Normalize Associations between Genomic Signals and Target Regions into a Matrix</i> |
|-------------------|--|

---

**Description**

Normalize Associations between Genomic Signals and Target Regions into a Matrix

**Usage**

```
normalizeToMatrix(signal, target, extend = 5000, w = max(extend)/50,
  value_column = NULL, mapping_column = NULL, background = ifelse(smooth, NA, 0), empty_value = NULL,
  mean_mode = c("absolute", "weighted", "w0", "coverage"), include_target = any(width(target) > 1),
  target_ratio = min(c(0.4, mean(width(target))/(sum(extend) + mean(width(target))))),
  k = min(c(20, min(width(target))))), smooth = FALSE, smooth_fun = default_smooth_fun,
  keep = c(0, 1), trim = NULL, flip_upstream = FALSE)
```

**Arguments**

|                |  |
|----------------|--|
| signal         | A <a href="#">GRanges-class</a> object.  |
| target         | A <a href="#">GRanges-class</a> object.  |
| extend         | Extended base pairs to the upstream and/or downstream of target. It can be a vector of length one or two. Length one means same extension to the upstream and downstream.  |
| w              | Window size for splitting upstream and downstream, measured in base pairs  |
| value_column   | Column index in signal that is mapped to colors. If it is not set, it assumes values for all signal regions are 1.   |
| mapping_column | Mapping column to restrict overlapping between signal and target. By default it tries to look for all regions in signal that overlap with every target.  |
| background     | Values for windows that don't overlap with signal.   |
| empty_value    | Deprecated, please use background instead.   |
| mean_mode      | When a window is not perfectly overlapped to signal, how to summarize values to the window. See 'Details' section for a detailed explanation.  |
| include_target | Whether include target in the heatmap? If the width of all regions in target is 1, include_target is enforced to FALSE.  |
| target_ratio   | The ratio of target columns in the normalized matrix. If the value is 1, extend will be reset to 0.  |
| k              | Number of windows only when target_ratio = 1 or extend == 0, otherwise ignored.  |
| smooth         | Whether apply smoothing on rows in the matrix?   |
| smooth_fun     | The smoothing function that is applied to each row in the matrix. This self-defined function accepts a numeric vector (may contain NA values) and returns a vector with same length. If the smoothing is failed, the function should call <a href="#">stop</a> to throw errors so that <a href="#">normalizeToMatrix</a> can catch how many rows are failed in smoothing. See the default <a href="#">default_smooth_fun</a> for example.  |
| keep           | Percentiles in the normalized matrix to keep. The value is a vector of two percent values. Values less than the first percentile is replaces with the first percentile and values larger than the second percentile is replaced with the second percentile.  |
| trim           | Deprecated, please use keep instead.   |
| flip_upstream  | Sometimes whether the signals are on the upstream or the downstream of the targets are not important and users only want to show the relative distance to targets. If the value is set to TRUE, the upstream part in the normalized matrix is flipped and added to the downstream part The flipping is only allowed when the targets are single-point targets or the targets are excluded in the normalized matrix (by setting include_target = FALSE). If the extension for the upstream and downstream is not equal, the smaller extension is used for the final matrix. |

**Details**

In order to visualize associations between signal and target, the data is transformed into a matrix and visualized as a heatmap by [EnrichedHeatmap](#) afterwards.

Upstream and downstream also with the target body are splitted into a list of small windows and overlap to signal. Since regions in signal and small windows do not always 100 percent overlap, there are four different averaging modes:

Following illustrates different settings for mean\_mode (note there is one signal region overlapping with other signals):

```

    40      50      20      values in signal regions
+++++++  +++   +++++  signal regions
          30      values in signal region
          +++++  signal region
=====
          4 6 3      3      a window (17bp), there are 4bp not overlapping to any signal regions.
                                overlap

```

absolute:  $(40 + 30 + 50 + 20)/4$

weighted:  $(40*4 + 30*6 + 50*3 + 20*3)/(4 + 6 + 3 + 3)$

w0:  $(40*4 + 30*6 + 50*3 + 20*3)/(4 + 6 + 3 + 3 + 4)$

coverage:  $(40*4 + 30*6 + 50*3 + 20*3)/17$

### Value

A matrix with following additional attributes:

upstream\_index column index corresponding to upstream of target

target\_index column index corresponding to target

downstream\_index column index corresponding to downstream of target

extend extension on upstream and downstream

smooth whether smoothing was applied on the matrix

failed\_rows index of rows which are failed after smoothing

The matrix is wrapped into a simple normalizeToMatrix class.

### Author(s)

Zuguang Gu <z.gu@dkfz.de>

### Examples

```

signal = GRanges(seqnames = "chr1",
  ranges = IRanges(start = c(1, 4, 7, 11, 14, 17, 21, 24, 27),
    end = c(2, 5, 8, 12, 15, 18, 22, 25, 28)),
  score = c(1, 2, 3, 1, 2, 3, 1, 2, 3))
target = GRanges(seqnames = "chr1", ranges = IRanges(start = 10, end = 20))
normalizeToMatrix(signal, target, extend = 10, w = 2)
normalizeToMatrix(signal, target, extend = 10, w = 2, include_target = TRUE)
normalizeToMatrix(signal, target, extend = 10, w = 2, value_column = "score")

```



---

```
print.normalizedMatrix
```

*Print the Normalized Matrix*

---

**Description**

Print the Normalized Matrix

**Usage**

```
## S3 method for class 'normalizedMatrix'  
print(x, ...)
```

**Arguments**

|     |   |
|-----|---|
| x   | The normalized matrix returned by <a href="#">normalizeToMatrix</a> . |
| ... | Other arguments.  |

**Value**

No value is returned.

**Author(s)**

Zuguang Gu <z.gu@dkfz.de>

**Examples**

```
# There is no example  
NULL
```

---

```
rbind.normalizedMatrix
```

*Bind Matrix by Rows*

---

**Description**

Bind Matrix by Rows

**Usage**

```
## S3 method for class 'normalizedMatrix'  
rbind(..., deparse.level = 1)
```

**Arguments**

|               |           |
|---------------|-----------|
| ...           | Matrices  |
| deparse.level | Not used. |

**Value**

A normalizedMatrix class object.

**Author(s)**

z.gu@dkfz.de

**Examples**

```
# There is no example
NULL
```

---

[.normalizedMatrix     *Subset normalized matrix by rows*

---

**Description**

Subset normalized matrix by rows

**Usage**

```
## S3 method for class 'normalizedMatrix'
x[i, j, drop = FALSE]
```

**Arguments**

|      |   |
|------|---|
| x    | the normalized matrix returned by <a href="#">normalizeToMatrix</a> |
| i    | row index   |
| j    | column index  |
| drop | whether drop the dimension  |

**Value**

A normalizedMatrix class object.

**Author(s)**

Zuguang Gu <z.gu@dkfz.de>

**Examples**

```
# There is no example
NULL
```

# Index

[.normalizedMatrix, 18

anno\_enriched, 2, 8

as.normalizedMatrix, 3

copyAttr, 5

default\_axis\_param, 2

default\_smooth\_fun, 4, 5, 15

discretize, 6

dist, 7

dist\_by\_closeness, 7

enriched\_score, 8, 10

EnrichedHeatmap, 3, 8, 9, 13, 15

Extract.normalizedMatrix  
([.normalizedMatrix), 18

extract\_anno\_enriched, 11

failed\_rows, 11

getSignalsFromList, 12, 12

grid.newpage, 11

Heatmap, 9

locfit, 6

loess, 6

makeWindows, 13

normalizeToMatrix, 4–6, 8, 10–13, 14, 15,  
17, 18

print.normalizedMatrix, 17

rbind.normalizedMatrix, 17

stop, 4, 15