

Package ‘POMA’

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Title Tools for Omics Data Analysis

Version 1.14.0

Description The POMA package offers a comprehensive toolkit designed for omics data analysis, streamlining the process from initial visualization to final statistical analysis. Its primary goal is to simplify and unify the various steps involved in omics data processing, making it more accessible and manageable within a single, intuitive R package. Emphasizing on reproducibility and user-friendliness, POMA leverages the standardized SummarizedExperiment class from Bioconductor, ensuring seamless integration and compatibility with a wide array of Bioconductor tools. This approach guarantees maximum flexibility and replicability, making POMA an essential asset for researchers handling omics datasets. See <https://github.com/pcastellanoescuder/POMAShiny>. Paper: Castellano-Escuder et al. (2021) <[doi:10.1371/journal.pcbi.1009148](https://doi.org/10.1371/journal.pcbi.1009148)> for more details.

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box_cox_transformation
Box-Cox Transformation

Description

Compute Box-Cox normalization.

Usage

`box_cox_transformation(data)`

Arguments

`data` A single variable.

cor_pmat *Correlation P-Values*

Description

Compute correlation p-values.

Usage

`cor_pmat(x, method)`

Arguments

`x` A data matrix.
`method` Character indicating which correlation coefficient has to be computed. Options are "pearson" (default), "kendall" and "spearman".

| | |
|-----------------|------------------------|
| detect_decimals | <i>Detect decimals</i> |
|-----------------|------------------------|

Description

Detect decimal variables.

Usage

```
detect_decimals(data)
```

Arguments

| | |
|------|----------------------------------|
| data | A data matrix (samples in rows). |
|------|----------------------------------|

| | |
|-------------------|-----------------------------------|
| flattenCorrMatrix | <i>Flatten Correlation Matrix</i> |
|-------------------|-----------------------------------|

Description

Flatten Correlation Matrix

Usage

```
flattenCorrMatrix(cormat, pmat)
```

Arguments

| | |
|--------|-----------------------|
| cormat | Output from cor. |
| pmat | Output from cor_pmat. |

| | |
|-----------|-------------------------|
| PomaBatch | <i>Batch Correction</i> |
|-----------|-------------------------|

Description

PomaBatch performs batch correction on a SummarizedExperiment object given a batch factor variable.

Usage

```
PomaBatch(data, batch, mod = NULL)
```

Arguments

| | |
|-------|---|
| data | A SummarizedExperiment object. |
| batch | Character. The name of the column in colData that contains the batch information. |
| mod | Character vector. Indicates the names of colData columns to be included as covariates. Default is NULL (no covariates). |

Value

A SummarizedExperiment object with batch-corrected data.

Author(s)

Pol Castellano-Escuder

References

Leek JT, Johnson WE, Parker HS, Fertig EJ, Jaffe AE, Zhang Y, Storey JD, Torres LC (2023). sva: Surrogate Variable Analysis. doi:10.18129/B9.bioc.sva <https://doi.org/10.18129/B9.bioc.sva>

Examples

```
data("st000284")

st000284 %>%
  PomaImpute(method = "knn") %>%
  PomaBatch(batch = "gender")
```

PomaBoxplots

Boxplots and Violin Plots

Description

PomaBoxplots generates boxplots and violin plots for samples and features. This function can be used for data exploration (e.g., comparison between pre and post normalized datasets).

Usage

```
PomaBoxplots(
  data,
  x = "samples",
  violin = FALSE,
  feature_name = NULL,
  theme_params = list(legend_title = FALSE, axis_x_rotate = TRUE)
)
```

Arguments

| | |
|--------------|---|
| data | A SummarizedExperiment object. |
| x | Character. Options are "samples" (to visualize sample boxplots) and "features" (to visualize feature boxplots). Default is "samples". |
| violin | Logical. Indicates if violin plots should be displayed instead of boxplots. Default is FALSE. |
| feature_name | Character vector. Indicates the feature/s to display. Default is NULL (all features will be displayed). |
| theme_params | List. Indicates theme_poma parameters. |

Value

A ggplot object.

Author(s)

Pol Castellano-Escuder

Examples

```

data("st000284")

# Sample boxplots
st000284 %>%
PomaNorm() %>%
PomaBoxplots(theme_params = list(axistext = "y"))

# Sample violin plots
st000284 %>%
PomaNorm() %>%
PomaBoxplots(violin = TRUE, theme_params = list(axistext = "y"))

# All feature boxplots
st000284 %>%
PomaNorm() %>%
PomaBoxplots(x = "features", theme_params = list(axis_x_rotate = TRUE))

# Specific feature boxplots
st000284 %>%
PomaNorm() %>%
PomaBoxplots(x = "features",
              feature_name = c("ornithine", "orotate"))

# Specific feature violin plots
st000284 %>%
PomaNorm() %>%
PomaBoxplots(x = "features",
              violin = TRUE,
              feature_name = c("ornithine", "orotate"))

```

Description

PomaClust performs a k-means clustering and plots the results in a classical multidimensional scaling (MDS) plot.

Usage

```
PomaClust(  
  data,  
  method = "euclidean",  
  k = NA,  
  k_max = floor(min(dim(data))/2),  
  show_clusters = TRUE,  
  labels = FALSE  
)
```

Arguments

| | |
|---------------|--|
| data | A SummarizedExperiment object. |
| method | Character. Indicates the distance method to perform MDS. Options are "euclidean", "maximum", "manhattan", "canberra" and "minkowski". See ?dist(). |
| k | Numeric. Indicates the number of clusters (default is NA). The optimal number of clusters will be used by default. |
| k_max | Numeric. Indicates the number of clusters among which the optimal k will be selected. |
| show_clusters | Logical. Indicates if clusters should be plotted or not. |
| labels | Logical. Indicates if sample names should be plotted or not. |

Value

A list with results including plots and tables.

Author(s)

Pol Castellano-Escuder

Examples

```
data("st000284")  
  
PomaClust(st000284)
```

PomaCorr

Correlation Analysis

Description

PomaCorr computes all pairwise correlations in the data.

Usage

```
PomaCorr(data, method = "pearson", label_size = 8, theme_params = list())
```

Arguments

| | |
|--------------|---|
| data | A SummarizedExperiment object. |
| method | Character. Indicates which correlation coefficient has to be computed. Options are "pearson" (default), "kendall" and "spearman". |
| label_size | Numeric. Indicates plot label size. |
| theme_params | List. Indicating theme_poma parameters. |

Value

A list with the results.

Author(s)

Pol Castellano-Escuder

References

Jerome Friedman, Trevor Hastie and Rob Tibshirani (2019). *glasso: Graphical Lasso: Estimation of Gaussian Graphical Models*. R package version 1.11. <https://CRAN.R-project.org/package=glasso>

Examples

```
data("st000284")

# Pearson correlation
PomaCorr(st000284)$correlations

## Gaussian graphical model
# library(ggraph)
# PomaCorr(st000284, corr_type = "glasso")
```

PomaCreateObject *Create a SummarizedExperiment Object*

Description

PomaCreateObject creates a SummarizedExperiment object from data frames.

Usage

```
PomaCreateObject(metadata = NULL, features = NULL, factor_levels = 10)
```

Arguments

| | |
|---------------|--|
| metadata | Metadata variables structured in columns. Sample ID must be the first column. |
| features | Matrix of features. Each feature is a column. |
| factor_levels | Numeric. Integer variables with more levels than indicated by this parameter will be treated as factors. |

Value

A SummarizedExperiment object.

Author(s)

Pol Castellano-Escuder

References

Morgan M, Obenchain V, Hester J, Pagès H (2021). SummarizedExperiment: SummarizedExperiment container. R package version 1.24.0, <https://bioconductor.org/packages/SummarizedExperiment>.

Examples

```
data(iris)

# Create metadata: Data frame with sample names and a group factor
metadata <- data.frame(ID = 1:150, Group = iris$Species)

# Create features: `p` column data frame with features
features <- iris[, 1:4]

# Create a `SummarizedExperiment` object with `POMA`
object <- PomaCreateObject(metadata = metadata, features = features)
```

PomaDensity

*Density Plots***Description**

PomaDensity generates a density plot for samples and features. This function can be used for data exploration (e.g., comparison between pre and post normalized datasets).

Usage

```
PomaDensity(
  data,
  x = "samples",
  feature_name = NULL,
  theme_params = list(legend_title = FALSE)
)
```

Arguments

| | |
|--------------|---|
| data | A SummarizedExperiment object. |
| x | Character. Options are "samples" (to visualize sample density plots) and "features" (to visualize feature density plots). Default is "samples". |
| feature_name | Character vector. Indicates the feature/s to display. Default is NULL (all features will be displayed). |
| theme_params | List. Indicates theme_poma parameters. |

Value

A ggplot object.

Author(s)

Pol Castellano-Escuder

Examples

```
data("st000284")

# Sample density plots
st000284 %>%
PomaNorm() %>%
PomaDensity(theme_params = list(axistext = "y"))

# All feature density plots
st000284 %>%
PomaNorm() %>%
PomaDensity(x = "features", theme_params = list(legend_position = "none"))
```

```
# Specific feature density plots
st000284 %>%
PomaNorm() %>%
PomaDensity(x = "features",
            feature_name = c("ornithine", "orotate"))
```

| | |
|-----------|---|
| PomaDESeq | <i>Differential Expression Analysis Based on the Negative Binomial Distribution</i> |
|-----------|---|

Description

PomaDESeq estimates variance-mean dependence in count data from high-throughput sequencing assays and test for differential expression based on a model using the negative binomial distribution.

Usage

```
PomaDESeq(data, adjust = "fdr")
```

Arguments

| | |
|--------|---|
| data | A SummarizedExperiment object. |
| adjust | Character. Multiple comparisons correction method to adjust p-values. Available options are: "fdr" (false discovery rate), "holm", "hochberg", "hommel", "bonferroni", "BH" (Benjamini-Hochberg), and "BY" (Benjamini-Yekutieli). |

Value

A tibble with the results.

Author(s)

Pol Castellano-Escuder

References

Love, M.I., Huber, W., Anders, S. Moderated estimation of fold change and dispersion for RNA-seq data with DESeq2 Genome Biology 15(12):550 (2014)

PomaHeatmap

Heatmap Plot

Description

PomaHeatmap generates a heatmap.

Usage

```
PomaHeatmap(  
  data,  
  covs = NULL,  
  sample_names = TRUE,  
  feature_names = FALSE,  
  show_legend = TRUE  
)
```

Arguments

| | |
|---------------|---|
| data | A SummarizedExperiment object. |
| covs | Character vector. Indicates the names of colData columns to be included as covariates. Default is NULL (no covariates). |
| sample_names | Logical. Indicates if sample names should be displayed or not. Default is TRUE. |
| feature_names | Logical. Indicates if feature names should be displayed or not. Default is FALSE. |
| show_legend | Logical. Indicates if legend should be displayed or not. Default is TRUE. |

Value

A heatmap plot.

Author(s)

Pol Castellano-Escuder

Examples

```
data("st000284")  
  
# Basic heatmap  
st000284 %>%  
  PomaNorm() %>%  
  PomaHeatmap()  
  
# Heatmap with one covariate  
st000284 %>%  
  PomaNorm() %>%
```

```

PomaHeatmap(covs = "factors")

# Heatmap with two covariates
st000284 %>%
  PomaNorm() %>%
  PomaHeatmap(covs = c("factors", "smoking_condition"))

```

PomaImpute

Impute Missing Values

Description

PomaImpute performs missing value imputation on a dataset using various imputation methods.

Usage

```

PomaImpute(
  data,
  zeros_as_na = FALSE,
  remove_na = TRUE,
  cutoff = 20,
  group_by = TRUE,
  method = "knn"
)

```

Arguments

| | |
|-------------|--|
| data | A SummarizedExperiment object. |
| zeros_as_na | Logical. Indicates if the zeros in the data are missing values. Default is FALSE. |
| remove_na | Logical. Indicates if features with a percentage of missing values over the cutoff parameter should be removed. Default is TRUE. |
| cutoff | Numeric. Percentage of missing values allowed in each feature. |
| group_by | Logical. If metadata file is present and its first variable is a factor, it can be used to compute missing values per group and drop them accordingly. Features will be removed only if all of the groups contain more missing values than allowed. Default is TRUE. |
| method | Character. The imputation method to use. Options include "none" (no imputation, replace missing values by zeros), "half_min" (replace missing values with half of the minimum value), "median" (replace missing values with the median), "mean" (replace missing values with the mean), "min" (replace missing values with the minimum value), "knn" (replace missing values using k-nearest neighbors imputation), and "random_forest" (replace missing values using random forest imputation). |

Value

A SummarizedExperiment object without missing values.

Author(s)

Pol Castellano-Escuder

References

Armitage, E. G., Godzien, J., Alonso-Herranz, V., López-González, Á., & Barbas, C. (2015). Missing value imputation strategies for metabolomics data. *Electrophoresis*, 36(24), 3050-3060.

Examples

```
data("st000336")

PomaImpute(st000336, method = "knn")
```

| | |
|-----------|---|
| PomaLasso | <i>Lasso, Ridge, and Elasticnet Regularized Generalized Linear Models for Binary Outcomes</i> |
|-----------|---|

Description

PomaLasso performs LASSO, Ridge, and Elasticnet regression for feature selection and prediction purposes for binary outcomes.

Usage

```
PomaLasso(
  data,
  alpha = 1,
  ntest = NULL,
  nfolds = 10,
  lambda = NULL,
  labels = FALSE
)
```

Arguments

| | |
|--------|---|
| data | A SummarizedExperiment object. |
| alpha | Numeric. Indicates the elasticnet mixing parameter. alpha = 1 is the LASSO penalty and alpha = 0 is the Ridge penalty. |
| ntest | Numeric. Indicates the percentage of observations that will be used as test set. Default is NULL (no test set). |
| nfolds | Numeric. Indicates number of folds for cross-validation (default is 10). Although nfolds can be as large as the sample size (leave-one-out CV), it is not recommended for large datasets. Smallest value allowable is nfolds = 3. |
| lambda | Numeric. Indicates the user supplied lambda sequence. Typical usage is to have the program compute its own lambda sequence based on nlambda and lambda.min.ratio. See ?glmnet::glmnet(). |

labels Logical. Indicates if feature names should be plotted in coefficient plot or not. Default is FALSE.

Value

A list with results.

Author(s)

Pol Castellano-Escuder

References

Jerome Friedman, Trevor Hastie, Robert Tibshirani (2010). Regularization Paths for Generalized Linear Models via Coordinate Descent. *Journal of Statistical Software*, 33(1), 1-22. URL <http://www.jstatsoft.org/v33/i01/>.

Examples

```
data("st000336")

# lasso
st000336 %>%
  PomaImpute() %>%
  PomaNorm() %>%
  PomaLasso()

# elasticnet
st000336 %>%
  PomaImpute() %>%
  PomaNorm() %>%
  PomaLasso(alpha = 0.5)

# ridge
st000336 %>%
  PomaImpute() %>%
  PomaNorm() %>%
  PomaLasso(alpha = 0)
```

Description

PomaLimma uses the classical limma package to compute differential expression analysis.

Usage

```
PomaLimma(data, contrast = NULL, covs = NULL, adjust = "fdr", weights = FALSE)
```

Arguments

| | |
|----------|---|
| data | A SummarizedExperiment object. |
| contrast | Character. Indicates the comparison. For example, "Group1-Group2" or "control-intervention". |
| covs | Character vector. Indicates the names of colData columns to be included as covariates. Default is NULL (no covariates). If not NULL, a limma model will be fitted using the specified covariates. Note: The order of the covariates is important and should be listed in increasing order of importance in the experimental design. |
| adjust | Character. Indicates the multiple comparisons correction method. Options are: "fdr", "holm", "hochberg", "hommel", "bonferroni", "BH" and "BY". |
| weights | Logical. Indicates whether the limma model should estimate the relative quality weights for each group. See ?limma::arrayWeights(). |

Value

A tibble with the results.

Author(s)

Pol Castellano-Escuder

References

Matthew E. Ritchie, Belinda Phipson, Di Wu, Yifang Hu, Charity W. Law, Wei Shi, Gordon K. Smyth, limma powers differential expression analyses for RNA-sequencing and microarray studies, *Nucleic Acids Research*, Volume 43, Issue 7, 20 April 2015, Page e47, <https://doi.org/10.1093/nar/gkv007>

Examples

```
data("st000284")

st000284 %>%
  PomaNorm() %>%
  PomaLimma(contrast = "Healthy-CRC", adjust = "fdr")
```

PomaLM

Linear Models

Description

PomaLM performs a linear model on a SummarizedExperiment object.

Usage

```
PomaLM(data, x = NULL, y = NULL, adjust = "fdr")
```


Arguments

| | |
|---------------------|---|
| <code>data</code> | A SummarizedExperiment object. |
| <code>x</code> | Character vector. Indicates the names of independent variables. If it's NULL (default), all features will be used. |
| <code>y</code> | Character. Indicates the name of colData numeric columns to be used as dependent variable. If it's set to NULL, the first numeric variable in colData will be used as the dependent variable. |
| <code>adjust</code> | Character. Multiple comparisons correction method to adjust p-values. Available options are: "fdr" (false discovery rate), "holm", "hochberg", "hommel", "bonferroni", "BH" (Benjamini-Hochberg), and "BY" (Benjamini-Yekutieli). |

Value

A list with results including plots and tables.

Author(s)

Pol Castellano-Escuder

Examples

```
data("st000284")

# Perform linear model with all features
st000284 %>%
PomaLM()

# Perform linear model with two features
st000284 %>%
PomaLM(x = c("x1_methyladenosine", "x2_deoxyuridine"))
```

Description

PomaLMM performs linear mixed models on a SummarizedExperiment object.

Usage

```
PomaLMM(data, x = NULL, y = NULL, adjust = "fdr", clean_plot = FALSE)
```

Arguments

| | |
|-------------------------|---|
| <code>data</code> | A SummarizedExperiment object. |
| <code>x</code> | Character vector. Indicates the names of <code>colData</code> columns to be used as random and fixed effects (independent variables). If it's set to <code>NULL</code> (default), all variables in <code>colData</code> will be used. |
| <code>y</code> | Character vector. Indicates the names of dependent variables. If it's <code>NULL</code> (default), all features will be used. |
| <code>adjust</code> | Character. Multiple comparisons correction method to adjust p-values. Available options are: "fdr" (false discovery rate), "holm", "hochberg", "hommel", "bonferroni", "BH" (Benjamini-Hochberg), and "BY" (Benjamini-Yekutieli). |
| <code>clean_plot</code> | Logical. Indicates if remove intercept and linear mixed model residues boxplots from the plot. Default is <code>FALSE</code> . |

Value

A list with results including plots and tables. Table values indicate the percentage variance explained per variable.

Author(s)

Pol Castellano-Escuder

Examples

```
data("st000284")

# Perform linear mixed model with all features
st000284 %>%
PomaLMM()

# Perform linear mixed model with two features
st000284 %>%
PomaLMM(y = c("x1_methyladenosine", "x1_methylhistamine"))

# Perform linear mixed model with one random effect
st000284 %>%
PomaLMM(x = "smoking_condition")

# Perform linear mixed model with two random effects and two features
st000284 %>%
PomaLMM(x = c("smoking_condition", "gender"),
        y = c("x1_methyladenosine", "x1_methylhistamine"))

# Perform linear mixed model with no random effects and two features, therefore, a linear model will be fitted
st000284 %>%
PomaLMM(x = "age_at_consent", # Numerical, i.e., fixed effect
        y = c("x1_methyladenosine", "x1_methylhistamine"))

# Perform linear mixed model with no random effects and all features, therefore, a linear model will be fitted
st000284 %>%
```

```
PomaLMM(x = "age_at_consent") # Numerical i.e., fixed effect
```

PomaNorm

Normalize Data

Description

PomaNorm performs data normalization using various normalization methods.

Usage

```
PomaNorm(data, sample_norm = "none", method = "log_pareto")
```

Arguments

| | |
|-------------|--|
| data | A SummarizedExperiment object. |
| sample_norm | Character. Sample normalization method. Options include "none" (default), "sum", or "quantile". |
| method | Character. The normalization method to use. Options include "none" (no normalization), "auto_scaling" (autoscaling normalization, i.e., Z-score normalization), "level_scaling" (level scaling normalization), "log_scaling" (log scaling normalization), "log_transform" (log transformation normalization), "vast_scaling" (vast scaling normalization), "log_pareto" (log Pareto scaling normalization), "min_max" (min-max normalization), and "box_cox" (Box-Cox transformation). |

Value

A SummarizedExperiment object with normalized data.

Author(s)

Pol Castellano-Escuder

References

Van den Berg, R. A., Hoefsloot, H. C., Westerhuis, J. A., Smilde, A. K., & van der Werf, M. J. (2006). Centering, scaling, and transformations: improving the biological information content of metabolomics data. *BMC genomics*, 7(1), 142.

Examples

```
data("st000284")  
  
PomaNorm(st000284, method = "log_pareto")
```

PomaOddsRatio *Logistic Regression Model Odds Ratios*

Description

PomaOddsRatio calculates the Odds Ratios for each feature from a logistic regression model using the binary outcome (group/type must be a binary factor) as a dependent variable.

Usage

```
PomaOddsRatio(data, feature_name = NULL, covs = NULL, show_ci = TRUE)
```

Arguments

| | |
|--------------|--|
| data | A SummarizedExperiment object. |
| feature_name | Character vector. Indicates the name/s of feature/s that will be used to fit the model. If it's NULL (default), all variables will be included in the model. |
| covs | Character vector. Indicates the names of colData columns to be included as covariates. Default is NULL (no covariates). |
| show_ci | Logical. Indicates if the 95% confidence intervals will be plotted. Default is TRUE. |

Value

A list with results including plots and tables.

Author(s)

Pol Castellano-Escuder

Examples

```
data("st000336")

st000336 %>%
  PomaImpute() %>%
  PomaNorm() %>%
  PomaOddsRatio(feature_name = c("glutamic_acid", "glutamine", "glycine", "histidine"))
```

Description

PomaOutliers analyses and removes statistical outliers from the data.

Usage

```
PomaOutliers(  
  data,  
  method = "euclidean",  
  type = "median",  
  coef = 2,  
  labels = FALSE  
)
```

Arguments

| | |
|--------|---|
| data | A SummarizedExperiment object. |
| method | Character. Indicates the distance measure method to perform MDS. |
| type | Character. Indicates the type of outlier analysis to perform. Options are "median" (default) and "centroid". See <code>vegan::betadisper</code> . |
| coef | Numeric. Indicates the outlier coefficient. Lower values are more sensitive to outliers while higher values are less restrictive about outliers. |
| labels | Logical. Indicates if sample names should to be plotted. |

Value

A list with the results.

Author(s)

Pol Castellano-Escuder

Examples

```
data("st000336")  
  
# clean outliers  
st000336 %>%  
  PomaImpute() %>%  
  PomaNorm() %>%  
  PomaOutliers()
```

PomaPCA

Principal Components Analysis

Description

PomaPCA performs a principal components analysis on the given SummarizedExperiment object.

Usage

```
PomaPCA(  
  data,  
  center = TRUE,  
  scale = TRUE,  
  ncomp = 4,  
  labels = FALSE,  
  ellipse = FALSE,  
  load_length = 1  
)
```

Arguments

| | |
|-------------|---|
| data | A SummarizedExperiment object. |
| center | Logical. Indicates whether the variables should be shifted to be zero centered. Default is TRUE. |
| scale | Logical. Indicates whether the variables should be scaled to have unit variance before the analysis takes place. Default is TRUE. |
| ncomp | Numeric. Number of components to be included in the results. Default is 4. |
| labels | Logical. Indicates if sample names should be displayed. |
| ellipse | Logical. Indicates whether a 95 percent confidence interval ellipse should be displayed in score plot and biplot. Default is FALSE. |
| load_length | Numeric. Indicates the length of biplot loading arrows. Value between 1 and 2. Default is 1. |

Value

A list with results including plots and tables.

Author(s)

Pol Castellano-Escuder

Examples

```
data("st000336")

st000336 %>%
  PomaImpute() %>%
  PomaNorm() %>%
  PomaPCA()
```

| | |
|---------|--|
| PomaPCR | <i>Principal Components Regression</i> |
|---------|--|

Description

PomaPCR performs Principal Components Regression.

Usage

```
PomaPCR(data, center = TRUE, scale = TRUE, ncomp = 2, y = NULL, adjust = "fdr")
```

Arguments

| | |
|--------|---|
| data | A SummarizedExperiment object. |
| center | Logical. Indicates whether the variables should be shifted to be zero centered. Default is TRUE. |
| scale | Logical. Indicates whether the variables should be scaled to have unit variance before the analysis takes place. Default is TRUE. |
| ncomp | Numeric. Indicates the number of principal components used as predictors in the model. Default is 2. |
| y | Character. Indicates the name of colData columns to be used as dependent variable. If it's set to NULL, the first numeric variable in colData will be used as the dependent variable. |
| adjust | Character. Multiple comparisons correction method to adjust p-values. Available options are: "fdr" (false discovery rate), "holm", "hochberg", "hommel", "bonferroni", "BH" (Benjamini-Hochberg), and "BY" (Benjamini-Yekutieli). |

Value

A tibble with the results.

Author(s)

Pol Castellano-Escuder

Examples

```
data("st000284")

# PCR with 2 components
st000284 %>%
  PomaPCR(y = "age_at_consent")

# PCR with 20 components
st000284 %>%
  PomaPCR(ncomp = 20)
```

PomaPLS

Partial Least Squares Methods

Description

PomaPLS performs Partial Least Squares (PLS) regression, Partial Least Squares Discriminant Analysis (PLS-DA) to classify samples, and Sparse Partial Least Squares Discriminant Analysis (sPLS-DA) to classify samples (supervised analysis) and select variables.

Usage

```
PomaPLS(  
  data,  
  method = "pls",  
  y = NULL,  
  ncomp = 5,  
  labels = FALSE,  
  ellipse = TRUE,  
  cross_validation = FALSE,  
  validation = "Mfold",  
  folds = 5,  
  nrepeat = 10,  
  vip = 1,  
  num_features = 10,  
  theme_params = list()  
)
```

Arguments

| | |
|---------------------|---|
| <code>data</code> | A SummarizedExperiment object. |
| <code>method</code> | Character. PLS method. Options include "pls", "plsda", and "splsda". |
| <code>y</code> | Character. Indicates the name of colData columns to be used as dependent variable. If it's set to NULL, the first variable in colData will be used as the dependent variable. |
| <code>ncomp</code> | Numeric. Number of components in the model. Default is 5. |

| | |
|------------------|---|
| labels | Logical. Indicates if sample names should be displayed. |
| ellipse | Logical. Indicates whether a 95 percent confidence interval ellipse should be displayed. Default is TRUE. |
| cross_validation | Logical. Indicates if cross-validation should be performed for PLS-DA ("plsda") and sPLS-DA ("splstda") methods. Default is FALSE. |
| validation | Character. (Only for "plsda" and "splstda" methods). Indicates the cross-validation method. Options are "Mfold" and "loo" (Leave-One-Out). |
| folds | Numeric. (Only for "plsda" and "splstda" methods). Number of folds for "Mfold" cross-validation method (default is 5). If the validation method is "loo", this value is set to 1. |
| nrepeat | Numeric. (Only for "plsda" and "splstda" methods). Number of times the cross-validation process is repeated. |
| vip | Numeric. (Only for "plsda" method). Indicates the variable importance in the projection (VIP) cutoff. |
| num_features | Numeric. (Only for "splstda" method). Number of features to discriminate groups. |
| theme_params | List. Indicates theme_poma parameters. |

Value

A list with results including plots and tables.

Author(s)

Pol Castellano-Escuder

Examples

```
data("st000284")

# PLS
st000284 %>%
  PomaNorm() %>%
  PomaPLS(method = "pls")

data("st000336")

# PLSDA
st000336 %>%
  PomaImpute() %>%
  PomaNorm() %>%
  PomaPLS(method = "plsda")

# PLSDA with Cross-Validation
st000336 %>%
  PomaImpute() %>%
  PomaNorm() %>%
```

```

PomaPLS(method = "plsda", cross_validation = TRUE)

# sPLSDA
st000336 %>%
  PomaImpute() %>%
  PomaNorm() %>%
  PomaPLS(method = "splsda")

# sPLSDA with Cross-Validation
st000336 %>%
  PomaImpute() %>%
  PomaNorm() %>%
  PomaPLS(method = "splsda", ncomp = 3, cross_validation = TRUE)

```

PomaRandForest

Classification Random Forest

Description

PomaRandForest performs classification random forest. This method can be used both for prediction and variable selection.

Usage

```

PomaRandForest(
  data,
  ntest = NULL,
  ntree = 500,
  mtry = floor(sqrt(ncol(t(SummarizedExperiment::assay(data))))),
  nodesize = 1,
  nvar = 20
)

```

Arguments

| | |
|----------|--|
| data | A SummarizedExperiment object. |
| ntest | Numeric. Indicates the percentage of observations that will be used as test set. Default is NULL (no test set). |
| ntree | Numeric. Indicates the number of trees to grow. |
| mtry | Numeric. Indicates the number of variables randomly sampled as candidates at each split. This value is set \sqrt{p} (where p is number of variables in data) by default. |
| nodesize | Numeric. Indicates the minimum size of terminal nodes. Default is 1. |
| nvar | Numeric. Indicates the number of variables to show in the Gini Index plot. |

Value

A list with results including plots and tables.

Author(s)

Pol Castellano-Escuder

References

A. Liaw and M. Wiener (2002). Classification and Regression by randomForest. R News 2(3), 18–22.

Examples

```
data("st000336")

st000336 %>%
  PomaImpute() %>%
  PomaRandForest()
```

PomaRankProd

Rank Product/Rank Sum Analysis

Description

PomaRankProd performs the Rank Product (or Rank Sum) method to identify differentially expressed genes.

Usage

```
PomaRankProd(data, logged = TRUE, paired = NA, cutoff = 0.05, method = "pfp")
```

Arguments

| | |
|--------|---|
| data | A SummarizedExperiment object. |
| logged | Logical. Indicates if data should be log transformed first. |
| paired | Numeric. Indicates the number of random pairs generated in the function, if set to NA (default), the odd integer closer to the square of the number of replicates is used. |
| cutoff | Numeric. Indicates the pfp/pvalue threshold value used to select features. |
| method | Character. Indicates the method to identify features. "pfp" uses percentage of false prediction, which is a default setting. "pval" uses p-values which is less stringent than pfp. |

Value

A list with results including plots and tables.

Author(s)

Pol Castellano-Escuder

References

Breitling, R., Armengaud, P., Amtmann, A., and Herzyk, P.(2004) Rank Products: A simple, yet powerful, new method to detect differentially regulated genes in replicated microarray experiments, FEBS Letter, 57383-92

Hong, F., Breitling, R., McEntee, W.C., Wittner, B.S., Nemhauser, J.L., Chory, J. (2006). RankProd: a bioconductor package for detecting differentially expressed genes in meta-analysis Bioinformatics. 22(22):2825-2827

Del Carratore, F., Jankevics, A., Eisinga, R., Heskes, T., Hong, F. & Breitling, R. (2017). RankProd 2.0: a refactored Bioconductor package for detecting differentially expressed features in molecular profiling datasets. Bioinformatics. 33(17):2774-2775

Examples

```
data("st000336")

st000336 %>%
  PomaImpute() %>%
  PomaRankProd()
```

PomaUMAP

Dimensionality Reduction with UMAP

Description

PomaUMAP performs a dimension reduction of the data using the Uniform Manifold Approximation and Projection (UMAP) method. See `?uwot::umap()` for more.

Usage

```
PomaUMAP(
  data,
  n_neighbors = floor(sqrt(nrow(data))),
  n_components = 2,
  metric = "euclidean",
  pca = NULL,
  min_dist = 0.01,
  spread = 1,
  hdbSCAN_minpts = floor(nrow(data) * 0.05),
  show_clusters = TRUE,
  hide_noise = TRUE,
  labels = FALSE,
  theme_params = list(legend_title = TRUE, legend_position = "bottom")
)
```

Arguments

| | |
|-----------------------------|---|
| <code>data</code> | A SummarizedExperiment object. |
| <code>n_neighbors</code> | Numeric. Indicates the size of local neighborhood (sample points) used for manifold approximation. |
| <code>n_components</code> | Numeric. Indicates the dimension of the space to embed into. |
| <code>metric</code> | Character. Indicates the distance measure method to find nearest neighbors. Options are "euclidean", "cosine", "manhattan", "hamming" and "correlation". See <code>?uwot::umap()</code> . |
| <code>pca</code> | If not NULL (default), reduce data to this number of columns using PCA before UMAP. |
| <code>min_dist</code> | Numeric. Indicates the effective minimum distance between embedded points. |
| <code>spread</code> | Numeric. Indicates the effective scale of embedded points. |
| <code>hdbscan_minpts</code> | Numeric. Indicates the minimum size of clusters. See <code>?hdbscan::hdbscan()</code> . |
| <code>show_clusters</code> | Logical. Indicates if clusters computed with HDBSCAN method should be plotted or not. |
| <code>hide_noise</code> | Logical. Specifies whether to hide Cluster 0 in the plot. In HDBSCAN, Cluster 0 is typically regarded as "noise." |
| <code>labels</code> | Logical. Indicates if sample names should be plotted or not. |
| <code>theme_params</code> | List. Indicates <code>theme_poma</code> parameters. |

Value

A list with results including plots and tables.

Author(s)

Pol Castellano-Escuder

References

McInnes, L., Healy, J., & Melville, J. (2018). Umap: Uniform manifold approximation and projection for dimension reduction. arXiv preprint arXiv:1802.03426.

Campello, R. J., Moulavi, D., & Sander, J. (2013, April). Density-based clustering based on hierarchical density estimates. In Pacific-Asia conference on knowledge discovery and data mining (pp. 160-172). Springer, Berlin, Heidelberg.

Examples

```
data("st000284")

st000284 %>%
  PomaNorm() %>%
  PomaUMAP()
```

Description

PomaUnivariate performs parametric and non-parametric univariate statistical tests on a SummarizedExperiment object to compare groups or conditions. Available methods include T-test, ANOVA, ANCOVA, Mann Whitney U Test (Wilcoxon Rank Sum Test), and Kruskal-Wallis.

Usage

```
PomaUnivariate(
  data,
  method = "ttest",
  covs = NULL,
  error = NULL,
  paired = FALSE,
  var_equal = FALSE,
  adjust = "fdr",
  run_post_hoc = TRUE
)
```

Arguments

| | |
|--------------|---|
| data | A SummarizedExperiment object. |
| method | Character. The univariate statistical test to be performed. Available options include "ttest" (T-test), "anova" (analysis of variance), "mann" (Wilcoxon rank-sum test), and "kruskal" (Kruskal-Wallis test). |
| covs | Character vector. Indicates the names of colData columns to be included as covariates. Default is NULL (no covariates). If not NULL, an ANCOVA model will be fitted using the specified covariates. Note: The order of the covariates is important and should be listed in increasing order of importance in the experimental design. |
| error | Character vector. Indicates the name of a colData column to be included as an error term (e.g. replicates). Default is NULL (no error term). |
| paired | Logical. Indicates if the data is paired or not. Default is FALSE. |
| var_equal | Logical. Indicates if the data variances are assumed to be equal or not. Default is FALSE. |
| adjust | Character. Multiple comparisons correction method to adjust p-values. Available options are: "fdr" (false discovery rate), "holm", "hochberg", "hommel", "bonferroni", "BH" (Benjamini-Hochberg), and "BY" (Benjamini-Yekutieli). |
| run_post_hoc | Logical. Indicates if computing post-hoc tests or not. Setting this parameter to FALSE can save time for large datasets. |

Value

A list with the results.

Author(s)

Pol Castellano-Escuder

Examples

```
data("st000336")

# Perform T-test
st000336 %>%
PomaImpute() %>%
PomaUnivariate(method = "ttest")

# Perform Mann-Whitney U test
st000336 %>%
PomaImpute() %>%
PomaUnivariate(method = "mann", adjust = "fdr")

data("st000284")
# Perform Two-Way ANOVA
st000284 %>%
PomaUnivariate(method = "anova", covs = c("gender"))

# Perform Three-Way ANOVA
st000284 %>%
PomaUnivariate(method = "anova", covs = c("gender", "smoking_condition"))

# Perform ANCOVA with one numeric covariate and one factor covariate
# st000284 %>%
# PomaUnivariate(method = "anova", covs = c("age_at_consent", "smoking_condition"))

# Perform Kruskal-Wallis test
st000284 %>%
PomaUnivariate(method = "kruskal", adjust = "holm")
```

PomaVolcano

Volcano Plot

Description

PomaVolcano creates a volcano plot from a given dataset. This function is designed to visualize the statistical significance (p-value) against the magnitude of change (log2 fold change) for features.

Usage

```
PomaVolcano(
  data,
  pval_cutoff = 0.05,
  log2fc_cutoff = NULL,
  labels = FALSE,
  x_label = "log2 (Fold Change)",
  y_label = "-log10 (P-value)"
)
```

Arguments

| | |
|----------------------------|--|
| <code>data</code> | A data frame with at least three columns: feature names, statistical significance values, and magnitude of change values. These should be the first three columns of the data, in this exact order. |
| <code>pval_cutoff</code> | Numeric. Specifies the p-value threshold for significance in the volcano plot. The default is set to 0.05. This parameter determines the horizontal line in the plot indicating the threshold for statistical significance. |
| <code>log2fc_cutoff</code> | Numeric. Specifies the log2 fold change cutoff for the volcano plot. If not provided, the cutoff is set to the 75th percentile of the absolute log2 fold changes in the data. This parameter determines the vertical lines in the plot indicating the magnitude of change threshold. |
| <code>labels</code> | Logical. Indicates whether to plot labels for significant features. |
| <code>x_label</code> | Character. Custom label for the x-axis. |
| <code>y_label</code> | Character. Custom label for the y-axis. |

Value

A ggplot object representing the volcano plot.

Author(s)

Pol Castellano-Escuder

Examples

```
st000336 %>%
PomaImpute() %>%
PomaUnivariate() %>%
magrittr::extract2("result") %>%
dplyr::select(feature, fold_change, pvalue) %>%
PomaVolcano()
```

| | |
|------------|---|
| poma_pal_c | <i>Return function to interpolate a continuous POMA color palette</i> |
|------------|---|

Description

Return function to interpolate a continuous POMA color palette

Usage

```
poma_pal_c(palette = "nature")
```

Arguments

| | |
|---------|--|
| palette | Character name of palette in poma_palettes |
|---------|--|

| | |
|------------|---|
| poma_pal_d | <i>Return function to interpolate a discrete POMA color palette</i> |
|------------|---|

Description

Return function to interpolate a discrete POMA color palette

Usage

```
poma_pal_d(palette = "nature")
```

Arguments

| | |
|---------|--|
| palette | Character name of palette in poma_palettes |
|---------|--|

| | |
|---------------|--------------------------------------|
| quantile_norm | <i>Sample Quantile Normalization</i> |
|---------------|--------------------------------------|

Description

Compute quantile normalization.

Usage

```
quantile_norm(data)
```

Arguments

| | |
|------|----------------------------------|
| data | A data matrix (samples in rows). |
|------|----------------------------------|

scale_color_poma_c *Color scale constructor for continuous viridis "plasma" palette*

Description

Color scale constructor for continuous viridis "plasma" palette

Usage

```
scale_color_poma_c()
```

scale_color_poma_d *Color scale constructor for discrete viridis "plasma" palette*

Description

Color scale constructor for discrete viridis "plasma" palette

Usage

```
scale_color_poma_d()
```

scale_fill_poma_c *Fill scale constructor for continuous viridis "plasma" palette*

Description

Fill scale constructor for continuous viridis "plasma" palette

Usage

```
scale_fill_poma_c()
```

scale_fill_poma_d *Fill scale constructor for discrete viridis "plasma" palette*

Description

Fill scale constructor for discrete viridis "plasma" palette

Usage

```
scale_fill_poma_d()
```

st000284

Colorectal Cancer Detection Using Targeted Serum Metabolic Profiling

Description

Colorectal cancer (CRC) is one of the most prevalent and deadly cancers in the world. Despite an expanding knowledge of its molecular pathogenesis during the past two decades, robust biomarkers to enable screening, surveillance, and therapy monitoring of CRC are still lacking. In this study, we present a targeted liquid chromatography-tandem mass spectrometry-based metabolic profiling approach for identifying biomarker candidates that could enable highly sensitive and specific CRC detection using human serum samples. In this targeted approach, 158 metabolites from 25 metabolic pathways of potential significance were monitored in 234 serum samples from three groups of patients (66 CRC patients, 76 polyp patients, and 92 healthy controls). Partial least squares-discriminant analysis (PLS-DA) models were established, which proved to be powerful for distinguishing CRC patients from both healthy controls and polyp patients. Receiver operating characteristic curves generated based on these PLS-DA models showed high sensitivities (0.96 and 0.89, respectively, for differentiating CRC patients from healthy controls or polyp patients); good specificities (0.80 and 0.88), and excellent areas under the curve (0.93 and 0.95) were also obtained. Monte Carlo cross validation (MCCV) was also applied, demonstrating the robust diagnostic power of this metabolic profiling approach.

Usage

st000284

Format

A SummarizedExperiment object: 224 samples, 113 metabolites, 4 covariables and 3 groups (CRC, Healthy and Polyp).

metabolites 113 serum metabolites.

covariables Age at consent, Gender, Smoking Condition and Alcohol Consumption.

Source

https://www.metabolomicsworkbench.org/data/DRCCMetadata.php?Mode=Study&StudyID=ST000284&StudyType=MS&ResultType=1%20target=_blank

References

Colorectal Cancer Detection Using Targeted Serum Metabolic Profiling, J. Proteome. Res., 2014, 13, 4120-4130.

`st000336`*Targeted LC/MS of urine from boys with DMD and controls*

Description

Duchenne Muscular Dystrophy (DMD) is an X-linked recessive form of muscular dystrophy that affects males via a mutation in the gene for the muscle protein, dystrophin. Progression of the disease results in severe muscle loss, ultimately leading to paralysis and death. Steroid therapy has been a commonly employed method for reducing the severity of symptoms. This study aims to quantify the urine levels of amino acids and organic acids in patients with DMD both with and without steroid treatment. Track the progression of DMD in patients who have provided multiple urine samples.

Usage`st000336`**Format**

A SummarizedExperiment object: 57 samples, 31 metabolites, 1 covariable and 2 groups (Controls and DMD).

metabolites 31 urine metabolites.

covariables Steroid status.

Source

<https://www.metabolomicsworkbench.org/data/DRCCMetadata.php?Mode=Study&DataMode=AllData&StudyID=ST000336&StudyType=MS&ResultType=1#DataTabs>

`sum_norm`*Sample Sum Normalization*

Description

Compute sum normalization. Final unit is a percentage.

Usage`sum_norm(data)`**Arguments**

`data` A data matrix (samples in rows).

| | |
|------------|---|
| theme_poma | <i>A ggplot theme which allow custom yet consistent styling of plots in the POMA package and web app.</i> |
|------------|---|

Description

A ggplot theme which allow custom yet consistent styling of plots in the POMA package and web app.

Usage

```
theme_poma(  
  base_size = 15,  
  axistitle = "xy",  
  axistext = "xy",  
  legend_position = "bottom",  
  legend_title = TRUE,  
  axis_x_rotate = FALSE,  
  margin = 2  
)
```

Arguments

| | |
|-----------------|---|
| base_size | (integer) Base point size |
| axistitle | (string) Axis titles. Options include "none" or any combination of "X", "Y", "x" and "y". |
| axistext | (string) Axis text labels for values or groups. Options include "none" or any combination of "X", "Y", "x" and "y". |
| legend_position | Character. Legend position. See ggplot2 documentation. |
| legend_title | Logical. Include legend title. |
| axis_x_rotate | Logical. Rotate x-axis 45 degrees. |
| margin | (numeric) Should a margin of x be added to the plot? Defaults to 0 (no margin by default). |

Examples

```
## Not run:  
library(ggplot2)  
ggplot(diamonds, aes(cut)) + geom_bar() + theme_poma()  
  
## End(Not run)
```

%>%

Pipe operator

Description

See `magrittr::%>%` for details.

Usage

lhs %>% rhs

Value

Nothing. Just allow the use of magrittr pipe "%>%"

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