Package ‘Ecume’

Type Package
Title Equality of 2 (or k) Continuous Univariate and Multivariate Distributions
Version 0.9.1
Description We implement (or re-implements in R) a variety of statistical tools. They are focused on non-parametric two-sample (or k-sample) distribution comparisons in the univariate or multivariate case. See the vignette for more info.
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**classifier_test**

*Classifier k-sample test*

**Description**
Classifier k-sample test

**Usage**

```r
classifier_test(
  x,
  y,
  split = 0.7,
  thresh = 0,
  method = "knn",
  control = caret::trainControl(method = "cv"),
  ...
)
```

**Arguments**

- `x` Samples from the first distribution or a list of samples from k distribution
- `y` Samples from the second distribution. Only used if x is a vector.
- `split` How to split the data between training and test. Default to .7
- `thresh` Value to add to the null hypothesis. See details.
- `method` Which model(s) to use during training. Default to knn.
- `control` Control parameters when fitting the methods. See `trainControl`
- `...` Other parameters passed to `train`

**Details**

See Lopez-Paz et al for more background on those tests.

**Value**

A list containing the following components:

- `statistic` the value of the test statistic.
- `p.value` the p-value of the test.

**References**

Examples

```r
x <- matrix(c(runif(100, 0, 1),
              runif(100, -1, 1)),
             ncol = 2)
y <- matrix(c(runif(100, 0, 3),
              runif(100, -1, 1)),
             ncol = 2)
classifier_test(x, y)
```

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**ks_test**  
*Weighted KS Test*

**Description**
Weighted Kolmogorov-Smirnov Two-Sample Test with threshold

**Usage**
```r
ks_test(x, y, thresh = 0.05, w_x = rep(1, length(x)), w_y = rep(1, length(y)))
```

**Arguments**
- `x`: Vector of values sampled from the first distribution
- `y`: Vector of values sampled from the second distribution
- `thresh`: The threshold needed to clear between the two cumulative distributions
- `w_x`: The observation weights for x
- `w_y`: The observation weights for y

**Details**
The usual Kolmogorov-Smirnov test for two vectors `X` and `Y`, of size `m` and `n` rely on the empirical cdfs `E_x` and `E_y` and the test statistic

\[
D = \sup_{t \in (X,Y)} |E_x(t) - E_y(t) |
\]

This modified Kolmogorov-Smirnov test relies on two modifications.

- Using observation weights for both vectors `X` and `Y`: Those weights are used in two places, while modifying the usual KS test. First, the empirical cdfs are updates to account for the weights. Secondly, the effective sample sizes are also modified. This is inspired from [https://stackoverflow.com/a/55664242/13768995](https://stackoverflow.com/a/55664242/13768995), using Monahan (2011).
- Testing against a threshold: the test statistic is thresholded such that \( D = \max(D - \text{thresh}, 0) \). Since \( 0 \leq D \leq 1 \), the value of the threshold is also between 0 and 1, representing an effect size for the difference.
Value

A list with class "htest" containing the following components:

- **statistic** the value of the test statistic.
- **p.value** the p-value of the test.
- **alternative** a character string describing the alternative hypothesis.
- **method** a character string indicating what type of test was performed.
- **data.name** a character string giving the name(s) of the data.

References


Examples

```r
x <- runif(100)
y <- runif(100, min = .5, max = .5)
ks_test(x, y, thresh = .001)
```

mmd_test

Perform the Maximum Mean Discrepancy unbiased bootstrap test

Description

Maximum Mean Discrepancy Unbiased Test

Usage

```r
mmd_test(
	x,
	y,


type = ifelse(min(nrow(x), nrow(y)) < 1000, "unbiased", "linear"),


null = c("permutation", "exact"),

iterations = 10^3,

frac = 1,

...
)
```

Arguments

- **x** d-dimensional samples from the first distribution
- **y** d-dimensional samples from the first distribution
- **kernel** A character that must match a known kernel. See details.
mmd_test

**type**
Which statistic to use. One of 'unbiased' or 'linear'. See Gretton et al for details. Default to 'unbiased' if the two vectors are of length less than 1000 and to 'linear' otherwise.

**null**
How to assess the null distribution. This can only be set to exact if the type is 'unbiased' and the kernel is 'rbf'.

**iterations**
How many iterations to do to simulate the null distribution. Default to $10^4$. Only used if null is 'permutations'.

**frac**
For the linear statistic, how many points to sample. See details.

**...**
Further arguments passed to kernel functions

**Details**
This computes the MMD^2u unbiased statistic or the MMDl linear statistic from Gretton et al. The code relies on the pairwise_kernel function from the python module sklearn. To list the available kernels, see the examples.

**Value**
A list containing the following components:

- **statistic** the value of the test statistic.
- **p.value** the p-value of the test.

**References**

**Examples**
```r
x <- matrix(rnorm(1000, 0, 1), ncol = 10)
y <- matrix(rnorm(1000, 0, 2), ncol = 10)
mmd_test(x, y)
mmd_test(x, y, type = "linear")
x <- matrix(rnorm(1000, 0, 1), ncol = 10)
y <- matrix(rnorm(1000, 0, 1), ncol = 10)
# Set iterations to small number for runtime
# Increase for more accurate results
mmd_test(x, y, iterations = 10^2)
```
Description

Stouffer's Z-score method

Usage

`stouffer_zscore(pvals, weights = rep(1, seq_along(pvals)), side = "two")`

Arguments

- `pvals` A vector of p-values
- `weights` A vector of weights
- `side` How the p-values were generated. One of 'right', 'left' or 'two'.

Details

Given a set of i.i.d p-values and associated weights, it combines the p-values $p_i$. Letting $\phi$ be the standard normal cumulative distribution function and $Z_i = \phi^{-1}(1 - p_i)$, the meta-analysis Z-score is

$$Z = (\sum w_i Z_i) * (\sum (w_i)^2)^{-1/2}$$

Value

A list containing the following components:

- `statistic` the value of the test statistic.
- `p.value` the p-value of the test.

References


Examples

```r
pvals <- runif(100, 0, 1)
weights <- runif(100, 0, 1)
stouffer_zscore(pvals, weights)
```
Permutation test based on Wasserstein distance

Usage

wasserstein_permut(
  x,
  y,
  iterations = 10^4,
  fast = nrow(x) + nrow(y) > 10^3,
  S = NULL,
  ...
)

Arguments

x  Samples from the first distribution
y  Samples from the second distribution. Only used if x is a vector.
iterations How many iterations to do to simulate the null distribution. Default to 10^4.
fast If true, uses the subwasserstein approximate function. Default to true if there are more than 1,000 samples total.
S  Number of samples to use in approximate mode. Must be set if fast=TRUE. See subwasserstein.
... Other parameters passed to wasserstein or wasserstein1d

Value

A list containing the following components:

- statistic the Wasserstein distance between x and y.
- p.value the p-value of the permutation test.

Examples

x <- matrix(c(runif(100, 0, 1),
  runif(100, -1, 1)),
  ncol = 2)
y <- matrix(c(runif(100, 0, 3),
  runif(100, -1, 1)),
  ncol = 2)
# Set iterations to small number for runtime
# Increase for more accurate results
wasserstein_permut(x, y, iterations = 10^2)
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