Package ‘fxTWAPLS’

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Title  An Improved Version of WA-PLS
Version  0.1.3
Description  The goal of this package is to provide an improved version of WA-PLS (Weighted Averaging Partial Least Squares) by including the tolerances of taxa and the frequency of the sampled climate variable. This package also provides a way of leave-out cross-validation that removes both the test site and sites that are both geographically close and climatically close for each cycle, to avoid the risk of pseudo-replication.
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cv.pr.w  Pseudo-removed leave-out cross-validation

Description

Pseudo-removed leave-out cross-validation

Usage

cv.pr.w(
  modern_taxa,
  modern_climate,
  nPLS = 5,
  trainfun,
  predictfun,
  pseudo,
  usefx = FALSE,
  fx_method = "bin",
  bin = NA,
  cpus = 4,
  test_mode = TRUE,
  test_it = 5
)
Arguments

modern_taxa  The modern taxa abundance data, each row represents a sampling site, each column represents a taxon.
modern_climate  The modern climate value at each sampling site.
nPLS  The number of components to be extracted.
trainfun  Training function you want to use, either WAPLS.w or TWAPLS.w.
predictfun  Predict function you want to use: if trainfun is WAPLS.w, then this should be WAPLS.predict.w; if trainfun is TWAPLS.w, then this should be TWAPLS.predict.w.
pseudo  The geographically and climatically close sites to each test site, obtained from get_pseudo function.
usefx  Boolean flag on whether or not use fx correction.
fx_method  Binned or p-spline smoothed fx correction: if usefx = FALSE, this should be NA; otherwise, fx function will be used when choosing "bin"; fx_pspline function will be used when choosing "pspline".
bin  Binwidth to get fx, needed for both binned and p-splined method. if usefx = FALSE, this should be NA;
cpus  Number of CPUs for simultaneous iterations to execute, check parallel::detectCores() for available CPUs on your machine.
test_mode  Boolean flag to execute the function with a limited number of iterations, test_it, for testing purposes only.
test_it  Number of iterations to use in the test mode.

Value

Leave-one-out cross validation results.

See Also

fx, TWAPLS.w, TWAPLS.predict.w, WAPLS.w, and WAPLS.predict.w

Examples

```r
## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

# Extract taxa
taxaColMin <- which(colnames(moderne_pollen) == "taxa0")
taxaColMax <- which(colnames(moderne_pollen) == "taxaN")
taxa <- modern_pollen[, taxaColMin:taxaColMax]

point <- modern_pollen[, c("Long", "Lat")]
test_mode <- TRUE # It should be set to FALSE before running
dist <- fxTWAPLS::get_distance(
  point,
  cpus = 2, # Remove the following line
)```
test_mode = test_mode
)
pseudo_Tmin <- fxTWAPLS::get_pseudo(
  dist,
  modern_pollen$Tmin,
  cpus = 2, # Remove the following line
  test_mode = test_mode
)

cv_pr_tf_Tmin2 <- fxTWAPLS::cv.pr.w(
  taxa,
  modern_pollen$Tmin,
  nPLS = 5,
  fxTWAPLS::TWAPLS.w2,
  fxTWAPLS::TWAPLS.predict.w,
  pseudo_Tmin,
  usefx = TRUE,
  fx_method = "bin",
  bin = 0.02,
  cpus = 2, # Remove the following line
  test_mode = test_mode
)

# Run with progress bar
`%>%` <- magrittr::`%>%`
cv_pr_tf_Tmin2 <- fxTWAPLS::cv.pr.w(
  taxa,
  modern_pollen$Tmin,
  nPLS = 5,
  fxTWAPLS::TWAPLS.w2,
  fxTWAPLS::TWAPLS.predict.w,
  pseudo_Tmin,
  usefx = TRUE,
  fx_method = "bin",
  bin = 0.02,
  cpus = 2, # Remove the following line
  test_mode = test_mode
) %>%
f TWAPLS::pb()

## End(Not run)

---

**cv.w**

*Leave-one-out cross-validation*

---

**Description**

Leave-one-out cross-validation as rioja ([https://cran.r-project.org/package=rioja](https://cran.r-project.org/package=rioja)).
Usage

```r
cv.w(
  modern_taxa,
  modern_climate,
  nPLS = 5,
  trainfun,
  predictfun,
  usefx = FALSE,
  fx_method = "bin",
  bin = NA,
  cpus = 4,
  test_mode = FALSE,
  test_it = 5
)
```

Arguments

- `modern_taxa`: The modern taxa abundance data, each row represents a sampling site, each column represents a taxon.
- `modern_climate`: The modern climate value at each sampling site.
- `nPLS`: The number of components to be extracted.
- `trainfun`: Training function you want to use, either `WAPLS.w` or `TWAPLS.w`.
- `predictfun`: Predict function you want to use: if `trainfun` is `WAPLS.w`, then this should be `WAPLS.predict.w`; if `trainfun` is `TWAPLS.w`, then this should be `TWAPLS.predict.w`.
- `usefx`: Boolean flag on whether or not use `fx` correction.
- `fx_method`: Binned or p-spline smoothed `fx` correction: if `usefx = FALSE`, this should be `NA`; otherwise, `fx` function will be used when choosing "bin"; `fx_pspline` function will be used when choosing "pspline".
- `bin`: Binwidth to get `fx`, needed for both binned and p-splined method. if `usefx = FALSE`, this should be `NA`;
- `cpus`: Number of CPUs for simultaneous iterations to execute, check `parallel::detectCores()` for available CPUs on your machine.
- `test_mode`: boolean flag to execute the function with a limited number of iterations, `test_it`, for testing purposes only.
- `test_it`: number of iterations to use in the test mode.

Value

leave-one-out cross validation results

See Also

`fx`, `TWAPLS.w`, `TWAPLS.predict.w`, `WAPLS.w`, and `WAPLS.predict.w`
Examples

```r
## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

# Extract taxa
taxaColMin <- which(colnames(modern_pollen) == "taxa0")
taxaColMax <- which(colnames(modern_pollen) == "taxaN")
taxa <- modern_pollen[, taxaColMin:taxaColMax]

## LOOCV
test_mode <- TRUE # It should be set to FALSE before running
cv_tf_Tmin2 <- fxTWAPLS::cv.w(taxa, modern_pollen$Tmin, nPLS = 5, fxTWAPLS::TWAPLS.w2, fxTWAPLS::TWAPLS.predict.w, usefx = TRUE, fx_method = "bin", bin = 0.02, cpus = 2, # Remove the following line
test_mode = test_mode )

# Run with progress bar
`%>%` <- magrittr::>%>
cv_tf_Tmin2 <- fxTWAPLS::cv.w(taxa, modern_pollen$Tmin, nPLS = 5, fxTWAPLS::TWAPLS.w2, fxTWAPLS::TWAPLS.predict.w, usefx = TRUE, fx_method = "bin", bin = 0.02, cpus = 2, # Remove the following line
test_mode = test_mode ) %>% fxTWAPLS::pb()
```

## End(Not run)

---

**fx**

*Get frequency of the climate value*

---

**Description**

Function to get the frequency of the climate value, which will be used to provide fx correction for WA-PLS and TWA-PLS.
fx_pspline

Usage

fx(x, bin, show_plot = FALSE)

Arguments

x Numeric vector with the modern climate values.
bin Binwidth to get the frequency of the modern climate values.
show_plot Boolean flag to show a plot of fx ~ x.

Value

Numeric vector with the frequency of the modern climate values.

See Also

cv.w, cv.pr.w, and sse.sample

Examples

## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

# Get the frequency of each climate variable fx
fx_Tmin <- fxTWAPLS::fx(moderan_pollen$Tmin, bin = 0.02, show_plot = TRUE)
fx_gdd <- fxTWAPLS::fx(moderan_pollen$gdd, bin = 20, show_plot = TRUE)
fx_alpha <- fxTWAPLS::fx(moderan_pollen$alpha, bin = 0.002, show_plot = TRUE)

## End(Not run)

fx_pspline

Get frequency of the climate value with p-spline smoothing

Description

Function to get the frequency of the climate value, which will be used to provide fx correction for WA-PLS and TWA-PLS.

Usage

fx_pspline(x, bin, show_plot = FALSE)

Arguments

x Numeric vector with the modern climate values.
bin Binwidth to get the frequency of the modern climate values, the curve will be p-spline smoothed later
show_plot Boolean flag to show a plot of fx ~ x.
get_distance

Description

Get the distance between points, the output will be used in `get_pseudo`.

Usage

```r
get_distance(point, cpus = 4, test_mode = FALSE, test_it = 5)
```

Arguments

- **point**: Each row represents a sampling site, the first column is longitude and the second column is latitude, both in decimal format.
**get_pseudo**

```r
get_pseudo
```

**cpus**  
Number of CPUs for simultaneous iterations to execute, check `parallel::detectCores()` for available CPUs on your machine.

**test_mode**  
Boolean flag to execute the function with a limited number of iterations, `test_it`, for testing purposes only.

**test_it**  
Number of iterations to use in the test mode.

**Value**

Distance matrix, the value at the $i$-th row, means the distance between the $i$-th sampling site and the whole sampling sites.

**See Also**

- `get_pseudo`

**Examples**

```r
## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

point <- modern_pollen[, c("Long", "Lat")]
test_mode <- TRUE  # It should be set to FALSE before running
dist <- fxTWAPLS::get_distance(
  point,  
  cpus = 2, # Remove the following line
  test_mode = test_mode
)

# Run with progress bar
> dist <- fxTWAPLS::get_distance(
  point,  
  cpus = 2, # Remove the following line
  test_mode = test_mode
) >
  fxTWAPLS::pb()

## End(Not run)
```

---

**get_pseudo**  
*Get geographically and climatically close sites*

**Description**

Get the sites which are both geographically and climatically close to the test site, which could result in pseudo-replication and inflate the cross-validation statistics. The output will be used in `cv.pr.w`. 
get_pseudo

Usage

get_pseudo(dist, x, cpus = 4, test_mode = FALSE, test_it = 5)

Arguments

dist Distance matrix which contains the distance from other sites.
x The modern climate values.
cpus Number of CPUs for simultaneous iterations to execute, check parallel::detectCores() for available CPUs on your machine.
test_mode Boolean flag to execute the function with a limited number of iterations, test_it, for testing purposes only.
test_it Number of iterations to use in the test mode.

Value

The geographically and climatically close sites to each test site.

See Also

get_distance

Examples

## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

point <- modern_pollen[, c("Long", "Lat")]
test_mode <- TRUE  # It should be set to FALSE before running
dist <- fxTWAPLS::get_distance(
    point,
    cpus = 2, # Remove the following line
    test_mode = test_mode
)
pseudo_Tmin <- fxTWAPLS::get_pseudo(
    dist,
    modern_pollen$Tmin,
    cpus = 2, # Remove the following line
    test_mode = test_mode
)
# Run with progress bar
`%>%` <- magrittr::`%>%`
pseudo_Tmin <- fxTWAPLS::get_pseudo(
    dist,
    modern_pollen$Tmin,
    cpus = 2, # Remove the following line
    test_mode = test_mode
)
`%>%`
fTWAPLS::pb()
pb

## End(Not run)

---

pb  

*Show progress bar*

---

**Description**

Show progress bar

**Usage**

```r
pb(expr, ...)
```

**Arguments**

- `expr`  
  - R expression.
- `...`  
  - Arguments passed on to `progressr::with_progress`
- `cleanup`  
  - If TRUE, all progression handlers will be shutdown at the end regardless of the progression is complete or not.
- `delay_terminal`  
  - If TRUE, output and conditions that may end up in the terminal will delayed.
- `delay_stdout`  
  - If TRUE, standard output is captured and relayed at the end just before any captured conditions are relayed.
- `delay_conditions`  
  - A character vector specifying `base::condition` classes to be captured and relayed at the end after any captured standard output is relayed.
- `interrupts`  
  - Controls whether interrupts should be detected or not. If TRUE and a interrupt is signaled, progress handlers are asked to report on the current amount progress when the evaluation was terminated by the interrupt, e.g. when a user pressed Ctrl-C in an interactive session, or a batch process was interrupted because it ran out of time. Note that it’s optional for a progress handler to support this and only some do.
- `interval` (numeric)  
  - The minimum time (in seconds) between successive progress updates from handlers.
- `enable` (logical)  
  - If FALSE, then progress is not reported. The default is to report progress in interactive mode but not batch mode. See below for more details.

**Value**

Return data from the function called.
plot_residuals  

Plot the residuals

Description
Plot the residuals, the black line is 0 line, the red line is the locally estimated scatterplot smoothing, which shows the degree of local compression.

Usage
plot_residuals(train_output, col)

Arguments
- train_output: Training output, can be the output of WA-PLS, WA-PLS with fx correction, TWA-PLS, or TWA-PLS with fx correction
- col: Choose which column of the fitted value to plot, in other words, how many number of components you want to use.

Value
Plotting status.

See Also
TWAPLS.w and WAPLS.w

Examples
```r
## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

# Extract taxa
taxaColMin <- which(colnames(modern_pollen) == "taxa0")
taxaColMax <- which(colnames(modern_pollen) == "taxaN")
taxa <- modern_pollen[, taxaColMin:taxaColMax]

fit_tf_Tmin2 <- fxTWAPLS::TWAPLS.w2(
taxa,
modern_pollen$Tmin,
nPLS = 5,
usefx = TRUE,
fx_method = "bin",
bin = 0.02
)

nsig <- 3 # This should be got from the random t-test of the cross validation
fxTWAPLS::plot_residuals(fit_tf_Tmin2, nsig)
```
plot_train

## End(Not run)

---

**plot_train**  
*Plot the training results*

### Description

Plot the training results, the black line is the 1:1 line, the red line is the linear regression line to fitted and $x$, which shows the degree of overall compression.

### Usage

```r
plot_train(train_output, col)
```

### Arguments

- **train_output**: Training output, can be the output of WA-PLS, WA-PLS with $fx$ correction, TWA-PLS, or TWA-PLS with $fx$ correction.
- **col**: Choose which column of the fitted value to plot, in other words, how many number of components you want to use.

### Value

Plotting status.

### See Also

`TWAPLS.w` and `WAPLS.w`

### Examples

```r
## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

# Extract taxa
taxaColMin <- which(colnames(modern_pollen) == "taxa0")
taxaColMax <- which(colnames(modern_pollen) == "taxaN")
taxa <- modern_pollen[, taxaColMin:taxaColMax]

fit_tf_Tmin2 <- fxTWAPLS::TWAPLS.w2(
taxa, 
modern_pollen$Tmin, 
nPLS = 5, 
usefx = TRUE, 
fx_method = "bin", 
bin = 0.02
```
nsig <- 3 # This should be got from the random t-test of the cross validation
fxTWAPLS::plot_train(fit_tf_Tmin2, nsig)

## End(Not run)

---

**rand.t.test.w**

**Random t-test**

**Description**

Do a random t-test to the cross-validation results.

**Usage**

```r
rand.t.test.w(cvoutput, n.perm = 999)
```

**Arguments**

- `cvoutput` Cross-validation output either from `cv.w` or `cv.pr.w`.
- `n.perm` The number of permutation times to get the p value, which assesses whether using the current number of components is significantly different from using one less.

**Value**

A matrix of the statistics of the cross-validation results. Each component is described below:

- **R2** the coefficient of determination (the larger, the better the fit).
- **Avg.Bias** average bias.
- **Max.Bias** maximum bias.
- **Min.Bias** minimum bias.
- **RMSEP** root-mean-square error of prediction (the smaller, the better the fit).
- **delta.RMSEP** the percent change of RMSEP using the current number of components than using one component less.
- **p** assesses whether using the current number of components is significantly different from using one component less, which is used to choose the last significant number of components to avoid over-fitting.

- The degree of overall compression is assessed by doing linear regression to the cross-validation result and the observed climate values.
  - **Compre.b0**: the intercept.
  - **Compre.b1**: the slope (the closer to 1, the less the overall compression).
  - **Compre.b0.se**: the standard error of the intercept.
  - **Compre.b1.se**: the standard error of the slope.
See Also

cv.w and cv.pr.w

Examples

## Not run:

## Random t-test
rand_pr_tf_Tmin2 <- fxTWAPLS::rand.t.test.w(cv_pr_tf_Tmin2, n_perm = 999)

# note: choose the last significant number of components based on the p-value,
# see details at Liu Mengmeng, Prentice Iain Colin, ter Braak Cajo J. F.,
# Harrison Sandy P.. 2020 An improved statistical approach for reconstructing
# <https://doi.org/10.1098/rspa.2020.0346>

## End(Not run)

---

sse.sample  Calculate Sample Specific Errors

Description

Calculate Sample Specific Errors

Usage

sse.sample(
  modern_taxa,
  modern_climate,
  fossil_taxa,
  trainfun,
  predictfun,
  nboot,
  nPLS,
  nsig,
  usefx = FALSE,
  fx_method = "bin",
  bin = NA,
  cpus = 4,
  seed = NULL,
  test_mode = FALSE,
  test_it = 5
)
Arguments

modern_taxa  The modern taxa abundance data, each row represents a sampling site, each column represents a taxon.

modern_climate  The modern climate value at each sampling site.

fossil_taxa  Fossil taxa abundance data to reconstruct past climates, each row represents a site to be reconstructed, each column represents a taxon.

trainfun  Training function you want to use, either WAPLS.w or TWAPLS.w.

predictfun  Predict function you want to use: if trainfun is WAPLS.w, then this should be WAPLS.predict.w; if trainfun is TWAPLS.w, then this should be TWAPLS.predict.w.

nboot  The number of bootstrap cycles you want to use.

nPLS  The number of components to be extracted.

nsig  The significant number of components to use to reconstruct past climates, this can be obtained from the cross-validation results.

usefx  Boolean flag on whether or not use fx correction.

fx_method  Binned or p-spline smoothed fx correction: if usefx = FALSE, this should be NA; otherwise, fx function will be used when choosing "bin"; fx_pspline function will be used when choosing "pspline".

bin  Binwidth to get fx, needed for both binned and p-splined method. if usefx = FALSE, this should be NA;

cpus  Number of CPUs for simultaneous iterations to execute, check parallel::detectCores() for available CPUs on your machine.

seed  Seed for reproducibility.

test_mode  Boolean flag to execute the function with a limited number of iterations, test_it, for testing purposes only.

test_it  Number of iterations to use in the test mode.

Value

The bootstrapped standard error for each site.

See Also

fx, TWAPLS.w, TWAPLS.predict.w, WAPLS.w, and WAPLS.predict.w

Examples

```r
## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

# Extract taxa
taxaColMin <- which(colnames(moderan_pollen) == "taxa0")
taxaColMax <- which(colnames(moderan_pollen) == "taxaN")
taxa <- modern_pollen[, taxaColMin:taxaColMax]
```
# Load reconstruction data
Holocene <- read.csv("/path/to/Holocene.csv")
taxaColMin <- which(colnames(Holocene) == "taxa0")
taxaColMax <- which(colnames(Holocene) == "taxaN")
core <- Holocene[, taxaColMin:taxaColMax]

## SSE
nboot <- 5 # Recommended 1000
nsig <- 3 # This should be got from the random t-test of the cross validation
sse_tf_Tmin2 <- fxTWAPLS::sse.sample(
  modern_taxa = taxa,
  modern_climate = modern_pollen$Tmin,
  fossil_taxa = core,
  trainfun = fxTWAPLS::TWAPLS.w2,
  predictfun = fxTWAPLS::TWAPLS.predict.w,
  nboot = nboot,
  nPLS = 5,
  nsig = nsig,
  usefx = TRUE,
  fx_method = "bin",
  bin = 0.02,
  cpus = 2,
  seed = 1
)

# Run with progress bar
`%>%` <- magrittr:: `%>%`
sse_tf_Tmin2 <- fxTWAPLS::sse.sample(
  modern_taxa = taxa,
  modern_climate = modern_pollen$Tmin,
  fossil_taxa = core,
  trainfun = fxTWAPLS::TWAPLS.w2,
  predictfun = fxTWAPLS::TWAPLS.predict.w,
  nboot = nboot,
  nPLS = 5,
  nsig = nsig,
  usefx = TRUE,
  fx_method = "bin",
  bin = 0.02,
  cpus = 2,
  seed = 1
) %>% fxTWAPLS::pb()

## End(Not run)
Description

TWA-PLS predict function

Usage

TWAPLS.predict.w(TWAPLSoutput, fossil_taxa)

Arguments

TWAPLSoutput  The output of the TWAPLS.w training function, either with or without fx correction.
fossil_taxa    Fossil taxa abundance data to reconstruct past climates, each row represents a site to be reconstructed, each column represents a taxon.

Value

A list of the reconstruction results. Each element in the list is described below:

fit  the fitted values using each number of components.
nPLS the total number of components extracted.

See Also

TWAPLS.w

Examples

## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

# Extract taxa
taxaColMin <- which(colnames(moderon_pollen) == "taxa0")
taxaColMax <- which(colnames(moderon_pollen) == "taxaN")
taxa <- modern_pollen[, taxaColMin:taxaColMax]

# Load reconstruction data
Holocene <- read.csv("/path/to/Holocene.csv")
taxaColMin <- which(colnames(Holocene) == "taxa0")
taxaColMax <- which(colnames(Holocene) == "taxaN")
core <- Holocene[, taxaColMin:taxaColMax]

## Train
fit_t_Tmin <- fxTWAPLS::TWAPLS.w(taxa, modern_pollen$Tmin, nPLS = 5)
fit_tf_Tmin <- fxTWAPLS::TWAPLS.w(taxa,
    modern_pollen$Tmin,
    nPLS = 5,
    usefx = TRUE,
    fx_method = "bin",
    bin = 0.02
fit_t_Tmin2 <- fxTWAPLS::TWAPLS.w2(taxa, modern_pollen$Tmin, nPLS = 5)
fit_tf_Tmin2 <- fxTWAPLS::TWAPLS.w2(taxa, modern_pollen$Tmin, nPLS = 5, usefx = TRUE, fx_method = "bin", bin = 0.02)

## Predict
fossil_t_Tmin <- fxTWAPLS::TWAPLS.predict.w(fit_t_Tmin, core)
fossil_tf_Tmin <- fxTWAPLS::TWAPLS.predict.w(fit_tf_Tmin, core)
fossil_t_Tmin2 <- fxTWAPLS::TWAPLS.predict.w(fit_t_Tmin2, core)
fossil_tf_Tmin2 <- fxTWAPLS::TWAPLS.predict.w(fit_tf_Tmin2, core)

## End(Not run)

---

TWAPLS.w  

TWA-PLS training function

Description

TWA-PLS training function, which can perform \( f^x \) correction. \( 1/f^2 \) correction will be applied at step 7.

Usage

TWAPLS.w(
  modern_taxa,  
  modern_climate,  
  nPLS = 5,  
  usefx = FALSE,  
  fx_method = "bin",  
  bin = NA
)

Arguments

- **modern_taxa**: The modern taxa abundance data, each row represents a sampling site, each column represents a taxon.
- **modern_climate**: The modern climate value at each sampling site.
- **nPLS**: The number of components to be extracted.
- **usefx**: Boolean flag on whether or not use \( f^x \) correction.
fx_method  Binned or p-spline smoothed fx correction: if usefx = FALSE, this should be NA; otherwise, fx function will be used when choosing "bin"; fx_pspline function will be used when choosing "pspline".

bin  Binwidth to get fx, needed for both binned and p-splined method. if usefx = FALSE, this should be NA;

Value

A list of the training results, which will be used by the predict function. Each element in the list is described below:

fit  the fitted values using each number of components.
x  the observed modern climate values.
taxon_name  the name of each taxon.
optimum  the updated taxon optimum
comp  each component extracted (will be used in step 7 regression).
u  taxon optimum for each component (step 2).
t  taxon tolerance for each component (step 2).
z  a parameter used in standardization for each component (step 5).
s  a parameter used in standardization for each component (step 5).
ort  a list that stores orthogonalization parameters (step 4).
alpha  a list that stores regression coefficients (step 7).
meanx  mean value of the observed modern climate values.
nPLS  the total number of components extracted.

See Also

fx, TWAPLS.predict.w, and WAPLS.w

Examples

```r
## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

# Extract taxa
taxaColMin <- which(colnames(modern_pollen) == "taxa0")
taxaColMax <- which(colnames(modern_pollen) == "taxaN")
taxa <- modern_pollen[, taxaColMin:taxaColMax]

# Training
fit_t_Tmin <- fxTWAPLS::TWAPLS.w(taxa, modern_pollen$Tmin, nPLS = 5)
fit_tf_Tmin <- fxTWAPLS::TWAPLS.w(taxa,
  modern_pollen$Tmin,
  nPLS = 5,
  usefx = TRUE,
```
TWAPLS.w2

## End(Not run)

---

**TWAPLS.w2**  
*TWA-PLS training function v2*

---

**Description**

TWA-PLS training function, which can perform $fx$ correction. $1/fx$ correction will be applied at step 2 and step 7.

**Usage**

```r
TWAPLS.w2(
    modern_taxa,
    modern_climate,
    nPLS = 5,
    usefx = FALSE,
    fx_method = "bin",
    bin = NA
)
```

**Arguments**

- **modern_taxa**: The modern taxa abundance data, each row represents a sampling site, each column represents a taxon.
- **modern_climate**: The modern climate value at each sampling site.
- **nPLS**: The number of components to be extracted.
- **usefx**: Boolean flag on whether or not use $fx$ correction.
- **fx_method**: Binned or p-spline smoothed $fx$ correction: if `usefx = FALSE`, this should be `NA`; otherwise, $fx$ function will be used when choosing "bin"; `fx_pspline` function will be used when choosing "pspline".
- **bin**: Binwidth to get $fx$, needed for both binned and p-splined method. if `usefx = FALSE`, this should be `NA`;

**Value**

A list of the training results, which will be used by the predict function. Each element in the list is described below:

- **fit**: the fitted values using each number of components.
- **x**: the observed modern climate values.
taxon_name the name of each taxon.

optimum the updated taxon optimum

comp each component extracted (will be used in step 7 regression).

u taxon optimum for each component (step 2).

t taxon tolerance for each component (step 2).

z a parameter used in standardization for each component (step 5).

s a parameter used in standardization for each component (step 5).

orth a list that stores orthogonalization parameters (step 4).

alpha a list that stores regression coefficients (step 7).

meanx mean value of the observed modern climate values.

nPLS the total number of components extracted.

See Also

fx, TWAPLS.predict.w, and WAPLS.w

Examples

```r
## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

# Extract taxa
taxaColMin <- which(colnames(moderne_pollen) == "taxa0")
taxaColMax <- which(colnames(moderne_pollen) == "taxaN")
taxa <- moderne_pollen[, taxaColMin:taxaColMax]

# Training
fit_t_Tmin2 <- fxTWAPLS::TWAPLS.w2(taxa, moderne_pollen$Tmin, nPLS = 5)
fit_tf_Tmin2 <- fxTWAPLS::TWAPLS.w2(
  taxa,
  moderne_pollen$Tmin,
  nPLS = 5,
  usefx = TRUE,
  fx_method = "bin",
  bin = 0.02
)

## End(Not run)
```
WAPLS.predict.w

WA-PLS predict function

Description

WA-PLS predict function

Usage

WAPLS.predict.w(WAPLSoutput, fossil_taxa)

Arguments

- **WAPLSoutput**: The output of the WAPLS.w training function, either with or without fx correction.
- **fossil_taxa**: Fossil taxa abundance data to reconstruct past climates, each row represents a site to be reconstructed, each column represents a taxon.

Value

A list of the reconstruction results. Each element in the list is described below:

- **fit**: The fitted values using each number of components.
- **nPLS**: The total number of components extracted.

See Also

WAPLS.w

Examples

```r
## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

# Extract taxa
taxaColMin <- which(colnames(modern_pollen) == "taxa0")
taxaColMax <- which(colnames(modern_pollen) == "taxaN")
taxa <- modern_pollen[, taxaColMin:taxaColMax]

# Load reconstruction data
Holocene <- read.csv("/path/to/Holocene.csv")
taxaColMin <- which(colnames(Holocene) == "taxa0")
taxaColMax <- which(colnames(Holocene) == "taxaN")
core <- Holocene[, taxaColMin:taxaColMax]

## Train
fit_Tmin <- fxTWAPLS::WAPLS.w(taxa, modern_pollen$Tmin, nPLS = 5)
fit_f_Tmin <- fxTWAPLS::WAPLS.w(
```
taxa, 
modern_pollen$Tmin,
nPLS = 5,
usefx = TRUE,
fx_method = "bin",
bin = 0.02 
)
fit_Tmin2 <- fxTWAPLS::WAPLS.w2(taxa, modern_pollen$Tmin, nPLS = 5)
fit_f_Tmin2 <- fxTWAPLS::WAPLS.w2(
taxa, 
modern_pollen$Tmin,
nPLS = 5,
usefx = TRUE,
fx_method = "bin",
bin = 0.02 
)
## Predict
fossil_Tmin <- fxTWAPLS::WAPLS.predict.w(fit_Tmin, core)
fossil_f_Tmin <- fxTWAPLS::WAPLS.predict.w(fit_f_Tmin, core)
fossil_Tmin2 <- fxTWAPLS::WAPLS.predict.w(fit_Tmin2, core)
fossil_f_Tmin2 <- fxTWAPLS::WAPLS.predict.w(fit_f_Tmin2, core)
## End(Not run)

WAPLS.w  
WA-PLS training function

Description

WA-PLS training function, which can perform fx correction. \( \frac{1}{f^2} \) correction will be applied at step 7.

Usage

WAPLS.w(
  modern_taxa, 
  modern_climate, 
  nPLS = 5, 
  usefx = FALSE, 
  fx_method = "bin", 
  bin = NA 
)

Arguments

modern_taxa    The modern taxa abundance data, each row represents a sampling site, each column represents a taxon.
modern_climate The modern climate value at each sampling site.
nPLS: The number of components to be extracted.
usefx: Boolean flag on whether or not use fx correction.
fx_method: Binned or p-spline smoothed fx correction: if usefx = FALSE, this should be NA; otherwise, fx function will be used when choosing "bin"; fx_pspline function will be used when choosing "pspline".
bin: Binwidth to get fx, needed for both binned and p-splined method. if usefx = FALSE, this should be NA;

Value

A list of the training results, which will be used by the predict function. Each element in the list is described below:

- **fit**: the fitted values using each number of components.
- **x**: the observed modern climate values.
- **taxon_name**: the name of each taxon.
- **optimum**: the updated taxon optimum (u* in the WA-PLS paper).
- **comp**: each component extracted (will be used in step 7 regression).
- **u**: taxon optimum for each component (step 2).
- **z**: a parameter used in standardization for each component (step 5).
- **s**: a parameter used in standardization for each component (step 5).
- **orth**: a list that stores orthogonalization parameters (step 4).
- **alpha**: a list that stores regression coefficients (step 7).
- **meanx**: mean value of the observed modern climate values.
- **nPLS**: the total number of components extracted.

See Also

fx, TWAPLS.w, and WAPLS.predict.w

Examples

```r
## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

# Extract taxa
taxaColMin <- which(colnames(moderan_pollen) == "taxa0")
taxaColMax <- which(colnames(moderan_pollen) == "taxaN")
taxa <- modern_pollen[, taxaColMin:taxaColMax]

# Training
fit_Tmin <- fxTWAPLS::WAPLS.w(taxa, modern_pollen$Tmin, nPLS = 5)
fit_f_Tmin <- fxTWAPLS::WAPLS.w(taxa, modern_pollen$Tmin, nPLS = 5)
```
nPLS = 5,
usefx = TRUE,
fx_method = "bin",
bin = 0.02
)

## End(Not run)

### Description

WA-PLS training function, which can perform fx correction. $1/fx$ correction will be applied at step 2 and step 7.

### Usage

```r
WAPLS.w2(
  modern_taxa,
  modern_climate,
  nPLS = 5,
  usefx = FALSE,
  fx_method = "bin",
  bin = NA
)
```

### Arguments

- **modern_taxa**: The modern taxa abundance data, each row represents a sampling site, each column represents a taxon.
- **modern_climate**: The modern climate value at each sampling site.
- **nPLS**: The number of components to be extracted.
- **usefx**: Boolean flag on whether or not use $fx$ correction.
- **fx_method**: Binned or p-spline smoothed $fx$ correction: if usefx = FALSE, this should be NA; otherwise, $fx$ function will be used when choosing "bin"; $fx_pspline$ function will be used when choosing "pspline".
- **bin**: Binwidth to get $fx$, needed for both binned and p-splined method. if usefx = FALSE, this should be NA;

### Value

A list of the training results, which will be used by the predict function. Each element in the list is described below:

- **fit**: the fitted values using each number of components.
x  the observed modern climate values.
taxon_name  the name of each taxon.
optimum  the updated taxon optimum (u* in the WA-PLS paper).
comp  each component extracted (will be used in step 7 regression).
u  taxon optimum for each component (step 2).
z  a parameter used in standardization for each component (step 5).
s  a parameter used in standardization for each component (step 5).
orth  a list that stores orthogonalization parameters (step 4).
alpha  a list that stores regression coefficients (step 7).
meanx  mean value of the observed modern climate values.
nPLS  the total number of components extracted.

See Also

fx, TWAPLS.w, and WAPLS.predict.w

Examples

```r
## Not run:
# Load modern pollen data
modern_pollen <- read.csv("/path/to/modern_pollen.csv")

# Extract taxa
taxaColMin <- which(colnames(morden_pollen) == "taxa0")
taxaColMax <- which(colnames(morden_pollen) == "taxaN")
taxa <- modern_pollen[, taxaColMin:taxisColMax]

# Training
fit_Tmin2 <- fxTWAPL::WAPLS.w2(taxa, modern_pollen$Tmin, nPLS = 5)
fit_f_Tmin2 <- fxTWAPL::WAPLS.w2(
taxa,
  modern_pollen$Tmin,
  nPLS = 5,
  usefx = TRUE,
  fx_method = "bin",
  bin = 0.02
)

## End(Not run)
```
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