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BugReports https://github.com/egarpor/goffda

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goffda-package

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goffda-package

gofda – Goodness-of-Fit Tests for Functional Data

Description


Author(s)

Eduardo García-Portugués and Javier Álvarez-Liébana.

References


Description


Usage

\texttt{aemet_temp}

Format

A list with the following entries:

\textbf{temp} an \texttt{fdata} with 2892 temperature (in Celsius degrees) curves, discretized on 365 equispaced grid points (days) on $[0.5, 364.5]$. Each curve corresponds to the yearly records of a weather station.

\textbf{df} a dataframe with metadata for each curve:

- \texttt{ind}: identifier of the weather station.
- \texttt{name}: name of the weather station.
- \texttt{year}: year of the observation.

Details

For consistency with the \texttt{fda.usc-package}'s \texttt{aemet} dataset, the names and identifiers of the 73 weather stations are the same as in that dataset. Only a minor fix has been applied to the "A CORUÑA/ALVEDRO" station, whose identifier was the same as the "A CORUÑA" station, "1387". The former was set to "1387A".

Information about the province, altitude, longitude, and latitude of each weather station can be retrieved in \texttt{df} from the \texttt{fda.usc-package}'s \texttt{aemet} dataset.

The dataset is a curated version of a larger database of 115 stations. It excludes stations with inconsistent records or that were relocated, closed, or opened during the 40-year period. There are 9 stations with missing years. The total of missing years is 28.

In leap years, the daily-average temperature is computed as the average of February 28th and 29th.

Author(s)

Original data processing scripts by Manuel Febrero-Bande and Manuel Oviedo de la Fuente. Adaptations by Eduardo García-Portugués.

Source

The data was retrieved from the FTP of the Meteorological State Agency of Spain (AEMET) in 2014 using a processing script by the authors of the \texttt{fda.usc-package}. 
References


Examples

```r
# Data splitting

# Load raw data
data("aemet_temp")

# Partition the dataset in the first and last 20 years
with(aemet_temp, {
  ind_pred <- which((1974 <= df$year) & (df$year <= 1993))
  ind_resp <- which((1994 <= df$year) & (df$year <= 2013))
  aemet_temp_pred <<- list("df" = df[ind_pred, ], "temp" = temp[ind_pred])
  aemet_temp_resp <<- list("df" = df[ind_resp, ], "temp" = temp[ind_resp])
})

# Average the temperature on each period
mean_aemet <- function(x) {
  m <- tapply(X = 1:nrow(x$temp$data), INDEX = x$df$ind,
    FUN = function(i) colMeans(x$temp$data[i, , drop = FALSE],
    na.rm = TRUE))
  x$temp$data <- do.call(rbind, m)
  return(x$temp)
}

# Build predictor and response fdatas
aemet_temp_pred <- mean_aemet(aemet_temp_pred)
aemet_tempresp <- mean_aemet(aemet_temp_resp)

# Plot
old_par <- par(mfrow = c(1, 2))
plot(aemet_temp_pred)
plot(aemet_temp_resp)
par(old_par)

# Average daily temperatures
day_avg_pred <- func_mean(aemet_temp_pred)
day_avg_resp <- func_mean(aemet_temp_resp)

# Average yearly temperatures
avg_year_pred <- rowMeans(aemet_temp_pred$data)
avg_year_resp <- rowMeans(aemet_temp_resp$data)

# Test the linear model with functional response and predictor
(comp_flmfr <- flm_test(X = aemet_temp_pred, Y = aemet_temp_resp,
est_method = "fpcr_l1s")
beta0 <- diag(rep(1, length(aemet_temp_pred$argvals)))
```
cv_glmnet

Fitting of regularized linear models

description

Convenience function for fitting multivariate linear models with multivariate response by relying on cv.glmnet from the glmnet-package. The function fits the multivariate linear model

\[ \mathbf{Y} = \mathbf{X}\mathbf{B} + \mathbf{E}, \]

where \( \mathbf{X} \) is a \( p \)-dimensional vector, \( \mathbf{Y} \) and \( \mathbf{E} \) are two \( q \)-dimensional vectors, and \( \mathbf{B} \) is a \( p \times q \) matrix. If \( \mathbf{X} \) and \( \mathbf{Y} \) are centered (i.e., have zero-mean columns), the function estimates \( \mathbf{B} \) by solving, for the sample \( \langle \mathbf{X}_1, \mathbf{Y}_1 \rangle, \ldots, \langle \mathbf{X}_n, \mathbf{Y}_n \rangle \), the elastic-net optimization problem

\[
\min_{\mathbf{B} \in \mathbb{R}^{p \times q}} \frac{1}{2n} \sum_{i=1}^{n} \| \mathbf{Y}_i - \mathbf{X}_i\mathbf{B} \|^2 + \lambda \left[ (1 - \alpha)\|\mathbf{B}\|_F^2 / 2 + \alpha \sum_{j=1}^{p} \| \mathbf{B}_j \|_2 \right],
\]

where \( \|\mathbf{B}\|_F \) stands for the Frobenious norm of the matrix \( \mathbf{B} \) and \( \| \mathbf{B}_j \|_2 \) for the Euclidean norm of the \( j \)-th row of \( \mathbf{B} \). The choice \( \alpha = 0 \) in the elastic-net penalization corresponds to ridge regression,
whereas $\alpha = 1$ yields a lasso-type estimator. The unpenalized least-squares estimator is obtained with $\lambda = 0$.

**Usage**

```r
cv_glmnet(x, y, alpha = c("lasso", "ridge")[1], lambda = NULL,
  intercept = TRUE, thresh = 1e-10, cv_1se = TRUE, cv_nlambda = 50,
  cv_folds = NULL, cv_grouped = FALSE, cv_lambda = 10^seq(2, -3,
    length.out = cv_nlambda), cv_second = TRUE, cv_tol_second = 0.025,
  cv_log10_exp = c(-0.5, 3), cv_thresh = 1e-05, cv_parallel = FALSE,
  cv_verbose = FALSE, ...)
```

**Arguments**

- **x**
  - input matrix of size $c(n,p)$, or a vector of length $n$.
- **y**
  - response matrix of size $c(n,q)$, or a vector of length $n$.
- **alpha**
  - elastic net mixing argument in `glmnet`, with $0 \leq \alpha \leq 1$. Alternatively, a character string indicating whether the "ridge" ($\alpha = 0$) or "lasso" ($\alpha = 1$) fit is to be performed.
- **lambda**
  - scalar giving the regularization parameter $\lambda$. If NULL (default), the optimal $\lambda$ is searched by cross-validation. If lambda is provided, then cross-validation is skipped and the fit is performed for the given lambda.
- **intercept**
  - flag passed to the intercept argument in `glmnet` to indicate if the intercept should be fitted (default; does not assume that the data is centered) or set to zero (the optimization problem above is solved as-is). Defaults to TRUE.
- **thresh**
  - convergence threshold of the coordinate descending algorithm, passed to the thresh argument in `glmnet`. Defaults to 1e-10.
- **cv_1se**
  - shall the optimal lambda be the lambda.1se, as returned by `cv.glmnet`? This favors sparser fits. If FALSE, then the optimal lambda is lambda.min, the minimizer of the cross-validation loss. Defaults to TRUE.
- **cv_nlambda**
  - the length of the sequence of $\lambda$ values, passed to the nlambda argument in `cv.glmnet` for the cross-validation search. Defaults to 50.
- **cv_folds**
  - number of folds to perform cross-validation. If NULL (the default), then cv_folds <- n internally, that is, leave-one-out cross-validation is performed. Passed to the nfolds argument in `cv.glmnet`.
- **cv_grouped**
  - passed to the grouped argument in `cv.glmnet`. Defaults to FALSE.
- **cv_lambda**
  - passed to the lambda argument in `cv.glmnet`. Defaults to $10^{seq(2,-3, length.out = cv_nlambda)}$.
- **cv_second**
  - flag to perform a second cross-validation search if the optimal $\lambda$ was found at the extremes of the first $\lambda$ sequence (indicating that the minimum may not be reliable). Defaults to TRUE.
- **cv_tol_second**
  - tolerance for performing a second search if second = TRUE. If the minimum is found at the $100 \times cv_tol_second$ lower/upper percentile of search interval, then the search interval is expanded for a second search. Defaults to 0.025.
cv_glmnet

- **cv_log10_exp**: expansion of the $\lambda$ sequence if the minimum is found close to its upper extreme. If that is the case, the sequence for the is set as $10^{\log_{10}(\lambda_{\text{min}}) + \text{cv_log10_exp}}$, where $\lambda_{\text{min}}$ is the minimum obtained in the first sequence. If the minimum is found close to the lower extreme of the sequence, then $-\text{rev}(\text{cv_log10_exp})$ is considered. Defaults to $c(-0.5, 5)$.

- **cv_thresh**: convergence threshold used during cross-validation in cv.glmnet. Defaults to $1e^{-5}$.

- **cv_parallel**: passed to the parallel argument in cv.glmnet. Defaults to FALSE.

- **cv_verbose**: flag to display information about the cross-validation search with plots and messages. More useful if second = TRUE. Defaults to FALSE.

- ... further parameters to be passed to glmnet to perform the final model fit.

**Details**

If $\alpha = 1$, then the lasso-type fit shrinks to zero, simultaneously, all the elements of certain rows of $B$, thus helping the selection of the $p$ most influential variables in $X$ for explaining/predicting $Y$.

The function first performs a cross-validation search for the optimal $\lambda$ if lambda = NULL (using cv_thresh to control the convergence threshold). After the optimal penalization parameter is determined, a second fit (now with convergence threshold thresh) using the default $\lambda$ sequence in glmnet is performed. The final estimate is obtained via predict.glmnet from the optimal $\lambda$ determined in the first step.

Due to its cross-validatory nature, cv_glmnet can be computationally demanding. Approaches for reducing the computation time include: considering a smaller number of folds than $n$, such as cv_folds = 10 (but will lead to random partitions of the data); decrease the tolerance of the coordinate descending algorithm by increasing cv_thresh; reducing the number of candidate $\lambda$ values with nlambda; setting second = FALSE to avoid a second cross-validation; or considering cv_parallel = TRUE to use a parallel backend (must be registered before hand; see examples).

By default, the $\lambda$ sequence is used with standardized $X$ and $Y$ (both divided by their columnwise variances); see glmnet and the standardized argument. Therefore, the optimal selected $\lambda$ value assumes standardization and must be used with care if the variables are not standardized. For example, when using the ridge analytical solution, a prior change of scale that depends on $q$ needs to be done. See the examples for the details.

**Value**

A list with the following entries:

- **beta_hat**: the estimated $B$, a matrix of size $c(p, q)$.
- **lambda**: the optimal $\lambda$ obtained by cross-validation and according to cv_1se.
- **cv**: if lambda = NULL, the result of the cross-validation search for the optimal $\lambda$. Otherwise, NULL.
- **fit**: the glmnet fit, computed with thresh threshold and with an automatically chosen $\lambda$ sequence.

**Author(s)**

Eduardo García-Portugués. Initial contributions by Gonzalo Álvarez-Pérez.
References


Examples

```r
## Quick example for multivariate linear model with multivariate response

# Simulate data
n <- 100
p <- 10; q <- 5
set.seed(123456)
x <- matrix(rnorm(n * p, sd = rep(1:p, each = n)), nrow = n, ncol = p)
e <- matrix(rnorm(n * q, sd = rep(q:1, each = n)), nrow = n, ncol = q)
beta <- matrix(((1:p - 1) / p)^2, nrow = p, ncol = q)
y <- x %*% beta + e

# Fit lasso (model with intercept, the default)
cv_glmnet(x = x, y = y, alpha = "lasso", cv_verbose = TRUE)$beta_hat

## Multivariate linear model with multivariate response

# Simulate data
n <- 100
p <- 10; q <- 5
set.seed(123456)
x <- matrix(rnorm(n * p, sd = rep(1:p, each = n)), nrow = n, ncol = p)
e <- matrix(rnorm(n * q, sd = rep(q:1, each = n)), nrow = n, ncol = q)
beta <- matrix(((1:p - 1) / p)^2, nrow = p, ncol = q)
y <- x %*% beta + e

# Fit ridge
cv0 <- cv_glmnet(x = x, y = y, alpha = "ridge", intercept = FALSE, cvVerbose = TRUE)
cv0$beta_hat

# Same fit for the chosen lambda
cv_glmnet(x = x, y = y, alpha = "ridge", lambda = cv0$lambda, intercept = FALSE)$beta_hat

# Fit lasso (model with intercept, the default)
cv1 <- cv_glmnet(x = x, y = y, alpha = "lasso", cvVerbose = TRUE)
cv1$beta_hat

# Use cv_1se = FALSE
cv1_min <- cv_glmnet(x = x, y = y, alpha = "lasso", cvVerbose = TRUE, cv_1se = FALSE)
cv1$beta_hat

# Compare it with ridge analytical solution. Observe the factor in lambda,
# necessary since lambda is searched for the standardized data. Note also
# that, differently to the case q = 1, no standardization with respect to
```
# y happens
sd_x <- apply(x, 2, function(x) sd(x)) * sqrt((n - 1) / n)
cv_glmnet(x = x, y = y, alpha = "ridge", lambda = cv0$lambda,
    thresh = 1e-20, intercept = FALSE)$beta_hat
solve(crossprod(x) + diag(cv0$lambda * sd_x^2 * n)) %*% t(x) %*% y

# If x is standardized, the match between glmnet and usual ridge
# analytical expression does not require scaling of lambda
x_std <- scale(x, scale = sd_x, center = TRUE)
cv_glmnet(x = x_std, y = y, alpha = "ridge", lambda = cv0$lambda,
    intercept = FALSE, thresh = 1e-20)$beta_hat
solve(crossprod(x_std) + diag(rep(cv0$lambda * n, p))) %*% t(x_std) %*% y

## Simple linear model

# Simulate data
n <- 100
p <- 1; q <- 1
set.seed(123456)
x <- matrix(rnorm(n * p), nrow = n, ncol = p)
e <- matrix(rnorm(n * q), nrow = n, ncol = q)
beta <- 2
y <- x * beta + e

# Fit by ridge (model with intercept, the default)
cv0 <- cv_glmnet(x = x, y = y, alpha = "ridge", cv_verbose = TRUE)
cv0$beta_hat

cv0$intercept

# Comparison with linear model with intercept
lm(y ~ 1 + x)$coefficients

# Fit by ridge (model without intercept)
cv0 <- cv_glmnet(x = x, y = y, alpha = "ridge", cv_verbose = TRUE,
    intercept = FALSE)
cv0$beta_hat

# Comparison with linear model without intercept
lm(y ~ 0 + x)$coefficients

# Same fit for the chosen lambda (and without intercept)
cv_glmnet(x = x, y = y, alpha = "ridge", lambda = cv0$lambda,
    intercept = FALSE)$beta_hat

# Same for lasso (model with intercept, the default)
cv1 <- cv_glmnet(x = x, y = y, alpha = "lasso")
cv1$beta_hat

## Multivariate linear model (p = 3, q = 1)

# Simulate data
n <- 50
p <- 10; q <- 1
set.seed(123456)
x <- matrix(rnorm(n * p, mean = 1, sd = rep(1:p, each = n)),
          nrow = n, ncol = p)
e <- matrix(rnorm(n * q), nrow = n, ncol = q)
beta <- ((1:p - 1) / p)^2
y <- x %*% beta + e

# Fit ridge (model without intercept)
cv0 <- cv_glmnet(x = x, y = y, alpha = "ridge", intercept = FALSE,
                  cv_verbose = TRUE)
cv0$beta_hat

# Same fit for the chosen lambda
cv_glmnet(x = x, y = y, alpha = "ridge", lambda = cv0$lambda,
          intercept = FALSE)$beta_hat

# Compare it with ridge analytical solution. Observe the factor in lambda,
# necessary since lambda is searched for the standardized data
sd_x <- apply(x, 2, function(x) sd(x)) * sqrt((n - 1) / n)
sd_y <- sd(y) * sqrt((n - 1) / n)
cv_glmnet(x = x, y = y, alpha = "ridge", lambda = cv0$lambda,
          intercept = FALSE, thresh = 1e-20)$beta_hat
solve(crossprod(x) + diag(cv0$lambda * sd_x^2 / sd_y * n)) %*% t(x) %*% y

# If x and y are standardized, the match between glmnet and usual ridge
# analytical expression does not require scaling of lambda
x_std <- scale(x, scale = sd_x, center = TRUE)
y_std <- scale(y, scale = sd_y, center = TRUE)
cv_glmnet(x = x_std, y = y_std, alpha = "ridge", lambda = cv0$lambda,
          intercept = FALSE, thresh = 1e-20)$beta_hat
solve(crossprod(x_std) + diag(rep(cv0$lambda * n, p))) %*% t(x_std) %*% y_std

# Fit lasso (model with intercept, the default)
cv1 <- cv_glmnet(x = x, y = y, alpha = "lasso", cv_verbose = TRUE)
cv1$beta_hat

# Parallelization
#
# Parallel
# doMC::registerDoMC(cores = 2)
# microbenchmark::microbenchmark(
#   cv_glmnet(x = x, y = y, nlambda = 100, cv_parallel = TRUE),
#   cv_glmnet(x = x, y = y, nlambda = 100, cv_parallel = FALSE),
#   times = 10)

Covariate, error, and kernel of a functional linear model with functional response
Description

Simulation of \( X \), a random variable in the Hilbert space of square-integrable functions in \([a, b]\), \( L^2([a, b]) \), and \( \varepsilon \), a random variable in \( L^2([c, d]) \). Together with the bivariate kernel \( \beta \), they are the necessary elements for sampling a Functional Linear Model with Functional Response (FLMFR):

\[
Y(t) = \int_a^b X(s) \beta(s,t) ds + \varepsilon(t).
\]

The next functions sample \( X \) and \( \varepsilon \), and construct \( \beta \), using different proposals in the literature:

- \( r_{cm2013_flmfr} \) is based on the numerical example given in Section 3 of Crambes and Mas (2013). Termed as S1 in Section 2 of García-Portugués et al. (2021).
- \( r_{ik2018_flmfr} \) is based on the numerical example given in Section 4 of Imaizumi and Kato (2018), but zeroing the first Functional Principal Components (FPC) coefficients of \( \beta \) (so the first FPC are not adequate for estimation). S3 in Section 2 of García-Portugués et al. (2021).
- \( r_{gof2021_flmfr} \) gives a numerical example in Section 2 of García-Portugués et al. (2021), denoted therein as S2.

Usage

\[
r_{cm2013_flmfr}(n, s = \text{seq}(0, 1, \text{len} = 101), t = \text{seq}(0, 1, \text{len} = 101), \text{std_error} = 0.15, n_{\text{fpc}} = 50, \text{concurrent} = \text{FALSE})
\]

\[
r_{ik2018_flmfr}(n, s = \text{seq}(0, 1, l = 101), t = \text{seq}(0, 1, l = 101), \text{std_error} = 1.5, \text{parameters} = \text{c}(1.75, 0.8, 2.4, 0.25), n_{\text{fpc}} = 50, \text{concurrent} = \text{FALSE})
\]

\[
r_{gof2021_flmfr}(n, s = \text{seq}(0, 1, \text{len} = 101), t = \text{seq}(0, 1, \text{len} = 101), \text{std_error} = 0.35, \text{concurrent} = \text{FALSE})
\]

Arguments

- \( n \) number of trajectories to sample.
- \( s, t \) grid points where functional covariates and responses are valued, respectively.
- \( \text{std_error} \) standard deviation of the random variables involved in the generation of the functional error \( \text{error_fdata} \). Defaults to 0.15.
- \( n_{\text{fpc}} \) number of FPC to be taken into account for the data generation. Must be greater than 4 when \( r_{ik2018_flmfr} \) is applied, since the first 4 FPC are null. Defaults to 50.
- \( \text{concurrent} \) flag to consider a concurrent FLMFR (degenerate case). Defaults to FALSE.
- \( \text{parameters} \) vector of parameters, only required for \( r_{ik2018_flmfr} \). Defaults to \( \text{c}(1.75, 0.8, 2.4, 0.25) \).

Details

Descriptions of the processes \( X \) and \( \varepsilon \), and of \( \beta \) can be seen in the references.
Value

A list with the following elements:

- X_fdata: functional covariates, an fdata object of length n.
- error_fdata: functional errors, an fdata object of length n.
- beta: either the matrix with $\beta(s, t)$ evaluated at the argvals of X_fdata and Y_fdata (if concurrent = FALSE) or a vector with $\beta(t)$ evaluated at the argvals of X_fdata (if concurrent = TRUE).

Author(s)

Javier Álvarez-Liébana.

References


Examples

# FLMFR based on Imaizumi and Kato (2018) adopting different Hilbert spaces
s <- seq(0, 1, l = 201)
t <- seq(2, 4, l = 301)
r_ik2018 <- r_ik2018_flmfr(n = 50, s = s, t = t, std_error = 1.5,
parameters = c(1.75, 0.8, 2.4, 0.25), n_fpc = 50)
plot(r_ik2018$X_fdata)
plot(r_ik2018$error_fdata)
image(x = s, y = t, z = r_ik2018$beta, col = viridisLite::viridis(20))

# FLMFR based on Cardot and Mas (2013) adopting different Hilbert spaces
r_cm2013 <- r_cm2013_flmfr(n = 50, s = s, t = t, std_error = 0.15,
n_fpc = 50)
plot(r_cm2013$X_fdata)
plot(r_cm2013$error_fdata)
image(x = s, y = t, z = r_cm2013$beta, col = viridisLite::viridis(20))

# FLMFR in García-Portugués et al. (2021) adopting different Hilbert spaces
r_gof2021 <- r_gof2021_flmfr(n = 50, s = s, t = t, std_error = 0.35,
concurrent = FALSE)
plot(r_gof2021$X_fdata)
plot(r_gof2021$error_fdata)
image(x = s, y = t, z = r_gof2021$beta, col = viridisLite::viridis(20))

# Concurrent model in García-Portugués et al. (2021)
r_gof2021 <- r_gof2021_flmfr(n = 50, s = s, t = s, std_error = 0.35,
flm_est

Estimation of functional linear models

Description

Estimation of the linear operator relating a functional predictor \( X \) with a functional response \( Y \) in the linear model

\[
Y(t) = \int_a^b \beta(s, t) X(s) \, ds + \varepsilon(t),
\]

where \( X \) is a random variable in the Hilbert space of square-integrable functions in \([a, b]\), \( L^2([a, b])\), \( Y \) and \( \varepsilon \) are random variables in \( L^2([c, d]) \), and \( s \in [a, b] \) and \( t \in [c, d] \).

The linear, Hilbert–Schmidt, integral operator is parametrized by the bivariate kernel \( \beta \in L^2([a, b] \otimes L^2([c, d])) \). Its estimation is done through the truncated expansion of \( \beta \) in the tensor product of the data-driven bases of the Functional Principal Components (FPC) of \( X \) and \( Y \), and through the fitting of the resulting multivariate linear model. The FPC basis for \( X \) is truncated in \( p \) components, while the FPC basis for \( Y \) is truncated in \( q \) components. Automatic selection of \( p \) and \( q \) is detailed below.

The particular cases in which either \( X \) or \( Y \) are constant functions give either a scalar predictor or response. The simple linear model arises if both \( X \) and \( Y \) are scalar, for which \( \beta \) is a constant.

Usage

\[
\text{flm_est}(X, Y, \text{est_method} = \"fpcr_l1s\", p = \text{NULL}, q = \text{NULL},
\text{thre_p} = 0.99, \text{thre_q} = 0.99, \text{lambda} = \text{NULL}, \text{X_fpc} = \text{NULL},
\text{Y_fpc} = \text{NULL}, \text{compute_residuals} = \text{TRUE}, \text{centered} = \text{FALSE},
\text{int_rule} = \text{\"trapezoid\"}, \text{cv_verbose} = \text{FALSE}, \ldots)
\]

Arguments

\( X, Y \)  
samples of functional/scalar predictors and functional/scalar response. Either \textit{fdata} objects (for functional variables) or vectors of length \( n \) (for scalar variables).

\textit{est_method}  
either \"fpcr\" (Functional Principal Components Regression; FPCR), \"fpcr_l2\" (FPCR with ridge penalty), \"fpcr_l1\" (FPCR with lasso penalty) or \"fpcr_l1s\" (FPCR with lasso-selected FPC). If \( X \) is scalar, \textit{flm_est} only considers \"fpcr\" as estimation method. See details below. Defaults to \"fpcr_l1s\".

\( p, q \)  
index vectors indicating the specific FPC to be considered for the truncated bases expansions of \( X \) and \( Y \), respectively. If a single number for \( p \) is provided, then \( p <= 1 \times \max(p) \) internally (analogously for \( q \)) and the first \( \max(p) \) FPC are considered. If \text{NULL} (default), then a data-driven selection of \( p \) and \( q \) is done. See details below.
flm_est

thre_p, thre_q  thresholds for the proportion of variance that is explained, at least, by the first $p$ and $q$ FPC of $X$ and $Y$, respectively. These thresholds are employed for an (initial) automatic selection of $p$ and $q$. Default to 0.99. thre_p (thre_q) is ignored if $p$ ($q$) is provided.

lambda  regularization parameter $\lambda$ for the estimation methods "fpcr_l2", "fpcr_l1", and "fpcr_l1s". If NULL (default), it is chosen with cv_glmnet.

X_fpc, Y_fpc  FPC decompositions of $X$ and $Y$, as returned by fpc. Computed if not provided.

compute_residuals  whether to compute the fitted values $Y_{\hat{}}$ and its $Y_{\hat{}\text{scores}}$, and the residuals of the fit and its residuals_scores. Defaults to TRUE.

centered  flag to indicate if $X$ and $Y$ have been centered or not, in order to avoid their recentering. Defaults to FALSE.

int_rule  quadrature rule for approximating the definite unidimensional integral: trapezoidal rule (int_rule = "trapezoid") and extended Simpson rule (int_rule = "Simpson") are available. Defaults to "trapezoid".

cv_verbose  flag to display information about the estimation procedure (passed to cv_glmnet). Defaults to FALSE.

...  further parameters to be passed to cv_glmnet (and then to cv.glmnet) such as cv_1se, cv_nlambda or cv_parallel, among others.

Details

flm_est deals seamlessly with either functional or scalar inputs for the predictor and response. In the case of scalar inputs, the corresponding dimension-related arguments ($p$, $q$, thre_p or thre_q) will be ignored as in these cases either $p = 1$ or $q = 1$.

The function translates the functional linear model into a multivariate model with multivariate response and then estimates the $p \times q$ matrix of coefficients of $\beta$ in the tensor basis of the FPC of $X$ and $Y$. The following estimation methods are implemented:

- "fpcr": Functional Principal Components Regression (FPCR); see details in Ramsay and Silverman (2005).
- "fpcr_l2": FPCR, with ridge penalty on the associated multivariate linear model.
- "fpcr_l1": FPCR, with lasso penalty on the associated multivariate linear model.
- "fpcr_l1s": FPCR, with FPC selected by lasso regression on the associated multivariate linear model.

The last three methods are explained in García-Portugués et al. (2021).

The $p$ FPC of $X$ and $q$ FPC of $Y$ are determined as follows:

- If $p = \text{NULL}$, then $p$ is set as $\text{p\_thre} <1: \text{j\_thre}$, where $\text{j\_thre}$ is the $j$-th FPC of $X$ for which the cumulated proportion of explained variance is greater than thre_p. If $p \neq \text{NULL}$, then $\text{p\_thre} <\text{p}$.
- If $q = \text{NULL}$, then the same procedure is followed with thre_q, resulting q_thre.
Once \( p_{\text{thre}} \) and \( q_{\text{thre}} \) have been obtained, the methods "fpcr_l1" and "fpcr_l1s" perform a second selection of the FPC that are effectively considered in the estimation of \( \beta \). This subset of FPC (of \( p_{\text{thre}} \)) is encoded in \( p_{\text{hat}} \). No further selection of FPC is done for the methods "fpcr" and "fpcr_l2".

The flag `compute_residuals` controls if \( Y_{\text{hat}}, Y_{\text{hat}}_{\text{scores}}, \text{residuals}, \text{and residuals}\_scores \) are computed. If \texttt{FALSE}, they are set to \texttt{NULL}. \( Y_{\text{hat}} \) equals \( \hat{Y}_i(t) = \int_a^b \hat{\beta}(s,t)X_i(s)\,ds \) and \( \text{residuals} \) stands for \( \hat{\varepsilon}_i(t) = Y_i(t) - \hat{Y}_i(t) \), both for \( i = 1, \ldots, n \). \( Y_{\text{hat}}_{\text{scores}} \) and \( \text{residuals}\_scores \) are the \( n \times q \) matrices of coefficients (or scores) of these functions in the FPC of \( Y \).

Missing values on \( X \) and \( Y \) are automatically removed.

**Value**

A list with the following entries:

- \textbf{Beta\_hat} estimated \( \beta \), a matrix with values \( \hat{\beta}(s,t) \) evaluated at the grid points for \( X \) and \( Y \). Its size is \( c(\text{length}(X$argvals)\),\text{length}(Y$argvals)) \).
- \textbf{Beta\_hat\_scores} the matrix of coefficients of \( \text{Beta\_hat} \) (resulting from projecting it into the tensor basis of \( X_{\text{fpc}} \) and \( Y_{\text{fpc}} \)), with dimension \( c(p_{\text{hat}},q_{\text{thre}}) \).
- \textbf{H\_hat} hat matrix of the associated fitted multivariate linear model, a matrix of size \( c(n,n) \). \texttt{NULL} if \texttt{est\_method = "fpcr\_l1"}, since lasso estimation does not provide it explicitly.
- \textbf{p\_thre} index vector indicating the FPC of \( X \) considered for estimating the model. Chosen by \texttt{thre\_p} or equal to \( p \) if given.
- \textbf{p\_hat} index vector of the FPC considered by the methods "fpcr\_l1" and "fpcr\_l1s" methods after further selection of the FPC considered in \( p_{\text{thre}} \). For methods "fpcr" and "fpcr\_l2", \( p_{\text{hat}} \) equals \( p_{\text{thre}} \).
- \textbf{q\_thre} index vector indicating the FPC of \( Y \) considered for estimating the model. Chosen by \texttt{thre\_q} or equal to \( q \) if given. Note that zeroing by lasso procedure only affects in the rows.
- \textbf{est\_method} the estimation method employed.
- \textbf{Y\_hat} fitted values, either an \texttt{fdata} object or a vector, depending on \( Y \).
- \textbf{Y\_hat\_scores} the matrix of coefficients of \( \text{Y\_hat} \), with dimension \( c(n,q_{\text{thre}}) \).
- \textbf{residuals} residuals of the fitted model, either an \texttt{fdata} object or a vector, depending on \( Y \).
- \textbf{residuals\_scores} the matrix of coefficients of \( \text{residuals} \), with dimension \( c(n,q_{\text{thre}}) \).
- \textbf{X\_fpc, Y\_fpc} FPC of \( X \) and \( Y \), as returned by \texttt{fpc} with \( n_{\text{fpc}} = n \).
- \textbf{lambda} regularization parameter \( \lambda \) used for the estimation methods "fpcr\_l2", "fpcr\_l1", and "fpcr\_l1s".
- \textbf{cv} cross-validation object returned by \texttt{cv\_glmnet}.

**Author(s)**

Eduardo García-Portugués and Javier Álvarez-Liébana.
## Quick example of functional response and functional predictor

```r
# Generate data
set.seed(12345)
n <- 50
X_fdata <- r_ou(n = n, t = seq(0, 1, l = 201), sigma = 2)
epsilon <- r_ou(n = n, t = seq(0, 1, l = 201), sigma = 0.5)
Y_fdata <- 2 * X_fdata + epsilon

# Lasso-selection FPCR (p and q are estimated)
flm_est(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l1s")
```

## Functional response and functional predictor

```r
# Generate data
set.seed(12345)
n <- 50
X_fdata <- r_ou(n = n, t = seq(0, 1, l = 201), sigma = 2)
epsilon <- r_ou(n = n, t = seq(0, 1, l = 201), sigma = 0.5)
Y_fdata <- 2 * X_fdata + epsilon

# FPCR (p and q are estimated)
fpcr_1 <- flm_est(X = X_fdata, Y = Y_fdata, est_method = "fpcr")
fpcr_1$Beta_hat_scores
fpcr_1$p_thre
fpcr_1$q_thre

# FPCR (p and q are provided)
fpcr_2 <- flm_est(X = X_fdata, Y = Y_fdata, est_method = "fpcr",
p = c(1, 5, 2, 7), q = 2:1)
fpcr_2$Beta_hat_scores
fpcr_2$p_thre
fpcr_2$q_thre

# Ridge FPCR (p and q are estimated)
l2_1 <- flm_est(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l2")
l2_1$Beta_hat_scores
l2_1$p_hat

# Ridge FPCR (p and q are provided)
l2_2 <- flm_est(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l2",
p = c(1, 5, 2, 7), q = 2:1)
l2_2$Beta_hat_scores
```

### References


```r
l2_2$p_hat

# Lasso FPCR (p and q are estimated)
l1_1 <- flm_est(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l1")
l1_1$Beta_hat_scores
l1_1$p_thre
l1_1$p_hat

# Lasso estimator (p and q are provided)
l1_2 <- flm_est(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l1",
               p = c(1, 5, 2, 7), q = 2:1)
l1_2$Beta_hat_scores
l1_2$p_thre
l1_2$p_hat

# Lasso-selection FPCR (p and q are estimated)
l1s_1 <- flm_est(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l1s")
l1s_1$Beta_hat_scores
l1s_1$p_thre
l1s_1$p_hat

# Lasso-selection FPCR (p and q are provided)
l1s_2 <- flm_est(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l1s",
               p = c(1, 5, 2, 7), q = 1:4)
l1s_2$Beta_hat_scores
l1s_2$p_thre
l1s_2$p_hat

## Scalar response

# Generate data
set.seed(12345)
n <- 50
beta <- r_ou(n = 1, t = seq(0, 1, l = 201), sigma = 0.5, x0 = 3)
X_fdata <- fdata_cen(r_ou(n = n, t = seq(0, 1, l = 201), sigma = 2))
epsilon <- rnorm(n, sd = 0.25)
Y <- drop(inprod_fdata(X_fdata1 = X_fdata, X_fdata2 = beta)) + epsilon

# FPCR
fpcr_4 <- flm_est(X = X_fdata, Y = Y, est_method = "fpcr")
fpcr_4$p_hat

# Ridge FPCR
l2_4 <- flm_est(X = X_fdata, Y = Y, est_method = "fpcr_l2")
l2_4$p_hat

# Lasso FPCR
l1_4 <- flm_est(X = X_fdata, Y = Y, est_method = "fpcr_l1")
l1_4$p_hat

# Lasso-selection FPCR
l1s_4 <- flm_est(X = X_fdata, Y = Y, est_method = "fpcr_l1s")
l1s_4$p_hat
```

## Scalar predictor

# Generate data
set.seed(12345)
n <- 50
X <- rnorm(n)
epsilon <- r_ou(n = n, t = seq(0, 1, l = 201), sigma = 0.5)
beta <- r_ou(n = 1, t = seq(0, 1, l = 201), sigma = 0.5, x0 = 3)
beta$data <- matrix(beta$data, nrow = n, ncol = ncol(beta$data), byrow = TRUE)
Y_fdata <- beta * X + epsilon

# FPCR
fpcr_4 <- flm_est(X = X, Y = Y_fdata, est_method = "fpcr")
plot(beta, col = 2)
lines(beta$argvals, drop(fpcr_4$Beta_hat))

---

**flm_stat**

*Projected Cramér–von Mises test statistic for the goodness-of-fit test of functional linear models*

### Description

Computation of the Projected Cramér–von Mises (PCvM) test statistic and its associated $\mathbf{A}_\star$ matrix. For a sample of functional covariates $X_1, \ldots, X_n$, the test statistic is computed from $x_{1,p}, \ldots, x_{n,p}$, the coefficients (scores) of the sample in a $p$-truncated basis expansion, such as Functional Principal Components (FPC).

The PCvM statistic is defined as

$$ PCvM_{n,p,q} = c \cdot \text{tr}(\hat{\mathbf{E}}_q^\prime \mathbf{A}_\star \hat{\mathbf{E}}_q) $$

where

$$ c = 2\pi^{(p+q)/2-1}/(q\Gamma(p/2)\Gamma(q/2)n^2), $$

$\hat{\mathbf{E}}_q$ is the $n \times q$ matrix of multivariate residuals, and $\mathbf{A}_\star$ is a $n \times n$ matrix whose $ij$-th element is $\sum_{r=1}^n A_{ijr}$, for $A_{ijr}$ depending on $(x_{i,p}, x_{j,p}, x_{r,p})$. Its exact expression can be seen in Escanciano (2006) and García-Portugués et al. (2021).

### Usage

```r
flm_stat(E, p, Adot_vec, constant = TRUE)
Adot(X)
```
**Arguments**

- **E**
  - the matrix of multivariate residuals, with dimension $c(n,q)$. A vector if $q = 1$.
- **p**
  - dimension of the covariates space. Must be a positive integer.
- **Adot_vec**
  - output from `Adot`. A vector of length $n * (n -1) / 2 + 1$. This corresponds to the most expensive computation in the test statistic.
- **constant**
  - whether to include the constant of the PCvM test statistic, $c$, in its computation. Defaults to TRUE.
- **X**
  - a matrix of size $c(n,p)$ containing the coefficients (scores) of the functional data in a p-truncated orthonormal basis expansion, such as FPC. Must not contain repeated rows.

**Details**

`Adot` assumes that $X$ does not have repeated rows or otherwise NaNs will be present in the result. If $X$ has repeated rows, `Adot` will throw a warning.

The implementation of the PCvM test statistic for scalar response is addressed in García-Portugués et al. (2014), whereas García-Portugués et al. (2021) presents its multivariate extension and shows that $A \cdot$ induces a weighted quadratic norm (if there are no repetitions in the sample). The PCvM statistic is rooted in the proposal by Escanciano (2006).

Both `flm_stat` and `A_dot` are coded in C++.

**Value**

- **flm_stat**: the value of the test statistic, a scalar.
- **A_dot**: a vector of length $n * (n -1) / 2 + 1$. The first entry contains the common diagonal element of $A \cdot$. The remaining entries are the upper triangular matrix (excluding the diagonal) of $A \cdot$, stacked by columns.

**Author(s)**

Eduardo García-Portugués.

**References**


Examples

```r
## flm_stat

# Generate data
n <- 200
q <- 2
p <- 3
E <- matrix(rnorm(n * q), nrow = n, ncol = q)
X_fdata <- r_ou(n = n, t = seq(0, 1, l = 101))

# Compute FPC
X_fpc <- fpc(X_fdata)

# Adot
Adot_vec <- Adot(X = X_fpc["scores"])

# Check equality
constant <- n^(-2) * 2 * pi^((p / 2) - 1) / gamma(p / 2)
constant * .Fortran("pcvm_statistic", n = as.integer(n),
    Adot_vec = Adot_vec, residuals = E[, 2],
    statistic = 0)$statistic
flm_stat(E = E[, 2, drop = FALSE], p = p, Adot_vec = Adot_vec,
    constant = FALSE)

## Adot

# Generate data
n <- 200
X_fdata <- r_ou(n = n, t = seq(0, 1, l = 101))

# Compute FPC
X_fpc <- fpc(X_fdata)

# Using inprod_fdata and Adot
Adot_vec <- Adot(X = X_fpc["scores"])

# Check with fda.usc::Adot with adequate inprod
head(drop(Adot_vec))
head(fda.usc::Adot(X_fdata))

# Obtention of the entire Adot matrix
Ad <- diag(rep(Adot_vec[1], n))
Ad[upper.tri(Ad, diag = FALSE)] <- Adot_vec[-1]
head(Ad <- t(Ad) + Ad - diag(diag(Ad)))

# Positive definite
eigen(Ad)$values

# Warning if X contains repeated observations
# Adot(X = rbind(1:3, 1:3, 3:5))

# Comparison with .Fortran("adot", PACKAGE = "fda.usc")
```
```r
n <- as.integer(n)
a <- as.double(rep(0, (n * (n - 1) / 2 + 1)))
inprod <- X_fpc["scores"] %*% t(X_fpc["scores"])
inprod <- inprod[upper.tri(inprod, diag = TRUE)]
X <- X_fpc["scores"]
microbenchmark::microbenchmark(
  .Fortran("adot", n = n, inprod = inprod, Adot_vec = a,
           PACKAGE = "fda.usc"),
  Adot(X = X),
  times = 50, control = list(warmup = 10))
```

---

**flm_term**

*Functional linear model term with bivariate kernel*

**Description**

Computation of the functional linear term

\[
\int_a^b \beta(s,t)X(s) \, ds,
\]

of a Functional Linear Model with Functional Response (FLMFR), where \(X\) is a random variable in the Hilbert space of square-integrable functions in \([a,b]\), \(L^2([a,b])\), \(\beta\) is the bivariate kernel of the FLMFR, and \(\varepsilon\) is a random variable in \(L^2([c,d])\).

**Usage**

```r
flm_term(X_fdata, beta, t, int_rule = "trapezoid", equispaced = NULL, concurrent = FALSE)
```

**Arguments**

- **X_fdata**: sample of functional data as an `fdata` object of length \(n\).
- **beta**: matrix containing the values \(\beta(s,t)\), for each grid point \(s\) in \([a,b]\) and \(t\) in \([c,d]\). If `concurrent = TRUE`, a row/column vector must be introduced, valued in the same grid as `error_fdata`, with the same length as `length(X_fdata$argvals)`.
- **t**: vectors with grid points where functions are valued.
- **int_rule**: quadrature rule for approximating the definite unidimensional integral: trapezoidal rule (int_rule = "trapezoid") and extended Simpson rule (int_rule = "Simpson") are available. Defaults to "trapezoid".
- **equispaced**: flag to indicate if `X_fdata$data` is valued in an equispaced grid or not. Defaults to FALSE.
- **concurrent**: flag to consider a concurrent FLMFR (degenerate case). Defaults to FALSE.

**Value**

Functional linear model term as the integral (in \(s\)) between \(X\) and `beta`, as an `fdata` object of length \(n\).
Author(s)
Javier Álvarez-Liébana.

References

Examples
```r
## Generate a sample of functional responses via FLMFR

# Bivariate kernel beta(s,t) as an egg carton shape
s <- seq(0, 1, l = 101)
t <- seq(0, 1, l = 201)
beta <- outer(s, t, FUN = function(s, t) sin(6 * pi * s) + cos(6 * pi * t))

# Functional data as zero-mean Gaussian process with exponential variogram
X_fdata <- fda.usc::rproc2fdata(n = 50, t = s, sigma = "vexponential",
                               list = list(scale = 2.5))

# Functional error as an OU process with variance 0.075
sigma <- sqrt(0.075) * 2
error_fdata <- r_ou(n = 50, t = t, sigma = sigma)
Y_fdata <- flm_term(X_fdata = X_fdata, beta = beta, t = t) + error_fdata
plot(Y_fdata)

## Generate a sample of functional responses via concurrent model

# Function beta(t)
s <- seq(1, 3, l = 201)
t <- seq(2, 5, l = 201)
beta <- sin(pi * t) + cos(pi * t)

# Functional data as zero-mean Gaussian process with exponential variogram
X_fdata <- fda.usc::rproc2fdata(n = 50, t = s, sigma = "vexponential",
                               list = list(scale = 2.5))

# Functional error as an OU process with variance 0.075
sigma <- sqrt(0.075) * 2
error_fdata <- r_ou(n = 50, t = t, sigma = sigma)
Y_fdata <- flm_term(X_fdata = X_fdata, beta = beta, t = t, concurrent = TRUE) + error_fdata
plot(Y_fdata)
```

flm_test

Goodness-of-fit test for functional linear models
Description

Goodness-of-fit test of a functional linear model with functional response $Y \in L^2([c,d])$ and functional predictor $X \in L^2([a,b])$, where $L^2([a,b])$ is the Hilbert space of square-integrable functions in $[a,b]$.

The goodness-of-fit test checks the linearity of the regression model $m : L^2([a,b]) \to L^2([c,d])$ that relates $Y$ and $X$ by

$$Y(t) = m(X) + \varepsilon(t),$$

where $\varepsilon$ is a random variable in $L^2([c,d])$ and $t \in [c,d]$. The check is formalized as the test of the composite hypothesis

$$H_0 : m \in \{m_\beta : \beta \in L^2([a,b]) \otimes L^2([c,d])\},$$

where

$$m_\beta(X(s))(t) = \int_a^b \beta(s,t)X(s) \, ds$$

is the linear, Hilbert–Schmidt, integral operator parametrized by the bivariate kernel $\beta$. Its estimation is done by the truncated expansion of $\beta$ in the tensor product of the data-driven bases of Functional Principal Components (FPC) of $X$ and $Y$. The FPC basis for $X$ is truncated in $p$ components, while the FPC basis for $Y$ is truncated in $q$ components.

The particular cases in which either $X$ or $Y$ are constant functions give either a scalar predictor or response. The simple linear model arises if both $X$ and $Y$ are scalar, for which $\beta$ is a constant.

Usage

```r
flm_test(X, Y, beta0 = NULL, B = 500, est_method = "fpcr", p = NULL,
          q = NULL, thre_p = 0.99, thre_q = 0.99, lambda = NULL,
          boot_scores = TRUE, verbose = TRUE, plot_dens = TRUE,
          plot_proc = TRUE, plot_max_procs = 100, plot_max_p = 2,
          plot_max_q = 2, save_fit_flm = TRUE, save_boot_stats = TRUE,
          int_rule = "trapezoid", refit_lambda = FALSE, ...)
```

Arguments

- **X, Y**
  - samples of functional/scalar predictors and functional/scalar response. Either `fdata` objects (for functional variables) or vectors of length $n$ (for scalar variables).

- **beta0**
  - if provided (defaults to `NULL`), the simple null hypothesis $H_0 : m = m_{\beta_0}$ is tested. `beta0` must be a matrix of size $c(\text{length}(X$argvals$), \text{length}(Y$argvals$))$. If $X$ or $Y$ are scalar, `beta0` can be also an `fdata` object, with the same argvals as $X$ or $Y$. Can also be a constant (understood as a shorthand for a matrix with all its entries equal to the constant).

- **B**
  - number of bootstrap replicates. Defaults to 500.

- **est_method**
  - either "fpcr" (Functional Principal Components Regression; FPCR), "fpcr_l2" (FPCR with ridge penalty), "fpcr_l1" (FPCR with lasso penalty) or "fpcr_l1s" (FPCR with lasso-selected FPC). If $X$ is scalar, `flm_est` only considers "fpcr" as estimation method. See details below. Defaults to "fpcr_l1s".

p, q either index vectors indicating the specific FPC to be considered for the truncated bases expansions of X and Y, respectively. If a single number for p is provided, then \( p \leq 1: \max(p) \) internally (analogously for q) and the first \( \max(p) \) FPC are considered. If NULL (default), then a data-driven selection of p and q is done. See details below.

thre_p, thre_q thresholds for the proportion of variance that is explained, at least, by the first \( p \) and \( q \) FPC of X and Y, respectively. These thresholds are employed for an (initial) automatic selection of \( p \) and \( q \). Default to 0.99. \( \text{thre}_p \) (\( \text{thre}_q \)) is ignored if \( p \) (\( q \)) is provided.

lambda regularization parameter \( \lambda \) for the estimation methods “fpcr_l2”, "fpcr_l1", and "fpcr_l1s". If NULL (default), it is chosen with \texttt{cv.glmnet}.

boot_scores flag to indicate if the bootstrap shall be applied to the scores of the residuals, rather than to the functional residuals. This improves the computational expediency notably. Defaults to TRUE.

verbose flag to show information about the testing progress. Defaults to TRUE.

plot_dens flag to indicate if a kernel density estimation of the bootstrap statistics shall be plotted. Defaults to TRUE.

plot_proc whether to display a graphical tool to identify the degree of departure from the null hypothesis. If TRUE (default), the residual marked empirical process, projected in several FPC directions of X and Y, is shown, together with bootstrap analogues. The FPC directions are ones selected at the estimation stage.

plot_max_procs maximum number of bootstrapped processes to plot in the graphical tool. Set as the minimum of \( \text{plot\_max\_procs} \) and \( B \). Defaults to 100.

plot_max_p, plot_max_q maximum number of FPC directions to be considered in the graphical tool. They limit the resulting plot to be at most of size \( c(\text{plot\_max\_p},\text{plot\_max\_q}) \). Default to 2.

save_fit_flm, save_boot_stats flag to return \texttt{fit\_flm} and \texttt{boot\_*}. If FALSE, these memory-expensive objects are set to NA. Default to TRUE.

int_rule quadrature rule for approximating the definite unidimensional integral: trapezoidal rule (\texttt{int\_rule = "trapezoid"}) and extended Simpson rule (\texttt{int\_rule = "Simpson"}) are available. Defaults to "trapezoid".

refit_lambda flag to reselct \texttt{lambda} in each bootstrap replicate, incorporating its variability in the bootstrap calibration. Much more time consumig. Defaults to FALSE.

... further parameters to be passed to \texttt{cv.glmnet} (and then to \texttt{cv.glmnet}) such as \texttt{cv_1se}, \texttt{cv_nlambda} or \texttt{cv_parallel}, among others.

Details

The function implements the bootstrap-based goodness-of-fit test for the functional linear model with functional/scalar response and functional/scalar predictor, as described in Algorithm 1 in García-Portugués et al. (2021). The specifics are detailed there.

By default \texttt{cv_1se = TRUE} for \texttt{cv.glmnet} is considered, unless it is changed via ... This is the recommended choice for conducting the goodness-of-fit test based on regularized estimators, as
the oversmoothed estimate of the regression model under the null hypothesis notably facilitates the calibration of the test (see García-Portugués et al., 2021).

The graphical tool obtained with plot_proc = TRUE is based on an extension of the tool described in García-Portugués et al. (2014).

Repeated observations on \( X \) are internally removed, as otherwise they would cause NaNs in \( \text{Adot} \). Missing values on \( X \) and \( Y \) are also automatically removed.

Value

An object of the htest class with the following elements:

- **statistic**: test statistic.
- **p.value**: \( p \)-value of the test.
- **boot_statistics**: the bootstrapped test statistics, a vector of length \( B \).
- **method**: information on the type of test performed.
- **parameter**: a vector with the dimensions \( p \) and \( q \) considered in the test statistic. These are the lengths of the outputs \( p \) and \( q \).
- **p**: the index of the FPC considered for \( X \).
- **q**: the index of the FPC considered for \( Y \).
- **fit_flm**: the output resulted from calling \texttt{flm_est}.
- **boot_lambda**: bootstrapped \( \lambda \).
- **boot_p**: a list with the bootstrapped indexes of the FPC considered for \( X \).
- **data.name**: name of the value of data.

Author(s)

Eduardo García-Portugués.

References


Examples

```r
## Quick example for functional response and predictor

# Generate data under H0
n <- 100
set.seed(987654321)
X_fdata <- r_ou(n = n, t = seq(0, 1, l = 101), sigma = 2)
```
epsilon <- r_ou(n = n, t = seq(0, 1, l = 101), sigma = 0.5)
Y_fdata <- epsilon

# Test the FLMFR
flm_test(X = X_fdata, Y = Y_fdata)

# Simple hypothesis
flm_test(X = X_fdata, Y = Y_fdata, beta0 = 0)

# Generate data under H1
n <- 100
set.seed(987654321)
sample_frm_fr <- r_frm_fr(n = n, scenario = 3, s = seq(0, 1, l = 101),
t = seq(0, 1, l = 101), nonlinear = "quadratic")
X_fdata <- sample_frm_fr["X_fdata"]
Y_fdata <- sample_frm_fr["Y_fdata"]

# Test the FLMFR
flm_test(X = X_fdata, Y = Y_fdata)

## Functional response and predictor

# Generate data under H0
n <- 50
B <- 100
set.seed(987654321)
t <- seq(0, 1, l = 201)
X_fdata <- r_ou(n = n, t = t, sigma = 2)
epsilon <- r_ou(n = n, t = t, sigma = 0.5)
Y_fdata <- epsilon

# With boot_scores = TRUE
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr", B = B)
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l2", B = B)
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l1s", B = B)

# With boot_scores = FALSE
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr",
        boot_scores = FALSE, B = B)
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l2",
        boot_scores = FALSE, B = B)
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l1",
        boot_scores = FALSE, B = B)
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l1s",
        boot_scores = FALSE, B = B)

# Simple hypothesis
flm_test(X = X_fdata, Y = Y_fdata, beta0 = 2, est_method = "fpcr", B = B)
flm_test(X = X_fdata, Y = Y_fdata, beta0 = 0, est_method = "fpcr", B = B)
flm_test(X = X_fdata, Y = Y_fdata, beta0 = 0, est_method = "fpcr_l1s", B = B)

# Generate data under H1
n <- 50
```r
B <- 100
set.seed(987654321)
sample_frm_fr <- r_frm_fr(n = n, scenario = 3, s = t, t = t,
  nonlinear = "quadratic")
X_fdata <- sample_frm_fr$X_fdata
Y_fdata <- sample_frm_fr$Y_fdata

# With boot_scores = TRUE
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr", B = B)
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l2", B = B)
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l1s", B = B)

# With boot_scores = FALSE
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr",
  boot_scores = FALSE, B = B)
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l2",
  boot_scores = FALSE, B = B)
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l1",
  boot_scores = FALSE, B = B)
flm_test(X = X_fdata, Y = Y_fdata, est_method = "fpcr_l1s",
  boot_scores = FALSE, B = B)

## Scalar response and functional predictor

# Generate data under H0
n <- 50
B <- 100
set.seed(987654321)
t <- seq(0, 1, l = 201)
X_fdata <- r_ou(n = n, t = t, sigma = 2)
beta <- r_ou(n = 1, t = t, sigma = 0.5, x0 = 2)
epsilon <- rnorm(n = n)
Y <- drop(inprod_fdata(X_fdata1 = X_fdata, X_fdata2 = beta) + epsilon)

# With boot_scores = TRUE
flm_test(X = X_fdata, Y = Y, est_method = "fpcr", B = B)
flm_test(X = X_fdata, Y = Y, est_method = "fpcr_l2", B = B)
flm_test(X = X_fdata, Y = Y, est_method = "fpcr_l1s", B = B)

# With boot_scores = FALSE
flm_test(X = X_fdata, Y = Y, est_method = "fpcr",
  boot_scores = FALSE, B = B)
flm_test(X = X_fdata, Y = Y, est_method = "fpcr_l2",
  boot_scores = FALSE, B = B)
flm_test(X = X_fdata, Y = Y, est_method = "fpcr_l1",
  boot_scores = FALSE, B = B)
flm_test(X = X_fdata, Y = Y, est_method = "fpcr_l1s",
  boot_scores = FALSE, B = B)

# Simple hypothesis
flm_test(X = X_fdata, Y = Y, beta0 = beta, est_method = "fpcr", B = B)
flm_test(X = X_fdata, Y = Y, beta0 = 0, est_method = "fpcr", B = B)
flm_test(X = X_fdata, Y = Y, beta0 = 0, est_method = "fpcr_l1s", B = B)
```
# Generate data under H1

n <- 50
B <- 100
set.seed(987654321)
X_fdata <- r_ou(n = n, t = t, sigma = 2)
beta <- r_ou(n = 1, t = t, sigma = 0.5)
epsilon <- rnorm(n = n)
Y <- drop(exp(inprod_fdata(X_fdata1 = X_fdata^2, X_fdata2 = beta)) + epsilon)

# With boot_scores = TRUE
flm_test(X = X_fdata, Y = Y, est_method = "fpcr", B = B)
flm_test(X = X_fdata, Y = Y, est_method = "fpcr_l2", B = B)
flm_test(X = X_fdata, Y = Y, est_method = "fpcr_l1s", B = B)

# With boot_scores = FALSE
flm_test(X = X_fdata, Y = Y, est_method = "fpcr", boot_scores = FALSE, B = B)
flm_test(X = X_fdata, Y = Y, est_method = "fpcr_l2", boot_scores = FALSE, B = B)
flm_test(X = X_fdata, Y = Y, est_method = "fpcr_l1", boot_scores = FALSE, B = B)
flm_test(X = X_fdata, Y = Y, est_method = "fpcr_l1s", boot_scores = FALSE, B = B)

## Functional response and scalar predictor

# Generate data under H0

n <- 50
B <- 100
set.seed(987654321)
X <- rnorm(n)
X <- rnorm(n)
t <- seq(0, 1, l = 201)
X <- X
beta <- r_ou(n = 1, t = t, sigma = 0.5, x0 = 3)
beta$data <- matrix(beta$data, nrow = n, ncol = ncol(beta$data),
                        byrow = TRUE)
epsilon <- r_ou(n = n, t = t, sigma = 0.5)
Y_fdata <- X * beta + epsilon

# With boot_scores = TRUE
flm_test(X = X, Y = Y_fdata, est_method = "fpcr", B = B)

# With boot_scores = FALSE
flm_test(X = X, Y = Y_fdata, est_method = "fpcr", boot_scores = FALSE, B = B)

# Simple hypothesis
flm_test(X = X, Y = Y_fdata, beta0 = beta[1], est_method = "fpcr", B = B)
flm_test(X = X, Y = Y_fdata, beta0 = 0, est_method = "fpcr", B = B)

# Generate data under H1

n <- 50
B <- 100
set.seed(987654321)
```r
X <- rexp(n)
beta <- r_ou(n = 1, t = t, sigma = 0.5, x0 = 3)
beta$data <- matrix(beta$data, nrow = n, ncol = ncol(beta$data), byrow = TRUE)
epsilon <- r_ou(n = n, t = t, sigma = 0.5)
Y_fdata <- log(X * beta) + epsilon

# With boot_scores = TRUE
flm_test(X = X, Y = Y_fdata, est_method = "fpcr", B = B)

# With boot_scores = FALSE
flm_test(X = X, Y = Y_fdata, est_method = "fpcr", boot_scores = FALSE, B = B)
```

---

### fpc

**Computation of functional principal components**

#### Description

Computation of Functional Principal Components (FPC) for equispaced and non equispaced functional data.

#### Usage

```r
fpc(X_fdata, n_fpc = 3, centered = FALSE, int_rule = "trapezoid",
equispaced = FALSE, verbose = FALSE)
```

#### Arguments

- **X_fdata**: sample of functional data as an `fdata` object of length `n`.
- **n_fpc**: number of FPC to be computed. If `n_fpc > n`, `n_fpc` is set to `n`. Defaults to 3.
- **centered**: flag to indicate if `X_fdata` is centered or not. Defaults to `FALSE`.
- **int_rule**: quadrature rule for approximating the definite unidimensional integral: trapezoidal rule (`int_rule = "trapezoid"`) and extended Simpson rule (`int_rule = "Simpson"`) are available. Defaults to "trapezoid".
- **equispaced**: flag to indicate if `X_fdata$data` is valued in an equispaced grid or not. Defaults to `FALSE`.
- **verbose**: whether to show or not information about the fpc procedure. Defaults to `FALSE`.

#### Details

The FPC are obtained by performing the single value decomposition

\[
X W^{1/2} = UD(V'W^{1/2})
\]

where `X` is the matrix of discretized functional data, `W` is a diagonal matrix of weights (computed by `w_integral1D` according to `int_rule`), `D` is the diagonal matrix with singular values (standard deviations of FPC), `U` is a matrix whose columns contain the left singular vectors, and `V` is a matrix whose columns contain the right singular vectors (FPC).
Value

An "fpc" object containing the following elements:

- **d**: standard deviations of the FPC (i.e., square roots of eigenvalues of the empirical autocovariance estimator).
- **rotation**: orthonormal eigenfunctions (loadings or functional principal components), as an fdata class object.
- **scores**: rotated samples: inner products between X_fdata and eigenfunctions in rotation.
- **l**: vector of index of FPC.
- **equispaced**: equispaced flag.

Author(s)

Javier Álvarez-Liébana and Gonzalo Álvarez-Pérez.

References


Examples

```r
## Computing FPC for equispaced data

# Sample data
X_fdata1 <- r_ou(n = 200, t = seq(2, 4, l = 201))

# FPC with trapezoid rule
X_fpc1 <- fpc(X_fdata = X_fdata1, n_fpc = 50, equispaced = TRUE, int_rule = "trapezoid")

# FPC with Simpson's rule
X_fpc2 <- fpc(X_fdata = X_fdata1, n_fpc = 50, equispaced = TRUE, int_rule = "Simpson")

# Check if FPC are orthonormal
norms1 <- rep(0, length(X_fpc1$l))
for (i in X_fpc1$l) {
  norms1[i] <- integral1D(fx = X_fpc1$rotation$data[i, ]^2, t = X_fdata1$argvals)
}

norms2 <- rep(0, length(X_fpc2$l))
for (i in X_fpc2$l) {
  norms2[i] <- integral1D(fx = X_fpc2$rotation$data[i, ]^2, t = X_fdata1$argvals)
}
```
## Computing FPC for non equispaced data

# Sample data
X_fdata2 <- r_ou(n = 200, t = c(seq(0, 0.5, l = 201), seq(0.51, 1, l = 301)))

# FPC with trapezoid rule
X_fpc3 <- fpc(X_fdata = X_fdata2, n_fpc = 5, int_rule = "trapezoid",
equispaced = FALSE)

# Check if FPC are orthonormal
norms3 <- rep(0, length(X_fpc3$l))
for (i in X_fpc3$l) {
  norms3[i] <- integral1D(fx = X_fpc3$rotation$data[i, ]^2,
t = X_fdata2$argvals)
}

## Efficiency comparisons

# fpc() vs. fda.usc::fdata2pc()
data(phoneme, package = "fda.usc")
mlearn <- phoneme$learn[1:10, ]
res1 <- fda.usc::fdata2pc(mlearn, ncomp = 3)
res2 <- fpc(X_fdata = mlearn, n_fpc = 3)
plot(res1$x[, 1:3], col = 1)
points(res2$scores, col = 2)
microbenchmark::microbenchmark(fda.usc::fdata2pc(mlearn, ncomp = 3),
fpc(X_fdata = mlearn, n_fpc = 3), times = 1e3,
control = list(warmup = 20))

---

fpc_utils

Utilities for functional principal components

### Description

Computation of coefficients and reconstructions based on Functional Principal Components (FPC). The function `fpc_coefs` allows to project a functional data sample into a basis of FPC; the reconstruction of the sample from its projections and the FPC is done with `fpc_to_fdata`. The functions `beta_fpc_coefs` and `fpc_to_beta` do analogous operations but for the bivariate kernel $\beta$ and the tensor product of two FPC bases.

### Usage

```r
fpc_coefs(X_fdata, X_fpc, ind_X_fpc = 1:3, int_rule = "trapezoid")
```

```r
beta_fpc_coefs(beta, X_fpc, Y_fpc, ind_X_fpc = 1:3, ind_Y_fpc = 1:3,
```
int_rule = "trapezoid"

fpc_to_fdata(coefs, X_fpc, ind_coefs = 1:ncol(coefs))

fpc_to_beta(beta_coefs, X_fpc, Y_fpc, ind_coefs_X = 1:nrow(beta_coefs),
ind_coefs_Y = 1:ncol(beta_coefs))

Arguments

X_fdata sample of functional data as an \texttt{fdata} object of length \(n\).

X_fpc, Y_fpc "fpc" objects as resulted from calling \texttt{fpc}.

ind_X_fpc, ind_Y_fpc vectors giving the FPC indexes for whom the coefficients are computed. Their lengths must be smaller than the number of FPC in \(X_fpc\) and \(Y_fpc\), respectively. Default to 1:3.

int_rule quadrature rule for approximating the definite unidimensional integral: trapezoidal rule (\(\text{int_rule} = \text{"trapezoid"}\)) and extended Simpson rule (\(\text{int_rule} = \text{"Simpson"}\)) are available. Defaults to "trapezoid".

beta a matrix containing the bivariate kernel \(\beta\) evaluated on a grid. Must be of size \(c(\text{length}(X_fpc$rotation$argvals), \text{length}(Y_fpc$rotation$argvals))\).

coefs a vector of coefficients to combine linearly the FPC. Its length must be smaller than the number of FPC in \(X_fpc\).

ind_coefs, ind_coefs_X, ind_coefs_Y indexes of FPC to associate to the provided coefficients. By default, from the first FPC to the sizes of \(\text{coefs}\) or \(\beta\) \text{coefs}.

beta_coefs a matrix of coefficients to combine linearly the tensor products of FPC. Its size must be smaller than the number of FPC in \(X_fpc\) and \(Y_fpc\).

Value

fpc_coefs a vector of the same length as \text{coefs} containing the coefficients of \(X_fdata\) in the FPC of \(X_fpc\).

beta_fpc_coefs a matrix of the same size as \text{beta_coefs} containing the coefficients of \(\beta\) in the tensor product of the FPC in \(X_fpc\) and \(Y_fpc\).

fpc_to_fdata an \texttt{fdata} object of the same type as \(X_fpc$rotation\).

fpc_to_beta a matrix with the reconstructed kernel and size \(c(\text{length}(X_fpc$rotation$argvals), \text{length}(Y_fpc$rotation$argvals))\).

Author(s)

Eduardo García-Portugués.

References

Examples

```r
## Compute FPC coefficients and reconstruct data

# Sample data
X_fdata <- r_ou(n = 200, t = seq(2, 4, l = 201))

# Compute FPC
X_fpc <- fpc(X_fdata = X_fdata, n_fpc = 50)

# FPC coefficients are the same if the data is centered
fpc_coefs(X_fdata = fdata_cen(X_fdata), X_fpc = X_fpc)[1:4, ]
X_fpc$scores[1:4, 1:3]

# Reconstruct the first two curves for an increasing number of FPC
plot(X_fdata[1:2, ], col = 1)
n_fpc <- c(2, 5, 10, 25, 50)
for (j in 1:5) {
  lines(fpc_to_fdata(X_fpc = X_fpc, coefs = X_fpc$scores[, 1:n_fpc[j]])[1:2, ], col = j + 1)
}

## Project and reconstruct beta

# Surface
beta_fun <- function(s, t) sin(6 * pi * s) + cos(6 * pi * t)
s <- seq(0, 1, l = 101)
t <- seq(0, 1, l = 201)
beta_surf <- outer(s, t, FUN = beta_fun)

# Functional data as zero-mean Gaussian process with exponential variogram
X_fdata <- fda.usc::rproc2fdata(n = 100, t = s, sigma = "vexponential",
list = list(scale = 2.5))
Y_fdata <- flm_term(X_fdata = X_fdata, beta = beta_surf, t = t) +
r_ou(n = 100, t = t, sigma = sqrt(0.075) * 2)

# FPC
X_fpc <- fpc(X_fdata = X_fdata, n_fpc = 50)
Y_fpc <- fpc(X_fdata = Y_fdata, n_fpc = 50)

# Coefficients
beta_coefs <- beta_fpc_coefs(beta = beta_surf, X_fpc = X_fpc, Y_fpc = Y_fpc,
ind_X_fpc = 1:50, ind_Y_fpc = 1:50)

# Reconstruction
beta_surf1 <- fpc_to_beta(beta_coefs = beta_coefs[1:2, 1:5],
X_fpc = X_fpc, Y_fpc = Y_fpc)
beta_surf2 <- fpc_to_beta(beta_coefs = beta_coefs[1:15, 1:10],
X_fpc = X_fpc, Y_fpc = Y_fpc)
beta_surf3 <- fpc_to_beta(beta_coefs = beta_coefs[1:50, 1:50],
X_fpc = X_fpc, Y_fpc = Y_fpc)

# Show reconstructions
```
old_par <- par(mfrow = c(2, 2))
col <- viridisLite::viridis(20)
image(s, t, beta_surf, col = col, zlim = c(-2.5, 2.5), main = "Original")
image(s, t, beta_surf1, col = col, zlim = c(-2.5, 2.5), main = "2 x 5")
image(s, t, beta_surf2, col = col, zlim = c(-2.5, 2.5), main = "15 x 10")
image(s, t, beta_surf3, col = col, zlim = c(-2.5, 2.5), main = "50 x 50")
par(old_par)

ontario

Ontario temperature and electricity consumption during 2010–2014

Description

Real dataset employed Benatia et al. (2017). Contains the hourly electricity consumption and air temperature curves in the province of Ontario (Canada). It features a set of daily curves during the summer months of 2010–2014.

Usage

ontario

Format

A list with the following entries:

**temp** an fdata with 368 smoothed daily temperature (in Celsius degrees) curves of the Ontario province, discretized on 73 equispaced grid points on $[-24, 48]$ (see examples).

**elec** an fdata with the daily electricity consumption (in gigawatts) curves of the Ontario province. Discretized on 25 equispaced grid points on $[0, 24]$.

**df** a dataframe with time metadata for each curve:

- date: the date of the observation, a POSIXct object.
- weekday: the weekday of the observation.

Details

The summer months correspond to June 1st to September 15th. Weekend days and holidays are disregarded.

The smoothed temperature curves are constructed by a weighted average of the temperatures of 41 Ontarian cities that is afterwards smoothed with a local polynomial regression. The curves correspond to a 3-days window of the temperature (see examples). The temperature is standardized such that its original minimum, 6 °C, is subtracted.

The electricity consumption curves are discretized on the interval $[0, 24]$. That means that the last observation of the $i$-th curve is the same as the first observation of the $(i + 1)$-th curve if the curves correspond to consecutive days.

See more details about the construction of the dataset in Benatia et al. (2017).
Author(s)

Data gathered and processed by David Benatia, Marine Carrasco, and Jean-Pierre Florens. Javier Álvarez-Liébana and Eduardo García-Portugués imported the dataset and added temporal metadata.

Source

The dataset comes from the companion data to Benatia et al. (2017), which was retrieved from the first author’s website. The source of the electricity consumption data is the System operator’s website. The source of the preprocessed temperature values is the Environment Canada’s website.

References


Examples

```r
## Show data

# Load data
data("ontario")

# Plot
old_par <- par(mfrow = c(1, 2))
plot(ontario$temp)
plot(ontario$elec)
par(old_par)

# Observe the 3-day windows for each observation
plot(ontario$temp$argvals, ontario$temp$data[2, ], type = "o",
     xlim = c(-48, 72), ylim = c(7, 13), xlab = "Hours",
     ylab = "Electricity consumption", pch = 16)
points(ontario$temp$argvals - 24, ontario$temp$data[1, ], col = 3, pch = 2)
points(ontario$temp$argvals + 24, ontario$temp$data[3, ], col = 2, cex = 1.5)
abline(v = 24 * -2:3, lty = 2)
legend("top", legend = c("Curve 1", "Curve 2", "Curve 3"), col = c(3, 1, 2),
       pt.cex = c(1, 1, 1.5), pch = c(2, 16, 1))

# If the days are not consecutive, then the electricity consumptions at the
# end of one day and the beginning of the next do not match
head(abs(ontario$elec$data[-368, 25] - ontario$elec$data[-1, 1]))
head(diff(ontario$df$date))

## Test the linear model with functional response and predictor

(comp_flmfr <- flm_test(X = ontario$temp, Y = ontario$elec,
                         est_method = "fpcr_l1s"))
(simp_flmfr <- flm_test(X = ontario$temp, Y = ontario$elec,
                        beta0 = 0, est_method = "fpcr_l1s"))

# Visualize estimation
```
Simulation of an Ornstein–Uhlenbeck process

Description

Simulation of trajectories of the Ornstein–Uhlenbeck process \( \{ X_t \} \). The process is the solution to the stochastic differential equation

\[
dX_t = \alpha (X_t - \mu) dt + \sigma dW_t,
\]

whose stationary distribution is \( N(\mu, \sigma^2/(2\alpha)) \), for \( \alpha, \sigma > 0 \) and \( \mu \in \mathbb{R} \).

Given an initial point \( x_0 \) and the evaluation times \( t_1, \ldots, t_m \), a sample trajectory \( X_{t_1}, \ldots, X_{t_m} \) can be obtained by sampling the joint Gaussian distribution of \( (X_{t_1}, \ldots, X_{t_m}) \).

Usage

```r
r_ou(n, t = seq(0, 1, len = 201), mu = 0, alpha = 1, sigma = 1,
     x0 = rnorm(n, mean = mu, sd = sigma/sqrt(2 * alpha)))
```

Arguments

- \( n \) number of trajectories to sample.
- \( t \) evaluation times for the trajectories, a vector.
- \( \mu \) mean of the process, a scalar.
- \( \alpha \) strength of the drift, a positive scalar.
- \( \sigma \) diffusion coefficient, a positive scalar.
- \( x_0 \) a vector of length \( n \) giving the initial values of the Ornstein–Uhlenbeck trajectories. By default, \( n \) points are sampled from the stationary distribution. If a single scalar is passed, then the same \( x_0 \) is employed for all the trajectories.

Value

Random trajectories, an \texttt{fdata} object of length \( n \) and \( t \) as \texttt{argvals}.

Author(s)

Eduardo García-Portugués.
Sampling functional regression models with functional responses

Description

Simulation of a Functional Regression Model with Functional Response (FRMFR) comprised of an additive mix of a linear and nonlinear terms:

\[ Y(t) = \int_a^b X(s)\beta(s,t)ds + \Delta(X)(t) + \varepsilon(t), \]

where \( X \) is a random variable in the Hilbert space of square-integrable functions in \([a,b], L^2([a,b])\), \( \beta \) is the bivariate kernel of the FRMFR, \( \varepsilon \) is a random variable in \( L^2([c,d]) \), and \( \Delta(X) \) is a nonlinear term.

In particular, the scenarios considered in García-Portugués et al. (2021) can be easily simulated.

Usage

```
r_frm_fr(n, scenario = 3, X_fdata = NULL, error_fdata = NULL, beta = NULL, s = seq(0, 1, l = 101), t = seq(0, 1, l = 101), std_error = 0.15, nonlinear = NULL, concurrent = FALSE, int_rule = "trapezoid", n_fpc = 50, verbose = FALSE, ...)
nl_dev(X_fdata, t = seq(0, 1, l = 101), nonlinear = NULL, int_rule = "trapezoid", equispaced = equispaced, verbose = FALSE)
```

Arguments

- **n**: sample size, only required when scenario is given.
- **scenario**: an index from 1 to 3 (default) denoting one of the scenarios (S1, S2 or S3) simulated in García-Portugués et al. (2021) (see details below). If scenario = NULL, X_fdata, error_fdata, and beta have to be provided. Otherwise, X_fdata, error_fdata, and beta will be ignored.
- **X_fdata**: sample of functional covariates \( X(s) \) as `fdata` objects of length \( n \), with \( s \) in \([a,b]\). Defaults to NULL.
- **error_fdata**: sample of functional errors \( \varepsilon(t) \) as `fdata` objects of length \( n \), with \( t \) in \([c,d]\). If concurrent = TRUE, X_fdata and error_fdata must be valued in the same grid. Defaults to NULL.
beta matrix containing the values $\beta(s, t)$, for each grid point $s$ in $[a, b]$ and $t$ in $[c, d]$. If `concurrent = TRUE` (see details below), a row/column vector must be introduced, valued in the same grid as `error_fdata`. If `beta = NULL` (default), `scenario != NULL` is required.

$s, t$ grid points. If $X_fdata$, `error_fdata` and `beta` are provided, $s$ and $t$ are ignored. Default to $s = \text{seq}(0, 1, l = 101)$ and $t = \text{seq}(0, 1, l = 101)$, respectively.

`std_error` standard deviation of the random variables involved in the generation of the functional error `error_fdata`. Defaults to 0.15.

`nonlinear` nonlinear term. Either a character string ("exp", "quadratic" or "sin") or an `fdata` object of length n, valued in the same grid as `error_fdata`. If `nonlinear = NULL` (default), the nonlinear part is set to zero.

`concurrent` flag to consider a concurrent FLRFR (degenerate case). Defaults to FALSE.

`int_rule` quadrature rule for approximating the definite unidimensional integral: trapezoidal rule (`int_rule = "trapezoid"`) and extended Simpson rule (`int_rule = "Simpson"`) are available. Defaults to "trapezoid".

`n_fpc` number of components to be considered for the generation of functional variables. Defaults to 50.

`verbose` flag to display information about the sampling procedure. Defaults to FALSE.

... further parameters passed to `r_cm2013_flmfr`, `r_gof2021_flmfr` and `r_ik2018_flmfr`, depending on the chosen scenario.

`equispaced` flag to indicate if $X_fdata$data is valued in an equispaced grid or not. Defaults to FALSE.

Details

- `r_frm_fr` samples the above regression model, where the nonlinear term $\Delta(X)$ is computed by `nl_dev`. Functional covariates, errors, and $\beta$ are generated automatically from the scenarios in García-Portugués et al. (2021) when `scenario != NULL` (see the documentation of `r_gof2021_flmfr`). If `scenario = NULL`, covariates, errors and $\beta$ must be provided.

When `concurrent = TRUE`, the concurrent FRMFR

$$Y(t) = X(t)\beta(t) + \Delta(X)(t) + \varepsilon(t)$$

is considered.

- `nl_dev` computes a nonlinear deviation $\Delta(X)$: $\exp(\sqrt{X(a + (t - c)((b-a)/(d-c)))})$ (for "exp"), $(X^2(a + (t - c)((b-a)/(d-c))) - 1)$ ("quadratic") or $(\sin(2\pi t) - \cos(2\pi t))||X||^2$ ("sin"). Also, $\Delta(X)$ can be manually set as an `fdata` object of length n and valued in the same grid as `error_fdata`.

Value

A list with the following elements:

- `X_fdata` functional covariates, an `fdata` object of length n.
- `Y_fdata` functional responses, an `fdata` object of length n.
error_fdata functional errors, an \texttt{fdata} object of length $n$.

beta either the matrix with $\beta(s,t)$ evaluated at the argvals of $X_{\text{fdata}}$ and $Y_{\text{fdata}}$ (if \texttt{concurrent} = FALSE) or a vector with $\beta(t)$ evaluated at the argvals of $X_{\text{fdata}}$ (if \texttt{concurrent} = TRUE).

nl_dev nonlinear term, an \texttt{fdata} object of length $n$.

\textbf{Author(s)}

Javier Álvarez-Liébana.

\textbf{References}


\textbf{Examples}

```r
## Generate samples for the three scenarios

# Equispaced grids and Simpson's rule
s <- seq(0, 1, l = 101)
samp <- list()
old_par <- par(mfrow = c(3, 5))
for (i in 1:3) {
samp[[i]] <- r_frm_fr(n = 100, scenario = i, s = s, t = s,
                  int_rule = "Simpson")
plot(samp[[i]]$X_fdata)
plot(samp[[i]]$error_fdata)
plot(samp[[i]]$Y_fdata)
plot(samp[[i]]$nl_dev)
image(x = s, y = s, z = samp[[i]]$beta, col = viridisLite::viridis(20))
}
par(old_par)

## Linear term as a concurrent model

# The grids must be have the same number of grid points for a given
# nonlinear term and a given beta function
s <- seq(1, 2, l = 101)
t <- seq(0, 1, l = 101)
samp_c_1 <- r_frm_fr(n = 100, scenario = 3, beta = sin(t) - exp(t),
                  s = s, t = t, nonlinear = fda.usc::fdata(mdata =
                  t(matrix(rep(sin(t), 100), nrow = length(t))),
                  argvals = t),
                  concurrent = TRUE)
old_par <- par(mfrow = c(3, 2))
plot(samp_c_1$X_fdata)
plot(samp_c_1$error_fdata)
```

## Sample for given $X_{fdata}$, $error_{fdata}$, and $beta$

# Non equispaced grids with sinusoidal nonlinear term and intensity 0.5

s <- c(seq(0, 0.5, l = 50), seq(0.51, 1, l = 101))
t <- seq(2, 4, len = 151)

$X_{fdata}$ <- r_ou(n = 100, t = s, alpha = 2, sigma = 4, x0 = 1:100)
$error_{fdata}$ <- r_ou(n = 100, t = t, alpha = 1, sigma = 1, x0 = 1:100)

beta <- r_gof2021_flmfr(n = 100, s = s, t = t)$beta

$samp_{Xeps}$ <- r_frm_fr(scenario = NULL, $X_{fdata}$ = $X_{fdata}$, 
                          $error_{fdata}$ = $error_{fdata}$, $beta$ = beta, 
                          nonlinear = "exp", int_rule = "trapezoid")

old_par <- par(mfrow = c(3, 2))

plot(samp_Xeps$X_fdata)
plot(samp_Xeps$error_fdata)
plot(samp_Xeps$Y_fdata)
plot(samp_Xeps$nl_dev)

image(x = s, y = t, z = beta, col = viridisLite::viridis(20))
par(old_par)
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