

# Package ‘GET’

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**Description** Implementation of global envelopes for a set of general  $d$ -dimensional vectors  $T$  in various applications. A  $100(1-\alpha)\%$  global envelope is a band bounded by two vectors such that the probability that  $T$  falls outside this envelope in any of the  $d$  points is equal to  $\alpha$ . Global means that the probability is controlled simultaneously for all the  $d$  elements of the vectors. The global envelopes can be used for graphical Monte Carlo and permutation tests where the test statistic is a multivariate vector or function (e.g. goodness-of-fit testing for point patterns and random sets, functional analysis of variance, functional general linear model,  $n$ -sample test of correspondence of distribution functions), for central regions of functional or multivariate data (e.g. outlier detection, functional boxplot) and for global confidence and prediction bands (e.g. confidence band in polynomial regression, Bayesian posterior prediction). See Myllymäki and Mrkvička (2024) <doi:10.18637/jss.v111.i03>, Myllymäki et al. (2017) <doi:10.1111/rssb.12172>, Mrkvička and Myllymäki (2023) <doi:10.1007/s11222-023-10275-7>, Mrkvička et al. (2016) <doi:10.1016/j.spasta.2016.04.005>, Mrkvička et al. (2017) <doi:10.1007/s11222-016-9683-9>, Mrkvička et al. (2020) <doi:10.14736/kyb-2020-3-0432>, Mrkvička et al. (2021) <doi:10.1007/s11009-019-09756-y>, Myllymäki et al. (2021) <doi:10.1016/j.spasta.2020.100436>, Mrkvička et al. (2022) <doi:10.1002/sim.9236>, Dai et al. (2022) <doi:10.5772/intechopen.100124>, Dvořák and Mrkvička (2022) <doi:10.1007/s00180-021-01134-y>, Mrkvička et al. (2023) <doi:10.48550/arXiv.2309.04746>, and Konstantinou et al. (2024) <doi:10.1007/s00180-024-01569-z>.

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## Description

The **GET** package provides implementation of global envelopes for a set of general  $d$ -dimensional vectors  $T$  in various applications. A  $100(1-\alpha)\%$  the probability that  $T$  falls outside this envelope in any of the  $d$  points is equal to  $\alpha$ . Global means that the probability is controlled simultaneously for all the  $d$  elements of the vectors. The global envelopes can be used for central regions of functional or multivariate data (e.g. outlier detection, functional boxplot), for graphical Monte Carlo and permutation tests where the test statistic is a multivariate vector or function (e.g. goodness-of-fit testing for point patterns and random sets, functional ANOVA, functional GLM,  $n$ -sample test of correspondence of distribution functions), and for global confidence and prediction bands (e.g. confidence band in polynomial regression, Bayesian posterior prediction).

## Details

The **GET** package provides central regions (i.e. global envelopes) and global envelope tests with intrinsic graphical interpretation. The central regions can be constructed from (functional) data. The tests are Monte Carlo or permutation tests, which demand simulations from the tested null model. The methods are applicable for any multivariate vector data and functional data (after discretization).

To get an overview of the package, start R and type `library("GET")` and `vignette("GET")`.

To get examples of point pattern analysis, start R and type `library("GET")` and `vignette("pointpatterns")`.

To get examples of Mrkvička and Myllymäki (2022), start R and type `library("GET")` and `vignette("FDRenvelopes")`.

## Key functions in GET

- *Central regions or global envelopes or confidence bands:* `central_region`. E.g. 50% central region of growth curves of girls `growth`.
  - First create a `curve_set` of the growth curves, e.g.
 

```
cset <- curve_set(r = as.numeric(row.names(growth$hgtf)), obs = growth$hgtf)
```
  - Then calculate 50% central region (see `central_region` for further arguments)
 

```
cr <- central_region(cset, coverage = 0.5)
```
  - Plot the result (see `plot.global_envelope` for plotting options)
 

```
plot(cr)
```

It is also possible to do combined central regions for several sets of curves provided in a list for the function, see examples in `central_region`.

- *Global envelope tests:* `global_envelope_test` is the main function. E.g. A test of complete spatial randomness (CSR) for a point pattern  $X$ :
 

```
X <- spruces # an example pattern from spatstat
```

  - Use the function `envelope` of **spatstat** to create `nsim` simulations under CSR and to calculate the functions you want (below  $K$ -functions by Kest). Important: use the option `'savefuns=TRUE'` and specify the number of simulations `nsim`.

```
env <- envelope(X, nsim=999, savefuns = TRUE, fun = Kest, simulate = expression(runifpoint(ex
= X)))
```

- Perform the test (see [global\\_envelope\\_test](#) for further arguments)  
res <- global\_envelope\_test(env)
- Plot the result (see [plot.global\\_envelope](#) for plotting options)  
plot(res)

It is also possible to do combined global envelope tests for several sets of curves provided in a list for the function, see examples in [global\\_envelope\\_test](#). To obtain false discovery rate envelopes of Mrkvička and Myllymäki (2023) use the argument `typeone = "fdr"`.

- *Functional ordering*: [central\\_region](#) and [global\\_envelope\\_test](#) are based on different measures for ordering the functions (or vectors) from the most extreme to the least extreme ones. The core functionality of calculating the measures is in the function [forder](#), which can be used to obtain different measures for sets of curves. Usually there is no need to call [forder](#) directly.
- *Functional boxplots*: [fBoxplot](#)
- *Adjusted global envelope tests for composite null hypotheses*
  - [GET.composite](#), see a detailed example in [saplings](#)
- *One-way functional ANOVA*:
  - Graphical functional ANOVA tests: [graph.fanova](#)
  - Global rank envelope based on F-values: [frank.fanova](#)
- *Functional general linear model (GLM)*:
  - Graphical functional GLM: [graph.flm](#)
  - Global rank envelope based on F-values: [frank.flm](#)
  - For large data (not fitting comfortably in memory): [partial\\_forder](#)
- *Functional clustering*: [fclustering](#)
- *Global quantile regression*: [global\\_rq](#)
- Functions for performing global envelopes for other specific purposes:
  - Graphical n sample test of correspondence of distribution functions: [GET.distrequl](#)
  - Permutation-based tests of independence to samples from any bivariate distribution: [GET.distrindep](#)
  - Testing global and local dependence of point patterns on covariates: [GET.spatialF](#)
  - Testing local correlations: [GET.localcor](#)
  - Variogram and residual variogram with global envelopes: [GET.variogram](#)
- Deviation tests (for simple hypothesis): [deviation\\_test](#) (no graphical interpretation)
- Most functions accept the curves provided in a `curve_set` object. Use [curve\\_set](#) to create a `curve_set` object from the functions. Other formats to provide the curves to the above functions are also accepted, see the information on the help pages.

See the help files of the functions for examples.

### Workflow for (single hypothesis) tests based on single functions

To perform a test you always first need to obtain the test function  $T(r)$  for your data ( $T_1(r)$ ) and for each simulation ( $T_2(r), \dots, T_{s+1}(r)$ ) in one way or another. Given the set of the functions  $T_i(r), i = 1, \dots, s + 1$ , you can perform a test by [global\\_envelope\\_test](#).

1) The workflow when using your own programs for simulations:

- (Fit the model and) Create  $s$  simulations from the (fitted) null model.
- Calculate the functions  $T_1(r), T_2(r), \dots, T_{s+1}(r)$ .
- Use [curve\\_set](#) to create a `curve_set` object from the functions  $T_i(r), i = 1, \dots, s + 1$ .
- Perform the test  

```
res <- global_envelope_test(curve_set)
```

 where `curve_set` is the 'curve\_set'-object you created, and plot the result  

```
plot(res)
```

2) The workflow utilizing **spatstat**: start R, type `library("GET")` and `vignette("pointpatterns")`, which explains the workflow and gives many examples of point pattern analysis

### Functions for modifying sets of functions

It is possible to modify the curve set  $T_1(r), T_2(r), \dots, T_{s+1}(r)$  for the test.

- You can choose the interval of distances  $[r_{\min}, r_{\max}]$  by [crop\\_curves](#).
- For better visualisation, you can take  $T(r) - T_0(r)$  by [residual](#). Here  $T_0(r)$  is the expectation of  $T(r)$  under the null hypothesis.

### Example data (see references on the help pages of each data set)

- [abide\\_9002\\_23](#): see help page
- [adult\\_trees](#): a point pattern of adult rees
- [cgec](#): centred government expenditure centralization (GEC) ratios (see [graph.fanova](#))
- [fallen\\_trees](#): a point pattern of fallen trees
- [GDPTax](#): GDP per capita with country groups and other covariates
- [imageset3](#): a simulated set of images
- [rimov](#): water temperature curves in 365 days of the 36 years
- [saplings](#): a point pattern of saplings (see [GET.composite](#))

The data sets are used to show examples of the functions of the library.

### Number of functions

If the number of functions is low, the choice of the measure (or type or depth) plays a role, as explained in `vignette("GET")` (Section 2.4).

Note that the recommended minimum number of simulations for the rank envelope test (Myllymäki et al., 2017) based on a single function in spatial statistics is `nsim=2499`. When the number of argument values is large, also larger number simulations is needed in order to have a narrow p-interval. The "erl", "cont", "area", "qdir" and "st" global envelope tests and deviation tests can

be used with a lower number of simulations, although the Monte Carlo error is obviously larger with a lower number of simulations. For increasing the number of simulations, all the global rank envelopes approach the same curves.

Mrkvička et al. (2017) discussed the number of simulations for tests based on many functions.

## Documentation

Myllymäki and Mrkvička (2024) provides description of the package. The material can also be found in the corresponding vignette, which is available by starting R and typing `library("GET")` and `vignette("GET")`.

In the special case of spatial processes (spatial point processes, random sets), the functions are typically estimators of summary functions. The package supports the use of the R package **spatstat** for generating simulations and calculating estimators of the chosen summary function, but alternatively these can be done by any other way, thus allowing for any user-specified models/functions. To see examples of global envelopes for analysing point pattern data, start R, type `library("GET")` and `vignette("pointpatterns")`.

Mrkvička and Myllymäki (2023) developed false discovery rate (FDR) envelopes. Examples can be found by in associated vignette: start R, and type `library("GET")` and `vignette("pointpatterns")`.

Mrkvička et al. (2023a) proposed global quantile regression. An example of global quantile regression is given in the vignette `vignette("QuantileRegression")`.

The vignette `vignette("HotSpots")` illustrates the methodology proposed by Mrkvička et al. (2023b) for detecting hotspots on a linear network.

Type `citation("GET")` to get a full list of references.

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---

abide\_9002\_23

*Local brain activity at resting state*


---

## Description

Imaging measurements for local brain activity at resting state

## Usage

```
data("abide_9002_23")
```

## Format

A list of the `curve_set` containing the data, coordinates (x,y) where the data have been observed (third dimension is 23), the discrete factor `Group` (1=Autism; 2=Control), the discrete factor `Sex` (1=Male; 2=Female), and the continuous factor `Age`.



## Details

The data are a small part of ABIDE fALFF data available at ABIDE: [http://fcon\\_1000.projects.nitrc.org/indi/abide/fALFF](http://fcon_1000.projects.nitrc.org/indi/abide/fALFF): <http://fcp-indi.github.io/docs/user/alff.html> and distributed under the CC BY-NC-SA 3.0 license, <https://creativecommons.org/licenses/by-nc-sa/3.0/>.

The data are fractional Amplitude of Low Frequency Fluctuations (fALFF) (Zou et al. 2008) for Autism Brain Imaging Data Exchange collected resting state functional magnetic resonance imaging (R-fMRI) datasets (Di Martino et al. 2013). This data set in **GET** contains only a tiny part of the whole brain, namely the region 9002 (the right Cerebellum Crus 1) at slice 23 (see Figure 2 in Mrkvicka et al., 2019) for 514 individuals with the autism spectrum disorder (ASD) and 557 typical controls (TC) as specified in the given Group variable. Further the sex and age of each subject is given.

## References

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---

adult\_trees

*Adult trees data set*


---

## Description

Adult trees data set

## Usage

```
data("adult_trees")
```

## Format

A data.frame containing the locations (x- and y-coordinates) of 67 trees in an area of 75 m x 75 m.

## Details

A pattern of large trees (height > 25 m) originating from an uneven aged multi-species broadleaf nonmanaged forest in Kaluzhskie Zaseki, Russia.

The pattern is a sample part of data collected over 10 ha plot as a part of a research program headed by project leader Prof. O.V. Smirnova.

## References

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## See Also

[saplings](#)

## Examples

```
if(require("spatstat.geom", quietly=TRUE)) {
  data("adult_trees")
  adult_trees <- as.ppp(adult_trees, W = square(75))
  plot(adult_trees)
}
```

---

as.curve\_set

---

*Convert an envelope or fdata object to a curve\_set object*


---

## Description

If given an envelope object of **spatstat** or a fdata object of **fda.usc**, convert it into a curve\_set object. If given a curve\_set object, check its correctness and give it back.

## Usage

```
as.curve_set(curve_set, ...)
```

## Arguments

curve_set	An object to be converted to a <a href="#">curve_set</a> object. The envelope objects of <b>spatstat</b> and fdata objects of <b>fda.usc</b> are supported currently, besides curve_set objects.
...	Allows passing arguments to <a href="#">create_curve_set</a> .

**Value**

If an envelope object of **spatstat** or an **fdata** object is given, return a corresponding `curve_set` object. If a `curve_set` object was given, check it and return it unharmed.

---

central_region	<i>Central region / Global envelope</i>
----------------	---

---

**Description**

Provides central regions or global envelopes or confidence bands

**Usage**

```
central_region(
  curve_sets,
  type = "erl",
  coverage = 0.5,
  alternative = c("two.sided", "less", "greater"),
  probs = c(0.025, 0.975),
  quantile.type = 7,
  central = "median",
  nstep = 2,
  ...
)
```

**Arguments**

<code>curve_sets</code>	A <a href="#">curve_set</a> object or a list of <a href="#">curve_set</a> objects. Also envelope objects of <b>spatstat</b> and <b>fdata</b> of <b>fda.usc</b> are accepted instead of <code>curve_set</code> objects.
<code>type</code>	The type of the global envelope with current options for 'rank', 'erl', 'cont', 'area', 'qdir', 'st' and 'unscaled'. See details.
<code>coverage</code>	A number between 0 and 1. The 100*coverage% central region will be calculated. A vector of values can also be provided, leading to the corresponding number of central regions.
<code>alternative</code>	A character string specifying the alternative hypothesis. Must be one of the following: "two.sided" (default), "less" or "greater". The last two options only available for types 'rank', 'erl', 'cont' and 'area'.
<code>probs</code>	A two-element vector containing the lower and upper quantiles for the measure 'q' or 'qdir', in that order and on the interval [0, 1]. The default values are 0.025 and 0.975, suggested by Myllymäki et al. (2015, 2017).
<code>quantile.type</code>	As type argument of <a href="#">quantile</a> , how to calculate quantiles for 'q' or 'qdir'.
<code>central</code>	Either "mean" or "median". If the curve sets do not contain the component <code>theo</code> for the theoretical central function, then the central function (used for plotting only) is calculated either as the mean or median of functions provided in the curve sets. For 'qdir', 'st' and 'unscaled' only the mean is allowed as an option, due to their definition.

nstep	1 or 2 for how to construct a combined global envelope if list of curve sets is provided. 2 (default) for a two-step combining procedure, 1 for one-step.
...	Ignored.

## Details

Given a `curve_set` object, or an envelope object of `spatstat` or `fdata` object of `fda.usc`, the function `central_region` constructs a central region, i.e. a global envelope, from the given set of functions (or vectors).

Generally an envelope is a band bounded by the vectors (or functions)  $T_{low}$  and  $T_{hi}$ . A  $100(1 - \alpha)\%$  or `100*coverage%` global envelope is a set  $(T_{low}, T_{hi})$  of envelope vectors such that the probability that  $T_i$  falls outside this envelope in any of the  $d$  points of the vector  $T_i$  is less or equal to  $\alpha$ . The global envelopes can be constructed based on different measures that order the functions from the most extreme one to the least extreme one. We use such orderings of the functions for which we are able to construct global envelopes with intrinsic graphical interpretation.

The type of the global envelope can be chosen with the argument `type` and the options are given in the following. Further information about the measures, on which the global envelopes are based, can be found in Myllymäki and Mrkvíčka (2020, Section 2.).

- `'rank'`: The global rank envelope proposed by Myllymäki et al. (2017) based on the extreme rank defined as the minimum of pointwise ranks.
- `'erl'`: The global rank envelope based on the extreme rank length (Myllymäki et al., 2017, Mrkvíčka et al., 2018). This envelope is constructed as the convex hull of the functions which have extreme rank length measure that is larger or equal to the critical  $\alpha$  level of the extreme rank length measure.
- `'cont'`: The global rank envelope based on the continuous rank (Hahn, 2015; Mrkvíčka et al., 2019) based on minimum of continuous pointwise ranks. It is constructed as the convex hull in a similar way as the `'erl'` envelope.
- `'area'`: The global rank envelope based on the area rank (Mrkvíčka et al., 2019) which is based on area between continuous pointwise ranks and minimum pointwise ranks for those argument ( $r$ ) values for which pointwise ranks achieve the minimum (it is a combination of `erl` and `cont`). It is constructed as the convex hull in a similar way as the `'erl'` and `'area'` envelopes.
- `'qdir'`: The directional quantile envelope based on the directional quantile maximum absolute deviation (MAD) test (Myllymäki et al., 2017, 2015), which takes into account the unequal variances of the test function  $T(r)$  for different distances  $r$  and is also protected against asymmetry of distribution of  $T(r)$ .
- `'st'`: The studentised envelope based on the studentised MAD measure (Myllymäki et al., 2017, 2015), which takes into account the unequal variances of the test function  $T(r)$  for different distances  $r$ .
- `'unscaled'`: The unscaled envelope (Ripley, 1981), which leads to envelopes with constant width. It corresponds to the classical maximum deviation test without scaling. This test suffers from unequal variance of  $T(r)$  over the distances  $r$  and from the asymmetry of distribution of  $T(r)$ . We recommend to use the other alternatives instead. This unscaled global envelope is provided for reference.

The values of the chosen measure  $M$  are determined for each curve in the `curve_set`, and based on the chosen measure, the central region, i.e. the global envelope, is constructed for the given curves.

If a list of (suitable) objects are provided in the argument `curve_sets`, then by default (`nstep = 2`) the two-step combining procedure is used to construct the combined global envelope as described in Myllymäki and Mrkvička (2020, Section 2.2.). If `nstep = 1` and the lengths of the multivariate vectors in each component of the list are equal, then the one-step combining procedure is used where the functions are concatenated together into a one long vector (see again Myllymäki and Mrkvička, 2020, Section 2.2.).

## Value

Either an object of class `global_envelope` and or an `combined_global_envelope` object. The former class is obtained when a set of curves is provided, while the latter in the case that `curve_sets` is a list of objects. The `print` and `plot` function are defined for the returned objects (see examples).

The `global_envelope` object is essentially a data frame containing columns

- `r` = the vector of values of the argument `r` at which the test was made
- `lo` = the lower envelope based on the simulated functions; in case of a vector of coverage values, several 'lo' exist with names `paste0("lo.", 100*coverage)`
- `hi` = the upper envelope based on the simulated functions; in case of a vector of coverage values, several 'lo' exist with names `paste0("hi.", 100*coverage)`
- `central` = If the `curve_set` (or envelope object) contains a theoretical curve, then this function is used as the central curve and returned in this component. Otherwise, the central curve is the mean or median (according to the argument `central`) of the test functions  $T_i(r)$ ,  $i=2, \dots, s+1$ . Used for visualization only.

and potentially additionally

- `obs` = the data function, if there is only one data function in the given `curve_sets`. Otherwise not existing.

(Most often `central_region` is directly applied to functional data where all curves are observed.) Additionally, the returned object has some attributes, where

- `M` = A vector of the values of the chosen measure for all the function. If there is only one observed function, then `M[1]` gives the value of the measure for this.
- `M_alpha` = The critical value of `M` corresponding to the  $100(1-\alpha)\%$  global envelope (see Myllymäki and Mrkvička, 2020, Definition 1.1. IGI).

Further the object has some attributes for printing and plotting purposes, where `alternative`, `type`, `ties`, `alpha` correspond to those in the function call and `method` gives a name for the method. Attributes of an object `res` can be obtained using the function `attr`, e.g. `attr(res, "M")` for the values of the ordering measure.

If the given set of curves had the class envelope of `spatstat`, then the returned `global_envelope` object has also the class `fv` of `spatstat`, whereby one can utilize also the plotting functions of `spatstat`, see example in `plot.global_envelope`. However, the envelope objects are most often used with `global_envelope_test` and not with `central_region`. For an `fv` object, also some further attributes exists as required by `fv` of `spatstat`.

The `combined_global_envelope` is a list of `global_envelope` objects, where the components correspond to the components of `curve_sets`. The `combined_global_envelope` object constructed with `nstep = 2` contains, in addition to some conventional ones (`method`, `alternative`, `type`, `alpha`, `M`, `M_alpha`, see above), the second level envelope information as the attributes

- `level2_ge` = The second level envelope on which the envelope construction is based
- `level2_curve_set` = The second level `curve_set` from which `level2_ge` is constructed

In the case that the given curve sets are two-dimensional, i.e., their arguments values are two-dimensional, then the returned objects have in addition to the class `global_envelope` or `combined_global_envelope`, the class `global_envelope2d` or `combined_global_envelope2d`, respectively. This class is assigned for plotting purposes: For the 2d envelopes, also the default plots are 2d. Otherwise the 1d and 2d objects are similar.

## References

- Mrkvička, T., Myllymäki, M., Jilek, M. and Hahn, U. (2020) A one-way ANOVA test for functional data with graphical interpretation. *Kybernetika* 56(3), 432-458. doi: 10.14736/kyb-2020-3-0432
- Mrkvička, T., Myllymäki, M., Kuronen, M. and Narisetty, N. N. (2022) New methods for multiple testing in permutation inference for the general linear model. *Statistics in Medicine* 41(2), 276-297. doi: 10.1002/sim.9236
- Myllymäki, M., Grabarnik, P., Seijo, H. and Stoyan, D. (2015). Deviation test construction and power comparison for marked spatial point patterns. *Spatial Statistics* 11, 19-34. doi: 10.1016/j.spasta.2014.11.004
- Myllymäki, M., Mrkvička, T., Grabarnik, P., Seijo, H. and Hahn, U. (2017). Global envelope tests for spatial point patterns. *Journal of the Royal Statistical Society: Series B (Statistical Methodology)* 79, 381-404. doi: 10.1111/rssb.12172
- Myllymäki, M. and Mrkvička, T. (2024). GET: Global envelopes in R. *Journal of Statistical Software* 111(3), 1-40. doi: 10.18637/jss.v111.i03
- Ripley, B.D. (1981). *Spatial statistics*. Wiley, New Jersey.

## See Also

[forder](#), [global\\_envelope\\_test](#)

## Examples

```
## A central region of a set of functions
#-----
if(requireNamespace("fda", quietly=TRUE)) {
  cset <- curve_set(r=as.numeric(row.names(fda::growth$hgtf)),
                  obs=fda::growth$hgtf)
  plot(cset) + ggplot2::ylab("height")
  cr <- central_region(cset, coverage=0.50, type="erl")
  plot(cr)
}

## Confidence bands for linear or polynomial regression
#-----
# Simulate regression data according to the cubic model
```

```

# f(x) = 0.8x - 1.8x^2 + 1.05x^3 for x in [0,1]
par <- c(0,0.8,-1.8,1.05) # Parameters of the true polynomial model
res <- 100 # Resolution
x <- seq(0, 1, by=1/res); x2=x^2; x3=x^3;
f <- par[1] + par[2]*x + par[3]*x^2 + par[4]*x^3 # The true function
d <- f + rnorm(length(x), 0, 0.04) # Data
# Plot the true function and data
plot(f, type="l", ylim=range(d))
points(d)

# Estimate polynomial regression model
reg <- lm(d ~ x + x2 + x3)
ftheta <- reg$fitted.values
resid0 <- reg$residuals
s0 <- sd(resid0)

# Bootstrap regression
B <- 2000 # Number of bootstrap samples

ftheta1 <- array(0, c(B,length(x)))
s1 <- array(0,B)
for(i in 1:B) {
  u <- sample(resid0, size=length(resid0), replace=TRUE)
  reg1 <- lm((ftheta+u) ~ x + x2 + x3)
  ftheta1[i,] <- reg1$fitted.values
  s1[i] <- sd(reg1$residuals)
}

# Centering and scaling
meanftheta <- apply(ftheta1, 2, mean)
m <- array(0, c(B,length(x)))
for(i in 1:B) { m[i,] <- (ftheta1[i,]-meanftheta)/s1[i] }

# Central region computation
boot.cset <- curve_set(r=1:length(x), obs=ftheta+s0*t(m))
cr <- central_region(boot.cset, coverage=c(0.50, 0.80, 0.95), type="erl")

# Plotting the result
plot(cr) + ggplot2::labs(x=expression(italic(x)), y=expression(italic(f(x)))) +
  ggplot2::geom_point(data=data.frame(id=1:length(d), points=d),
    ggplot2::aes(x=id, y=points)) + # data points
  ggplot2::geom_line(data=data.frame(id=1:length(d), points=f),
    ggplot2::aes(x=id, y=points)) # true function

```

cgec

*Centred government expenditure centralization ratios***Description**

Centred government expenditure centralization (GEC) ratios

**Usage**

```
data("cgec")
```

**Format**

A list of two components. The first one is the `curve_set` object containing the observed values of centred GEC observed in year 1995-2016 for the above countries. The second component `group` gives the grouping.

**Details**

The data includes the government expenditure centralization (GEC) ratio in percent that has been centred with respect to country average in order to remove the differences in absolute values of GEC. The GEC ratio is the ratio of central government expenditure to the total general government expenditure. Data were collected from the Eurostat (2018) database. Only those European countries were included, where the data were available from 1995 to 2016 without interruption. Finally, 29 countries were classified into three groups in the following way:

- Group 1: Countries joining EC between 1958 and 1986 (Belgium, Denmark, France, Germany (until 1990 former territory of the FRG), Greece, Ireland, Italy, Luxembourg, Netherlands, Portugal, Spain, United Kingdom. These countries have long history of European integration, representing the core of integration process.
- Group 2: Countries joining the EU in 1995 (Austria, Sweden, Finland) and 2004 (Malta, Cyprus), except CEEC (separate group), plus highly economically integrated non-EU countries, EFTA members (Norway, Switzerland). Countries in this group have been, or in some case even still are standing apart from the integration mainstream. Their level of economic integration is however very high.
- Group 3: Central and Eastern European Countries (CEEC), having similar features in political and economic history. The process of economic and political integration have been initiated by political changes in 1990s. CEEC joined the EU in 2004 and 2007 (Bulgaria, Czech Republic, Estonia, Hungary, Latvia, Lithuania, Poland, Romania, Slovakia, Slovenia, data for Croatia joining in 2013 are incomplete, therefore not included).

This grouping is used in examples.

**References**

Eurostat (2018). "Government revenue, expenditure and main aggregates (gov10amain)". Retrieved from [https://ec.europa.eu/eurostat/data/database\(26/10/2018\)](https://ec.europa.eu/eurostat/data/database(26/10/2018)).

Mrkvička, T., Myllymäki, M., Jilek, M. and Hahn, U. (2020) A one-way ANOVA test for functional data with graphical interpretation. *Kybernetika* 56 (3), 432-458. doi: 10.14736/kyb-2020-3-0432

**See Also**

[graph.fanova](#)



## Examples

```
data("cgec")
# Plot data in groups
for(i in 1:3)
  assign(paste0("p", i), plot(subset(cgec$cgec, cgec$group == i)) +
    ggplot2::labs(title=paste("Group ", i, sep=""), y="Centred GEC"))
p3
if(require("patchwork", quietly=TRUE))
  p1 + p2 + p3
```

---

combined\_scaled\_MAD\_envelope\_test

*Combined global scaled maximum absolute difference (MAD) envelope tests*

---

## Description

Given a list of [curve\\_set](#) objects, a combined global scaled (directional quantile or studentized) MAD envelope test is performed with the test functions saved in the curve set objects. Details of this combined test can be found in Mrkvicka et al. (2017). The implementation of this test is provided here for historical reasons: we recommend now instead the use of [global\\_envelope\\_test](#) also for combined tests; these combined tests are there implemented as described in Myllymäki and Mrkvicka (2024).

## Usage

```
combined_scaled_MAD_envelope_test(
  curve_sets,
  type = c("qdir", "st"),
  alpha = 0.05,
  probs = c(0.025, 0.975),
  central = "mean",
  ...
)
```

## Arguments

curve_sets	A <a href="#">curve_set</a> object or a list of <a href="#">curve_set</a> objects containing a data function and simulated functions from which the envelope is to be constructed. Also envelope objects of <b>spatstat</b> are accepted instead of curve_set objects. If an envelope object is given, it must contain the summary functions from simulated patterns which can be achieved by setting savefuns = TRUE when calling the envelope function.
type	Either "qdir" for the direction quantile envelope test or "st" for the studentized envelope test.

alpha	The significance level. The 100(1-alpha)% global envelope will be calculated under the 'fwer' or 'fdr' control. If a vector of values is provided, the global envelopes are calculated for each value.
probs	A two-element vector containing the lower and upper quantiles for the measure 'q' or 'qdir', in that order and on the interval [0, 1]. The default values are 0.025 and 0.975, suggested by Myllymäki et al. (2015, 2017).
central	Either "mean" or "median". If the curve sets do not contain the component theo for the theoretical central function, then the central function (used for plotting only) is calculated either as the mean or median of functions provided in the curve sets. For 'qdir', 'st' and 'unscaled' only the mean is allowed as an option, due to their definition.
...	Additional parameters to be passed to <a href="#">central_region</a> .

## References

- Mrkvička, T., Myllymäki, M. and Hahn, U. (2017) Multiple Monte Carlo testing, with applications in spatial point processes. *Statistics & Computing* 27(5): 1239–1255. DOI: 10.1007/s11222-016-9683-9
- Myllymäki, M. and Mrkvička, T. (2024). GET: Global envelopes in R. *Journal of Statistical Software* 111(3), 1-40. doi: 10.18637/jss.v111.i03

## Examples

```
if(require("spatstat.explore", quietly=TRUE)) {
  # As an example test CSR of the saplings point pattern from spatstat by means of
  # L, F, G and J functions.
  data("saplings")
  X <- as.ppp(saplings, W=square(75))

  nsim <- 499 # Number of simulations for the tests

  # Specify distances for different test functions
  n <- 500 # the number of r-values
  rmin <- 0; rmax <- 20; rstep <- (rmax-rmin)/n
  rminJ <- 0; rmaxJ <- 8; rstepJ <- (rmaxJ-rminJ)/n
  r <- seq(0, rmax, by=rstep) # r-distances for Lest
  rJ <- seq(0, rmaxJ, by=rstepJ) # r-distances for Fest, Gest, Jest

  # Perform simulations of CSR and calculate the L-functions
  env_L <- envelope(X, nsim=nsim,
    simulate=expression(runifpoint(ex=X)),
    fun="Lest", correction="translate",
    transform=expression(.-r), # Take the L(r)-r function instead of L(r)
    r=r, # Specify the distance vector
    savefuns=TRUE, # Save the estimated functions
    savepatterns=TRUE) # Save the simulated patterns
  # Take the simulations from the returned object
  simulations <- attr(env_L, "simpatterns")
  # Then calculate the other test functions F, G, J for each simulated pattern
```

```

env_F <- envelope(X, nsim=nsim,
  simulate=simulations,
  fun="Fest", correction="Kaplan", r=rJ,
  savefuns=TRUE)
env_G <- envelope(X, nsim=nsim,
  simulate=simulations,
  fun="Gest", correction="km", r=rJ,
  savefuns=TRUE)
env_J <- envelope(X, nsim=nsim,
  simulate=simulations,
  fun="Jest", correction="none", r=rJ,
  savefuns=TRUE)

# Crop the curves to the desired r-interval I
curve_set_L <- crop_curves(env_L, r_min=rmin, r_max=rmax)
curve_set_F <- crop_curves(env_F, r_min=rminJ, r_max=rmaxJ)
curve_set_G <- crop_curves(env_G, r_min=rminJ, r_max=rmaxJ)
curve_set_J <- crop_curves(env_J, r_min=rminJ, r_max=rmaxJ)

# The combined directional quantile envelope test
res <- combined_scaled_MAD_envelope_test(
  curve_sets=list(L=curve_set_L, F=curve_set_F,
    G=curve_set_G, J=curve_set_J),
  type="qdir")
plot(res)
}

```

---

create\_image\_set

---

*Create a curve set of images*


---

## Description

Create a curve set consisting of a set of images, given a list containing the values of the 2d functions in the right form. Only 2d functions in a rectangular windows are supported; the values are provided in matrices (arrays). For more general 2d functions see [create\\_curve\\_set](#).

## Usage

```
create_image_set(image_set, ...)
```

## Arguments

image_set	A list containing elements <code>r</code> , <code>obs</code> , <code>sim_m</code> and <code>theo</code> . <code>r</code> , <code>sim_m</code> and <code>theo</code> are optional, <code>obs</code> needs to be provided always. If provided, <code>r</code> must be a list describing the argument values where the images have been observed (or simulated). It must consist of the following two or four components: a) "x" and "y" giving the equally spaced argument values for the x- and y-coordinates (first and second dimension of the 2d functions) where the data have been observed, b) "x", "y",
-----------	---

"width" and "height", where the width and height give the width and height of the pixels placed at x and y, or c) "xmin", "xmax", "ymin" and "ymax" giving the corner coordinates of the pixels where the data have been observed. If not given, r is set to be a list of values from 1 to the number of first/second dimension of 2d functions in obs. obs must be either a 2d matrix (dimensions matching the lengths of r vectors) or 3d array containing the observed 2d functions (the third dimension matching the number of functions). If obs is a 3d array, then sim\_m is ignored. If obs is a 2d array, then sim\_m must be a 3d array containing the simulated images (2d functions) (the third dimension matching the number of functions). If included, theo corresponds to the theoretical function (e.g., under the null hypothesis) and its dimensions must either match the dimensions of 2d functions in obs or it must be a constant.

... Do not use. (For internal use only.)

### Value

The given list as a curve\_set.

### Examples

```
a <- create_image_set(list(obs=array(runif(4*5*6), c(4,5,6))))
plot(a)
plot(a, idx=1:6)

a <- create_image_set(list(r=list(x=c(10,20,30,40), y=1:5*0.1),
                                obs=array(runif(4*5*6), c(4,5,6))))
plot(a)

a <- create_image_set(list(r=list(xmin=c(1, 2, 4, 7), xmax=c(2, 4, 7, 11),
                                ymin=c(1,1.1,2,2.1,3), ymax=c(1.1,2,2.1,3,3.1)),
                                obs=array(runif(4*5*6), c(4,5,6))))
plot(a)
plot(a, idx=1:5)
```

---

crop\_curves

*Crop the curves*

---

### Description

Crop the curves to a certain interval, or crop missing and infinite argument values from the curves

### Usage

```
crop_curves(curve_set, allfinite = TRUE, r_min = NULL, r_max = NULL)
```

**Arguments**

curve_set	A <a href="#">curve_set</a> object, or an envelope object of <b>spatstat</b> . If an envelope object is given, it must contain the summary functions from the simulated patterns which can be achieved by setting <code>savefuns = TRUE</code> when calling the envelope function.
allfinite	Logical. TRUE means that the argument values where any of the curves have missing or infinite values are removed. FALSE means that only <code>r_min</code> and <code>r_max</code> apply.
r_min	The minimum radius to include.
r_max	The maximum radius to include.

**Details**

The curves can be cropped to a certain interval defined by the arguments `r_min` and `r_max`. Also the argument values of the sets of curves which have missing or infinite values for any of the curves can be removed from the set (`allfinite = TRUE`). The interval should generally be chosen carefully for classical deviation tests.

**Value**

A `curve_set` object containing the cropped summary functions and the cropped radius vector.

---

curve_set	<i>Create a curve_set object</i>
-----------	----------------------------------

---

**Description**

Create a `curve_set` object out of data provided in the right form.

**Usage**

```
curve_set(
  obs,
  sim = NULL,
  r = NULL,
  theo = NULL,
  allfinite = FALSE,
  verbose = TRUE
)

create_curve_set(curve_set, allfinite = FALSE, verbose = TRUE)
```

## Arguments

obs	The observed data. See details.
sim	The simulated data. See details.
r	The argument values where the functions/vectors have been observed (or simulated). See details.
theo	The theoretical function. See details.
allfinite	Logical. TRUE requires that all values of the curves must be finite (not infinite and not missing, see <a href="#">is.finite</a> ). FALSE allows for infinite or missing values in the curves. These infinite and missing values can have consequences for the subsequent analyses. A warning is given if infinite or missing values exists (when verbose = TRUE).
verbose	Logical flag indicating whether to warn about the content.
curve_set	A list containing the element obs, and optionally the elements r, sim_m (same as sim, but in matrix format only) and theo. See details.

## Details

The functions are used to clump together the functional data in the form that can be handled by the other **GET** functions ([forder](#), [central\\_region](#), [global\\_envelope\\_test](#) etc.). The functions take care of checking the content of the data, and save relevant information of the curves for global envelope methods to be used in particular for plotting the results with graphical interpretation.

obs must be either

- a vector containing the data function/vector, or
- a matrix containing the s data functions/vectors, in which case it is assumed that each column corresponds to a data function/vector, or
- a list containing the s data functions/vectors.

If obs is a vector, sim must be either

- a matrix containing the simulated functions/vectors, each column corresponding to a function/vector and the number of rows matching the length of obs, or
- a list containing the simulated functions/vectors.

If obs is a matrix or a list, sim is ignored.

If given, r describes the 1- or 2-dimensional argument values where the functions/vectors have been observed (or simulated). It must be either

- a vector,
- a data.frame with columns "x", "y", "width" and "height", where the width and height give the width and height of the pixels placed at x and y, or
- a data.frame with columns "xmin", "xmax", "ymin" and "ymax" giving the corner coordinates of the pixels where the data have been observed.

If obs is a vector, theo can be given and it should then correspond to a theoretical function (e.g., under the null hypothesis). If present, its length must match the length of obs.

**Value**

An object of class `curve_set` containing the data. If the argument values are two-dimensional, then the `curve_set` is additionally a `curve_set2d` object.

**See Also**

[plot.curve\\_set](#), [plot.curve\\_set2d](#)

**Examples**

```
# 1d
cset <- curve_set(r = 1:10, obs = matrix(runif(10*5), ncol=5))
plot(cset)
# 2d
cset <- curve_set(r = data.frame(x=c(rep(1:3, 3), 4), y=c(rep(1:3, each=3), 1),
                                width=1, height=1),
                  obs = matrix(runif(10*5), ncol=5))
plot(cset)
```

---

deviation\_test

*Deviation test*


---

**Description**

Crop the curve set to the interval of distances `[r_min, r_max]`, calculate residuals, scale the residuals and perform a deviation test with a chosen deviation measure. The deviation tests are well known in spatial statistics; in **GET** they are provided for comparative purposes. Some (maximum type) of the deviation test have their corresponding envelope tests available, see Myllymäki et al., 2017 (and 'unscaled', 'st' and 'qdir' in [global\\_envelope\\_test](#)).

**Usage**

```
deviation_test(
  curve_set,
  r_min = NULL,
  r_max = NULL,
  use_theo = TRUE,
  scaling = "qdir",
  measure = "max",
  alternative = c("two.sided", "less", "greater"),
  savedevs = FALSE
)
```

### Arguments

curve_set	A residual curve_set object. Can be obtained by using residual().
r_min	The minimum radius to include.
r_max	The maximum radius to include.
use_theo	Whether to use the theoretical summary function or the mean of the functions in the curve_set.
scaling	The name of the scaling to use. Options include 'none', 'q', 'qdir' and 'st'. 'qdir' is default.
measure	The deviation measure to use. Default is 'max'. Must be one of the following: 'max', 'int' or 'int2'.
alternative	A character string specifying the alternative hypothesis when measure = 'max'; otherwise ignored. Must be one of the following: "two.sided" (default), "less" or "greater".
savedevs	Logical. Should the global rank values k_i, i=1,...,nsim+1 be returned? Default: FALSE.

### Details

The deviation test is based on a test function  $T(r)$  and it works as follows:

1) The test function estimated for the data,  $T_1(r)$ , and for nsim simulations from the null model,  $T_2(r), \dots, T_{nsim+1}(r)$ , must be saved in 'curve\_set' and given to the deviation\_test function.

2) The deviation\_test function then

- Crops the functions to the chosen range of distances  $[r_{\min}, r_{\max}]$ .
- If the curve\_set does not consist of residuals (see [residual](#)), then the residuals  $d_i(r) = T_i(r) - T_0(r)$  are calculated, where  $T_0(r)$  is the expectation of  $T(r)$  under the null hypothesis. If use\_theo = TRUE, the theoretical value given in the curve\_set\$theo is used for as  $T_0(r)$ , if it is given. Otherwise,  $T_0(r)$  is estimated by the mean of  $T_j(r)$ ,  $j = 2, \dots, nsim + 1$ .
- Scales the residuals. Options are
  - 'none' No scaling. Nothing done.
  - 'q' Quantile scaling.
  - 'qdir' Directional quantile scaling.
  - 'st' Studentised scaling.

See for details Myllymäki et al. (2013).

- Calculates the global deviation measure  $u_i$ ,  $i = 1, \dots, nsim + 1$ , see options for 'measure'.
  - 'max' is the maximum deviation measure

$$u_i = \max_{r \in [r_{\min}, r_{\max}]} |w(r)(T_i(r) - T_0(r))|$$

If alternative = "greater", then instead

$$u_i = \max_{r \in [r_{\min}, r_{\max}]} [w(r)(T_i(r) - T_0(r))]$$

i.e. the largest values will have the largest  $u_i$ .



If alternative = "less", then instead

$$u_i = \max_{r \in [r_{\min}, r_{\max}]} [-w(r)(T_i(r) - T_0(r))]$$

i.e. the smallest values will have the largest  $u_i$ .

- 'int2' is the integral deviation measure

$$u_i = \int_{r_{\min}}^{r_{\max}} (w(r)(T_i(r) - T_0(r)))^2 dr$$

- 'int' is the 'absolute' integral deviation measure

$$u_i = \int_{r_{\min}}^{r_{\max}} |w(r)(T_i(r) - T_0(r))| dr$$

- Calculates the p-value.

## Value

If 'savedevs=FALSE' (default), the p-value is returned. If 'savedevs=TRUE', then a list containing the p-value and calculated deviation measures  $u_i$ ,  $i = 1, \dots, \text{nsim} + 1$  (where  $u_1$  corresponds to the data pattern) is returned.

## References

- Myllymäki, M., Grabarnik, P., Seijo, H. and Stoyan, D. (2015). Deviation test construction and power comparison for marked spatial point patterns. *Spatial Statistics* 11: 19-34. doi: 10.1016/j.spasta.2014.11.004
- Myllymäki, M., Mrkvička, T., Grabarnik, P., Seijo, H. and Hahn, U. (2017). Global envelope tests for spatial point patterns. *Journal of the Royal Statistical Society: Series B (Statistical Methodology)*, 79: 381–404. doi: 10.1111/rssb.12172

## Examples

```
## Testing complete spatial randomness (CSR)
#-----
if(require("spatstat.explore", quietly=TRUE)) {
  pp <- unmark(spruces)
  nsim <- 999

  # Generate nsim simulations under CSR, calculate L-function for the data and simulations
  env <- envelope(pp, fun="Lest", nsim=nsim, savefuns=TRUE, correction="translate")
  # The deviation test using the integral deviation measure
  res <- deviation_test(env, measure='int')
  res
  # or
  res <- deviation_test(env, r_min=0, r_max=7, measure='int2')
}
```

---

 fallen\_trees

*Fallen trees*


---

### Description

Fallen trees

### Usage

```
data("fallen_trees")
```

### Format

A **list** of two data frames, where `trees` contains the locations (x and y coordinates) and heights (=marks) of 232 trees in a window with polygonal boundary, and `window` species the polygonal window (see examples).

### Details

The dataset comprised the locations and heights of 232 trees, which fell during two large wind gusts (1967 and 1990) in the west of France (Pontailler et al., 1997). The study area was a biological reserve, which had been preserved for at least four centuries, with little human influence for a long period (Guinier, 1950). Thus, the forest stand followed almost natural dynamics. It was an uneven-aged beech stand with a few old oaks.

The data was analysed in Myllymäki et al. (2017, Supplementary material).

### References

Guinier, P. (1950) Foresterie et protection de la nature. l'exemple de fontainebleau. *Rev Forestière Fr.*, II, 703-717.

Pontailler, J.-Y., Faille, A. and Lemée, G. (1997) Storms drive successional dynamics in natural forests: a case study in fontainebleau forest (france). *Forest Ecol. Manag.*, 98, 1-15.

Myllymäki, M., Mrkvička, T., Grabarnik, P., Seijo, H. and Hahn, U. (2017). Global envelope tests for spatial point patterns. *Journal of the Royal Statistical Society: Series B (Statistical Methodology)*, 79: 381–404. doi: 10.1111/rssb.12172

### Examples

```
data("fallen_trees")
if(require("spatstat.geom", quietly=TRUE)) {
  fallen_trees <- as.ppp(fallen_trees$trees, W = owin(poly=fallen_trees>window))
  plot(fallen_trees)
}
```

fBoxplot

*Functional boxplot*

## Description

Functional boxplot based on central region computed by a specified measure. The options of the measures can be found in [central\\_region](#).

## Usage

```
fBoxplot(curve_sets, factor = 1.5, coverage = 0.5, ...)
```

## Arguments

curve_sets	A <a href="#">curve_set</a> object or a list of <a href="#">curve_set</a> objects. Also envelope objects of <b>spatstat</b> and fdata of <b>fda.usc</b> are accepted instead of curve_set objects.
factor	The constant factor to inflate the central region to produce a functional boxplot and determine fences for outliers. Default is 1.5 as in a classical boxplot.
coverage	A number between 0 and 1. The 100*coverage% central region will be calculated. A vector of values can also be provided, leading to the corresponding number of central regions.
...	Additional parameters to be passed to <a href="#">central_region</a> , which is responsible for calculating the central region (global envelope) on which the functional boxplot is based.

## Examples

```
if(requireNamespace("fda", quietly=TRUE)) {
  years <- paste(1:18)
  curves <- fda::growth[['hgtf']][years,]
  # Heights
  cset1 <- curve_set(r = as.numeric(years),
                    obs = curves)
  bp <- fBoxplot(cset1, coverage=0.50, type="area", factor=1)
  plot(bp)

  # Considering simultaneously heights and height differences
  cset2 <- curve_set(r = as.numeric(years[-1]),
                    obs = curves[-1,] - curves[-nrow(curves),])
  csets <- list(Height=cset1, Change=cset2)
  res <- fBoxplot(csets, type='area', factor=1.5)
  plot(res) + ggplot2::labs(x="Age (years)", y="")
}
```

---

fclustering	<i>Functional clustering</i>
-------------	------------------------------

---

## Description

Functional clustering based on a specified measure. The options of the measures can be found in [central\\_region](#).

## Usage

```
fclustering(
  curve_sets,
  k,
  type = c("area", "st", "erl", "cont"),
  triangineq = FALSE,
  ...
)
```

## Arguments

<code>curve_sets</code>	A <code>curve_set</code> object or a list of <code>curve_set</code> objects to which the functional clustering is to be applied. If list of <code>curve_set</code> objects is provided, then the joined functional clustering is applied, which provides an equal weight combination of <code>curve_set</code> objects, if the <code>curve_set</code> objects contain the same numbers of elements (same lengths of vector $r$ ).
<code>k</code>	The number of clusters.
<code>type</code>	The measure which is used to compute the dissimilarity matrix. The preferred options are "area" and "st", but "erl" and "cont" can be also used with caution.
<code>triangineq</code>	Logical. Whether or not to compute the proportion of combinations of functions which satisfies the triangular inequality, see 'Value'.
<code>...</code>	Additional parameters to be passed to <a href="#">central_region</a> , which is responsible for calculating the central region (global envelope) on which the functional clustering is based.

## Details

Functional clustering joins the list of `curve_set` objects in one `curve_set` with long functions and applies on the differences of all functions the specified measure. This provides a dissimilarity matrix which is used in partitioning around medoids procedure. The resulting clusters can then be shown by plotting the function respectively for each `curve_set`. Thus for each `curve_set`, the panel with all the medoids is shown followed by all clusters represented by central region, medoid and all curves belonging to it, when the result object is plotted.

If there are less than three curves in some of the groups, then the central region is not plotted. This leads to a warning message from `ggplot2`.

**Value**

An object having the class `fclust`, containing

- `curve_sets` = The set(s) of functions determined for clustering
- `k` = Number of clusters
- `type` = Type of clustering method
- `triangineq` = The proportion of combinations of functions which satisfies the triangular inequality. The triangular inequality must hold to ensure the chosen measure forms a metric. In some weird cases it does not hold for 'area' measure, therefore this check is provided to ensure the data forms metric with the 'area' measure. The `triangineq` must be 1 to ensure the inequality holds for all functions.
- `dis` = The joined dissimilarity matrix
- `pam` = Results of the partitioning around medoids (pam) method applied on the joined functions with the dissimilarity matrix (`dis`). See [pam](#).

**References**

Dai, W., Athanasiadis, S., Mrkvička, T. (2021) A new functional clustering method with combined dissimilarity sources and graphical interpretation. Intech open, London, UK. DOI: 10.5772/intechopen.100124

**See Also**

[central\\_region](#), [plot.fclust](#)

**Examples**

```
# Read raw data from population growth rdata
# with countries over million inhabitants
data("popgrowthmillion")

# Create centred data
m <- apply(popgrowthmillion, 2, mean) # Country-wise means
cpopgrowthmillion <- popgrowthmillion
for(i in 1:dim(popgrowthmillion)[1]) {
  cpopgrowthmillion[i,] <- popgrowthmillion[i,] - m
}

# Create scaled data
t2 <- function(v) { sqrt(sum(v^2)) }
s <- apply(cpopgrowthmillion, 2, t2)
spopgrowthmillion <- popgrowthmillion
for(i in 1:dim(popgrowthmillion)[1]) {
  spopgrowthmillion[i,] <- cpopgrowthmillion[i,]/s
}

# Create curve sets
r <- 1951:2015
```

```

cset1 <- curve_set(r = r, obs = popgrowthmillion)
cset2 <- curve_set(r = r, obs = spopgrowthmillion)
csets <- list(Raw = cset1, Shape = cset2)

# Functional clustering with respect to joined "st" difference measure
# and "joined" central regions of each group
res <- fclustering(csets, k=3, type="area")
p <- plot(res, plotstyle = "marginal", coverage = 0.5)
p[[1]] # Central functions
p[[2]] # Groups: central functions and regions
# To collect the two figures into one use, e.g., patchwork:
if(require("patchwork", quietly=TRUE)) {
  p[[1]] + p[[2]] + plot_layout(widths = c(1, res$k))
}
# Silhouette plot of pam
plot(res$pam)

```

fdr\_envelope

*The FDR envelope*

## Description

Calculate the FDR envelope based on the ATSE or IATSE algorithm of Mrkvička and Myllymäki (2023).

## Usage

```

fdr_envelope(
  curve_sets,
  alpha = 0.05,
  alternative = c("two.sided", "less", "greater"),
  algorithm = c("IATSE", "ATSE"),
  lower = NULL,
  upper = NULL
)

```

## Arguments

curve_sets	A <a href="#">curve_set</a> object or a list of <a href="#">curve_set</a> objects containing a data function and simulated functions from which the envelope is to be constructed. Also envelope objects of <b>spatstat</b> are accepted instead of curve_set objects. If an envelope object is given, it must contain the summary functions from simulated patterns which can be achieved by setting savefuncs = TRUE when calling the envelope function.
alpha	The significance level. The 100(1-alpha)% global envelope will be calculated under the 'fwer' or 'fdr' control. If a vector of values is provided, the global envelopes are calculated for each value.

alternative	A character string specifying the alternative hypothesis. Must be one of the following: "two.sided" (default), "less" or "greater". The last two options only available for types 'rank', 'erl', 'cont' and 'area'.
algorithm	The algorithm for the computation of the FDR envelope. Either "IATSE" or "ATSE" standing for the iteratively adaptive two-stage envelope and the adaptive two-stage envelope, respectively, see Mrkvička and Myllymäki (2023).
lower	A single number (or a vector of suitable length) giving a lower bound for the functions. Used only for the extension of the FDR envelope.
upper	A single number (or a vector of suitable length) giving an upper bound for the functions. Used only for the extension of the FDR envelope.

### Details

Typical use of this function is through other functions. `fdr_envelope(cset)` is the same as `global_envelope_test(cset, typeone = "fdr")`. Functions such as [graph.fanova](#), [graph.flm](#), [frank.flm](#) allow to use the FDR control by specifying `typeone = "fdr"` appropriately (passing this to `global_envelope_test`).

### References

Mrkvička and Myllymäki (2023). False discovery rate envelopes. *Statistics and Computing* 33, 109. <https://doi.org/10.1007/s11222-023-10275-7>

### Examples

```
# A GLM example
data(rimov)
nsim <- 1000 # Number of simulations

res <- graph.flm(nsim=nsim,
  formula.full = Y~Year,
  formula.reduced = Y~1,
  curve_sets = list(Y=rimov),
  factors = data.frame(Year = 1979:2014),
  GET.args = list(typeone = "fdr"))

plot(res)
```

### Description

Calculates different measures for ordering the functions (or vectors) from the most extreme to least extreme one

## Usage

```
forder(
  curve_sets,
  measure = "erl",
  scaling = "qdir",
  alternative = c("two.sided", "less", "greater"),
  use_theo = TRUE,
  probs = c(0.025, 0.975),
  quantile.type = 7
)
```

## Arguments

curve_sets	A <a href="#">curve_set</a> object or a list of <a href="#">curve_set</a> objects. Also envelope objects of <b>spatstat</b> and fdata of <b>fda.usc</b> are accepted instead of curve_set objects.
measure	The measure to use to order the functions from the most extreme to the least extreme one. Must be one of the following: 'rank', 'erl', 'cont', 'area', 'max', 'int', 'int2'. Default is 'erl'.
scaling	The name of the scaling to use if measure is 'max', 'int' or 'int2'. Options include 'none', 'q', 'qdir' and 'st', where 'qdir' is the default.
alternative	A character string specifying the alternative hypothesis. Must be one of the following: "two.sided" (default), "less" or "greater". The last two options only available for types 'rank', 'erl', 'cont' and 'area'.
use_theo	Logical. When calculating the measures 'max', 'int', 'int2', should the theoretical function from curve_set be used (if 'theo' provided), see <a href="#">deviation_test</a> .
probs	A two-element vector containing the lower and upper quantiles for the measure 'q' or 'qdir', in that order and on the interval [0, 1]. The default values are 0.025 and 0.975, suggested by Myllymäki et al. (2015, 2017).
quantile.type	As type argument of <a href="#">quantile</a> , how to calculate quantiles for 'q' or 'qdir'.

## Details

Given a [curve\\_set](#) object or an envelope object of **spatstat**, which contains curves  $T_1(r), \dots, T_s(r)$ , the functions are ordered from the most extreme one to the least extreme one by one of the following measures (specified by the argument measure). Note that 'erl', 'cont' and 'area' were proposed as a refinement to the extreme ranks 'rank', because the extreme ranks can contain many ties. All of these completely non-parametric measures are smallest for the most extreme functions and largest for the least extreme ones, whereas the deviation measures ('max', 'int' and 'int2') obtain largest values for the most extreme functions.

- 'rank': extreme rank (Myllymäki et al., 2017). The extreme rank  $R_i$  is defined as the minimum of pointwise ranks of the curve  $T_i(r)$ , where the pointwise rank is the rank of the value of the curve for a specific r-value among the corresponding values of the  $s$  other curves such that the lowest ranks correspond to the most extreme values of the curves. How the pointwise ranks are determined exactly depends on the whether a one-sided (alternative is "less" or "greater") or the two-sided test (alternative="two.sided") is chosen.



- 'erl': extreme rank length (Myllymäki et al., 2017). Considering the vector of pointwise ordered ranks  $\mathbf{R}_i$  of the  $i$ th curve, the extreme rank length measure  $R_i^{erl}$  is equal to

$$R_i^{erl} = \frac{1}{s} \sum_{j=1}^s \mathbf{1}(\mathbf{R}_j \prec \mathbf{R}_i)$$

where  $\mathbf{R}_j \prec \mathbf{R}_i$  if and only if there exists  $n \leq d$  such that for the first  $k$ ,  $k < n$ , pointwise ordered ranks of  $\mathbf{R}_j$  and  $\mathbf{R}_i$  are equal and the  $n$ 'th rank of  $\mathbf{R}_j$  is smaller than that of  $\mathbf{R}_i$ . The scaling by

$s$

is applied to normalize the ranks following Mrkvička et al. (2019) and Narisetty and Nair (2016).

- 'cont': continuous rank (Hahn, 2015; Mrkvička et al., 2019) based on minimum of continuous pointwise ranks
- 'area': area rank (Mrkvička et al., 2019) based on area between continuous pointwise ranks and minimum pointwise ranks for those argument ( $r$ ) values for which pointwise ranks achieve the minimum (it is a combination of erl and cont)
- 'max' and 'int' and 'int2': Further options for the measure argument that can be used together with scaling. See the help in [deviation\\_test](#) for these options of measure and scaling. These measures are largest for the most extreme functions and smallest for the least extreme ones. The arguments use\_theo and probs are relevant for these measures only (otherwise ignored).

For details see Myllymäki and Mrkvička et al. (2020, Section 2)

## Value

A vector containing one of the above mentioned measures  $k$  for each of the functions in the curve set. If the component obs in the curve set is a vector, then its measure will be the first component (named 'obs') in the returned vector.

## References

- Hahn U (2015). "A note on simultaneous Monte Carlo tests." Technical report, Centre for Stochastic Geometry and advanced Bioimaging, Aarhus University.
- Mrkvička, T., Myllymäki, M., Jilek, M. and Hahn, U. (2020) A one-way ANOVA test for functional data with graphical interpretation. *Kybernetika* 56(3), 432-458. doi: 10.14736/kyb-2020-3-0432
- Mrkvička, T., Myllymäki, M., Kuronen, M. and Narisetty, N. N. (2022) New methods for multiple testing in permutation inference for the general linear model. *Statistics in Medicine* 41(2), 276-297. doi: 10.1002/sim.9236
- Myllymäki, M., Grabarnik, P., Seijo, H. and Stoyan, D. (2015). Deviation test construction and power comparison for marked spatial point patterns. *Spatial Statistics* 11, 19-34. doi: 10.1016/j.spasta.2014.11.004
- Myllymäki, M., Mrkvička, T., Grabarnik, P., Seijo, H. and Hahn, U. (2017). Global envelope tests for spatial point patterns. *Journal of the Royal Statistical Society: Series B (Statistical Methodology)* 79, 381-404. doi: 10.1111/rssb.12172
- Narisetty, N. N. and Nair, V. J. (2016) Extremal depth for functional data and applications. *Journal of the American Statistical Association* 111, 1705-1714.

**See Also**[partial\\_forder](#)**Examples**

```

if(requireNamespace("fda", quietly = TRUE)) {
  # Consider ordering of the girls in the Berkeley Growth Study data
  # available from the R package fda, see ?growth, according to their
  # annual heights or/and changes within years.
  # First create sets of curves (vectors), for raw heights and
  # for the differences within the years
  years <- paste(1:18)
  curves <- fda::growth[['hgtf']][years,]
  cset1 <- curve_set(r = as.numeric(years),
                    obs = curves)
  cset2 <- curve_set(r = as.numeric(years[-1]),
                    obs = curves[-1,] - curves[-nrow(curves),])

  # Order the girls from most extreme one to the least extreme one, below using the 'area' measure
  # a) according to their heights
  forder(cset1, measure = 'area')
  # Print the 10 most extreme girl indices
  order(forder(cset1, measure = 'area'))[1:10]
  # b) according to the changes (print indices)
  order(forder(cset2, measure = 'area'))[1:10]
  # c) simultaneously with respect to heights and changes (print indices)
  csets <- list(Height = cset1, Change = cset2)
  order(forder(csets, measure = 'area'))[1:10]
}

```

frank.fanova

*Rank envelope F-test***Description**

A one-way functional ANOVA based on the rank envelope applied to F values

**Usage**

```

frank.fanova(
  nsim,
  curve_set,
  groups,
  variances = "equal",
  test.equality = c("mean", "var", "cov"),
  cov.lag = 1,
  savefuncs = FALSE,
  ...
)

```

**Arguments**

<code>nsim</code>	The number of random permutations.
<code>curve_set</code>	The original data (an array of functions) provided as a <a href="#">curve_set</a> object or a <code>fdata</code> object of <b>fda.usc</b> . The curve set should include the argument values for the functions in the component <code>r</code> , and the observed functions in the component <code>obs</code> .
<code>groups</code>	The original groups (a factor vector representing the assignment to groups).
<code>variances</code>	Either "equal" or "unequal". If "equal", then the traditional F-values are used. If "unequal", then the corrected F-values are used. The current implementation uses <a href="#">lm</a> to get the corrected F-values.
<code>test.equality</code>	A character with possible values mean (default), var and cov. If mean, the functional ANOVA is performed to compare the means in the groups. If var, then the equality of variances of the curves in the groups is tested by performing the graphical functional ANOVA test on the functions $Z_{ij}(r) = T_{ij}(r) - \bar{T}_j(r).$ <p>If cov, then the equality of lag cov.lag covariance is tested by performing the fANOVA with</p> $W_{ij}(r) = \sqrt{ V_{ij}(r)  \cdot \text{sign}(V_{ij}(r))},$ <p>where</p> $V_{ij}(r) = (T_{ij}(r) - \bar{T}_j(r))((T_{ij}(r + s) - \bar{T}_j(r + s))).$ <p>See Mrkvicka et al. (2020) for more details.</p>
<code>cov.lag</code>	The lag of the covariance for testing the equality of covariances, see <code>test.equality</code> .
<code>savefuns</code>	Logical. If TRUE, then the functions from permutations are saved to the attribute <code>simfuns</code> .
<code>...</code>	Additional parameters to be passed to <a href="#">global_envelope_test</a> . For example, the type of multiple testing control, FWER or FDR must be set by <code>typeone</code> . And, if <code>typeone = "fwer"</code> , the type of the global envelope can be chosen by specifying the argument <code>type</code> . See <a href="#">global_envelope_test</a> for the defaults and available options. (The test here uses <code>alternative="two.sided"</code> and <code>nstep=1</code> (when relevant), but all the other specifications are to be given in ...)

**Details**

The test assumes that there are  $J$  groups which contain  $n_1, \dots, n_J$  functions  $T_{ij}, i = \dots, J, j = 1, \dots, n_j$ . The functions should be given in the argument `x`, and the groups in the argument `groups`. The test assumes that there exists non random functions  $\mu(r)$  and  $\mu_i(r)$  such that

$$T_{ij}(r) = \mu(r) + \mu_i(r) + e_{ij}(r), i = 1, \dots, J, j = 1, \dots, n_j$$

where  $e_{ij}(r)$  are independent and normally distributed. The test vector is

$$\mathbf{T} = (F(r_1), F(r_2), \dots, F(r_K)),$$

where  $F(r_i)$  stands for the F-statistic. The simulations are performed by permuting the test functions. Further details can be found in Mrkvicka et al. (2020).

The argument `variances="equal"` means that equal variances across groups are assumed. The correction for unequal variances can be done by using the corrected F-statistic (option `variances="unequal"`).

Unfortunately this test is not able to detect which groups are different from each other.

## References

Mrkvička, T., Myllymäki, M., Jilek, M. and Hahn, U. (2020) A one-way ANOVA test for functional data with graphical interpretation. *Kybernetika* 56 (3), 432-458. doi: 10.14736/kyb-2020-3-0432

## See Also

`graph.fanova`

## Examples

```
data("rimov")
groups <- factor(c(rep(1, times=12), rep(2, times=12), rep(3, times=12)))
res <- frank.fanova(nsim = 2499, curve_set = rimov, groups = groups)

plot(res)

data("imageset3")
res2 <- frank.fanova(nsim = 19, # Increase nsim for serious analysis!
                    curve_set = imageset3$image_set,
                    groups = imageset3$Group)

plot(res2)
plot(res2, fixed scales=FALSE)
```

---

frank.flm

*F rank functional GLM*


---

## Description

Multiple testing in permutation inference for the general linear model (GLM)

## Usage

```
frank.flm(
  nsim,
  formula.full,
  formula.reduced,
  curve_sets,
  factors = NULL,
  savefuncs = TRUE,
  lm.args = NULL,
  GET.args = NULL,
  mc.cores = 1,
  mc.args = NULL,
```

```

    cl = NULL,
    method = c("best", "mlm", "lm", "ne")
)

```

## Arguments

<code>nsim</code>	The number of random permutations.
<code>formula.full</code>	The formula specifying the general linear model, see formula in <a href="#">lm</a> .
<code>formula.reduced</code>	The formula of the reduced model with nuisance factors only. This model should be nested within the full model.
<code>curve_sets</code>	A named list of sets of curves giving the dependent variable (Y), and possibly additionally factors whose values vary across the argument values of the functions. The dimensions of the elements should match with each other. Note that factors that are fixed across the functions can be given in the argument factors. Also <a href="#">fdata</a> objects allowed.
<code>factors</code>	A data frame of factors. An alternative way to specify factors when they are constant for all argument values of the functions. The number of rows of the data frame should be equal to the number of curves. Each column should specify the values of a factor.
<code>savefuns</code>	Logical or "return". If TRUE, then the functions from permutations are saved to the attribute <code>simfuns</code> . If "return", then the function returns the permutations in a <code>curve_set</code> , instead of the result of the envelope test on those; this can be used by <a href="#">partial_forder</a> .
<code>lm.args</code>	A named list of additional arguments to be passed to <a href="#">lm</a> . See details.
<code>GET.args</code>	A named list of additional arguments to be passed to <a href="#">global_envelope_test</a> , e.g. <code>typeone</code> specifies the type of multiple testing control, FWER or FDR. See <a href="#">global_envelope_test</a> for the defaults and available options.
<code>mc.cores</code>	The number of cores to use, i.e. at most how many child processes will be run simultaneously. Must be at least one, and parallelization requires at least two cores. On a Windows computer <code>mc.cores</code> must be 1 (no parallelization). For details, see <a href="#">mclapply</a> , for which the argument is passed. Parallelization can be used in generating simulations and in calculating the second stage tests.
<code>mc.args</code>	A named list of additional arguments to be passed to <a href="#">mclapply</a> . Only relevant if <code>mc.cores</code> is more than 1.
<code>cl</code>	Allows parallelization through the use of <a href="#">parLapply</a> (works also in Windows), see the argument <code>cl</code> there, and examples.
<code>method</code>	For advanced use.

## Details

The function `frank.flm` performs a nonparametric test of significance of a covariate in the functional GLM. Similarly as in the graphical functional GLM ([graph.flm](#)), the Freedman-Lane algorithm (Freedman and Lane, 1983) is applied to permute the functions (to obtain the simulations under the null hypothesis of "no effects"); consequently, the test achieves the desired significance

level only approximately. If the reduced model contains only a constant, then the algorithm corresponds to simple permutation of raw data. In contrast to the graphical functional GLM, the Frank functional GLM is based on the F-statistics that are calculated at each argument value of the functions. The global envelope test is applied to the observed and simulated F-statistics. The test is able to find if the factor of interest is significant and also which argument values of the functional domain are responsible for the potential rejection.

The specification of the full and reduced formulas is important. The reduced model should be nested within the full model. The full model should include in addition to the reduced model the interesting factors whose effects are under investigation.

There are different versions of the implementation depending on the application.

- If there are no extra arguments given by the user in `lm.args`, then a fast implementation by solving the normal equations is used to directly compute the F-statistics.
- If all the covariates are constant across the functions, but there are some extra arguments, then a linear model is fitted separately by least-squares estimation to the data at each argument value of the functions fitting a multiple linear model by `lm`. The possible extra arguments passed in `lm.args` to `lm` must be of the form that `lm` accepts for fitting a multiple linear model. In the basic case, no extra arguments are needed.
- If some of the covariates vary across the space and there are user specified extra arguments given in `lm.args`, then the implementation fits a linear model at each argument value of the functions using `lm`, which can be rather slow. The arguments `lm.args` are passed to `lm` for fitting each linear model.

By default the fastest applicable method is used. This can be changed by setting method argument. The cases above correspond to "ne", "mlm" and "lm". Changing the default can be useful for checking the validity of the implementation.

## Value

A `global_envelope` object, which can be printed and plotted directly.

## References

- Mrkvička, T., Myllymäki, M., Kuronen, M. and Narisetty, N. N. (2022) New methods for multiple testing in permutation inference for the general linear model. *Statistics in Medicine* 41(2), 276-297. doi: 10.1002/sim.9236
- Freedman, D., & Lane, D. (1983) A nonstochastic interpretation of reported significance levels. *Journal of Business & Economic Statistics* 1(4), 292-298. doi:10.2307/1391660

## Examples

```
data("GDPtax")
factors.df <- data.frame(Group = GDPtax$Group, Tax = GDPtax$Profittax)
nsim <- 999

res.tax_within_group <- frank.flm(nsim = nsim,
  formula.full = Y~Group+Tax+Group:Tax,
  formula.reduced = Y~Group+Tax,
  curve_sets = list(Y=GDPtax$GDP),
```

```

    factors = factors.df)
plot(res.tax_within_group)

# Image set examples
data("abide_9002_23")
iset <- abide_9002_23$curve_set

res.F <- frank.flm(nsim = 19, # Increase nsim for serious analysis!
  formula.full = Y ~ Group + Age + Sex,
  formula.reduced = Y ~ Age + Sex,
  curve_sets = list(Y = iset),
  factors = abide_9002_23[['factors']],
  GET.args = list(type = "area"))
plot(res.F)

```

GDP

*GDP***Description**

Gross domestic product (GDP)

**Usage**

```
data("GDP")
```

**Format**

A `curve_set` object containing the GDP values for different countries.

**Details**

The data here was constructed based on the following data: The GDP data are publicly available at <https://data.worldbank.org/indicator/NY.GDP.PCAP.CD>. The excel file that we downloaded was called `API_NY.GDP.PCAP.CD_DS2_en_excel_v2_3358980.xls`. The inflation rates are publicly available at <https://data.worldbank.org/indicator/NY.GDP.DEFL.KD.ZG>. The excel file that we downloaded was called `API_NY.GDP.DEFL.KD.ZG_DS2_en_excel_v2_3469555.xls`, from there we took only the inflation rates for United States. Both are distributed under the CC-BY 4.0 license (see <https://datacatalog.worldbank.org/public-licenses#cc-by>).

Then we discounted the GDP of every country in the study to the 1960 USD, and we extrapolated the missing values of the GDP of a country using the closest known ratio of the GDP of the country and the median GDP in that year. Further, the missing values of GDP were interpolated using linear interpolation of the two closest ratios. Appendix of the vignette(`FDRenvelopes`) includes the code to prepare the `curve_set` object.

GDPtax

*GDP per capita with country groups and profit tax***Description**

GDP per capita with country groups and profit tax

**Usage**`data("GDPtax")`**Format**

A list of a three components. The first one (GDP) is a `curve_set` object with components `r` and `obs` containing the years of observations and the GDP curves, i.e. the observed values of GDP in those years. Each column of `obs` contains the GDP for the years for a particular country (seen as column names). The country grouping is given in the list component `Group` and the profit tax in `Profit tax`.

**Details**

The data includes the GDP per capita (current US\$) for years 1980-2017 (World Bank national accounts data, and OECD National Accounts data files). The data have been downloaded from the webpage <https://datamarket.com/data/set/15c9/gdp-per-capita-current-us#!ds=15c9!hd1&display=line>, distributed under the CC-BY 4.0 (<https://datacatalog.worldbank.org/public-licenses#cc-by>). From the same webpage the profit tax in 2010 (World Bank, Doing Business Project (<http://www.doingbusiness.org/ExploreTopics/>) and Total tax rate ( were downloaded. Furthermore, different country groups were formed from countries for which the GDP was available for 1980-2017 and profit tax for 2010:

- Group 1 (Major Advanced Economies (G7)): "Canada", "France", "Germany", "Italy", "Japan"
- Group 2 (Euro Area excluding G7): "Austria", "Belgium", "Cyprus", "Finland", "Greece", "Ireland", "Luxembourg", "Netherlands", "Portugal", "Spain"
- Group 3 (Other Advanced Economies (Advanced Economies excluding G7 and Euro Area)): "Australia", "Denmark", "Iceland", "Norway", "Sweden", "Switzerland"
- Group 4 (Emerging and Developing Asia): "Bangladesh", "Bhutan", "China", "Fiji", "India", "Indonesia", "Malaysia", "Nepal", "Philippines", "Thailand", "Vanuatu"

**References**

World Bank national accounts data, and OECD National Accounts data files. URL: <https://data.worldbank.org/indicator/NY.GDP.PCAP.CD>  
 World Bank, Doing Business Project (<http://www.doingbusiness.org/ExploreTopics/PayingTaxes/>).  
 URL: <https://data.worldbank.org/indicator/IC.TAX.PRFT.CP.ZS>

**See Also**

[graph.flm](#)



**Examples**

```
data("GDPtax")
# Plot data in groups
for(i in 1:4)
  assign(paste0("p", i), plot(subset(GDPtax$GDP, GDPtax$Group == i)) +
    ggplot2::labs(title=paste("Group ", i, sep=""), y="GDP"))
p4
if(require("patchwork", quietly=TRUE))
  p1 + p2 + p3 + p4
```

---

geom\_central\_region      *Central region plot*

---

**Description**

geom\_central\_region and stat\_central\_region can be used to compute and plot central\_region from data arranged in a data.frame.

**Usage**

```
geom_central_region(
  mapping = NULL,
  data = NULL,
  stat = "CentralRegion",
  position = "identity",
  ...,
  coverage = 0.5,
  type = "erl",
  filled = TRUE,
  drawcenterline = TRUE,
  colours = grey.colors(length(coverage), start = 0.9, end = 0.5),
  na.rm = FALSE,
  show.legend = NA,
  inherit.aes = TRUE
)

stat_central_region(
  mapping = NULL,
  data = NULL,
  position = "identity",
  ...,
  coverage = 0.5,
  type = "erl",
  na.rm = FALSE,
  show.legend = NA,
  inherit.aes = TRUE
)
```

**Arguments**

mapping	Set of aesthetic mappings created by <a href="#">aes()</a> . If specified and <code>inherit.aes = TRUE</code> (the default), it is combined with the default mapping at the top level of the plot. You must supply mapping if there is no plot mapping.
data	<p>The data to be displayed in this layer. There are three options:</p> <p>If <code>NULL</code>, the default, the data is inherited from the plot data as specified in the call to <a href="#">ggplot()</a>.</p> <p>A <code>data.frame</code>, or other object, will override the plot data. All objects will be fortified to produce a data frame. See <a href="#">fortify()</a> for which variables will be created.</p> <p>A function will be called with a single argument, the plot data. The return value must be a <code>data.frame</code>, and will be used as the layer data. A function can be created from a formula (e.g. <code>~ head(.x, 10)</code>).</p>
stat	<p>The statistical transformation to use on the data for this layer. When using a <code>geom_*()</code> function to construct a layer, the <code>stat</code> argument can be used to override the default coupling between geoms and stats. The <code>stat</code> argument accepts the following:</p> <ul style="list-style-type: none"> <li>• A Stat ggproto subclass, for example <code>StatCount</code>.</li> <li>• A string naming the stat. To give the stat as a string, strip the function name of the <code>stat_</code> prefix. For example, to use <code>stat_count()</code>, give the stat as <code>"count"</code>.</li> <li>• For more information and other ways to specify the stat, see the <a href="#">layer stat</a> documentation.</li> </ul>
position	<p>A position adjustment to use on the data for this layer. This can be used in various ways, including to prevent overplotting and improving the display. The <code>position</code> argument accepts the following:</p> <ul style="list-style-type: none"> <li>• The result of calling a position function, such as <code>position_jitter()</code>. This method allows for passing extra arguments to the position.</li> <li>• A string naming the position adjustment. To give the position as a string, strip the function name of the <code>position_</code> prefix. For example, to use <code>position_jitter()</code>, give the position as <code>"jitter"</code>.</li> <li>• For more information and other ways to specify the position, see the <a href="#">layer position</a> documentation.</li> </ul>
...	<p>Other arguments passed on to <a href="#">layer()</a>'s <code>params</code> argument. These arguments broadly fall into one of 4 categories below. Notably, further arguments to the <code>position</code> argument, or aesthetics that are required can <i>not</i> be passed through .... Unknown arguments that are not part of the 4 categories below are ignored.</p> <ul style="list-style-type: none"> <li>• Static aesthetics that are not mapped to a scale, but are at a fixed value and apply to the layer as a whole. For example, <code>colour = "red"</code> or <code>linewidth = 3</code>. The geom's documentation has an <b>Aesthetics</b> section that lists the available options. The 'required' aesthetics cannot be passed on to the <code>params</code>. Please note that while passing unmapped aesthetics as vectors is technically possible, the order and required length is not guaranteed to be parallel to the input data.</li> </ul>

- When constructing a layer using a `stat_*`() function, the `...` argument can be used to pass on parameters to the `geom` part of the layer. An example of this is `stat_density(geom = "area", outline.type = "both")`. The `geom`'s documentation lists which parameters it can accept.
- Inversely, when constructing a layer using a `geom_*`() function, the `...` argument can be used to pass on parameters to the `stat` part of the layer. An example of this is `geom_area(stat = "density", adjust = 0.5)`. The `stat`'s documentation lists which parameters it can accept.
- The `key_glyph` argument of `layer()` may also be passed on through `...`. This can be one of the functions described as [key glyphs](#), to change the display of the layer in the legend.

coverage	A number between 0 and 1. The <code>100*coverage%</code> central region will be calculated. A vector of values can also be provided, leading to the corresponding number of central regions.
type	The options and details for type are given in <a href="#">central_region</a> .
filled	Boolean. Should the ribbon be filled?
drawcenterline	Boolean. Should the center line be drawn?
colours	Colours for different coverage levels
na.rm	If FALSE, the default, missing values are removed with a warning. If TRUE, missing values are silently removed.
show.legend	logical. Should this layer be included in the legends? NA, the default, includes if any aesthetics are mapped. FALSE never includes, and TRUE always includes. It can also be a named logical vector to finely select the aesthetics to display.
inherit.aes	If FALSE, overrides the default aesthetics, rather than combining with them. This is most useful for helper functions that define both data and aesthetics and shouldn't inherit behaviour from the default plot specification, e.g. <a href="#">borders()</a> .

## Details

Plots of central regions (global envelopes) with the specified coverage and type (see [central\\_region](#)). When splitting the set of functions to groups by aesthetics or facets, see examples, the central regions are constructed separately for each group, each having the specified coverage.

If  $N_{\text{func}} \cdot (1 - \text{coverage}) < 1$ , where  $N_{\text{func}}$  is the number of functions/curves, the curves are plotted instead of any region.

## Aesthetics

`geom_central_region` requires `x`, `y` and `curveid`. Additionally `geom_central_region` uses the same aesthetics as [geom\\_ribbon](#) if `filled==TRUE` and [geom\\_line](#) otherwise. For multiple coverage values additional aesthetics are not currently supported.

## Computed variables

`stat_central_region` computes `after_stat(ymax)` and `after_stat(ymin)` for the high and low value of the central region.

For multiple coverages the variables use the same names as [central\\_region](#), i.e. `hi.95` and `lo.95` for the region with 95% coverage.

**See Also**

[central\\_region](#) for the basic computation and, [geom\\_ribbon](#) for the default base geom.

**Examples**

```
require("ggplot2")
## Generate some data
#-----
# Simulate regression data according to the cubic model
#  $f(x) = 0.8x - 1.8x^2 + 1.05x^3$  for  $x$  in  $[0,1]$ 
par <- c(0,0.8,-1.8,1.05) # Parameters of the true polynomial model
res <- 100 # Resolution
x <- seq(0, 1, by=1/res); x2=x^2; x3=x^3;

f <- par[1] + par[2]*x + par[3]*x^2 + par[4]*x^3 # The true function
d <- f + rnorm(length(x), 0, 0.04) # Data

# Estimate polynomial regression model
reg <- lm(d ~ x + x2 + x3)
ftheta <- reg$fitted.values
resid0 <- reg$residuals

# Bootstrap regression
B <- 200 # Number of bootstrap samples
df <- NULL
for(i in 1:B) {
  u <- sample(resid0, size=length(resid0), replace=TRUE)
  reg1 <- lm((ftheta+u) ~ x + x2 + x3)
  df <- rbind(df, data.frame(y=reg1$fitted.values, x=x, i=i,
    g2=ifelse(i<14, "A", "B"), g2=ifelse(i<100, "A", "B")))
}

ggplot(df) + geom_line(aes(x, y, group=i))
ggplot(df) + geom_central_region(aes(x=x, y=y, curveid=i), coverage=0.50)
ggplot(df) + geom_central_region(aes(x=x, y=y, curveid=i), coverage=0.50, filled=FALSE)
# Central regions for two groups as specified by 'g2'
ggplot(df) + geom_central_region(aes(x=x, y=y, curveid=i, col=g2), coverage=0.90, filled=FALSE)
ggplot(df) + geom_central_region(aes(x=x, y=y, curveid=i), coverage=0.90) + facet_wrap(vars(g2))

# Central regions with multiple coverage levels
ggplot(df) + geom_central_region(aes(x=x, y=y, curveid=i), coverage=c(0.2,0.4,0.6)) +
  theme_minimal()
ggplot(df) + geom_central_region(aes(x=x, y=y, curveid=i), coverage=seq(0.1, 0.9, length=20),
  colours=rainbow(20))

# Colors for multiregions are not supported
ggplot(df) + geom_central_region(aes(x=x, y=y+0.1*(g2=="B"),
  curveid=i, col=as.factor(g2)), coverage=c(0.05, 0.2,0.4,0.6))

ggplot(df) + geom_central_region(aes(x=x, y=y, curveid=i),
```

```

coverage=c(0.05, 0.2,0.4,0.6)) + facet_wrap(vars(g2))

# Using stat_central_region with geom_linerange and geom_rect
ggplot(df) +
  geom_linerange(aes(curveid=i, x=x, y=y, ymax=after_stat(ymax), ymin=after_stat(ymin),
                    group=g2, col=factor(g2)),
                stat="central_region", coverage = 0.90, position=position_dodge(0.01))
ggplot(within(df, {x = x+0.004*(g2=="B")})) +
  geom_rect(aes(curveid=i, x=x, y=y, xmax=after_stat(x), xmin=after_stat(x+0.004),
                ymax=after_stat(ymax), ymin=after_stat(ymin), group=g2, fill=factor(g2)),
            stat="central_region", coverage = 0.90)

# Non-finite values are not supported
ggplot(within(df, {y = ifelse(runif(length(y)) < 0.001, Inf, y)})) +
  geom_central_region(aes(x=x, y=y, curveid=i))

```

---

GET.composite

*Adjusted global envelope tests*


---

## Description

Adjusted global envelope tests for composite null hypothesis.

## Usage

```

GET.composite(
  X,
  X.ls = NULL,
  nsim = 499,
  nsimsub = nsim,
  simfun = NULL,
  fitfun = NULL,
  calcfun = function(X) {
    X
  },
  testfuns = NULL,
  ...,
  type = "erl",
  alpha = 0.05,
  alternative = c("two.sided", "less", "greater"),
  probs = c(0.025, 0.975),
  r_min = NULL,
  r_max = NULL,
  take_residual = FALSE,
  save.cons.envelope = savefuns,
  savefuns = FALSE,
  verbose = TRUE,

```

```

    MrkvickaEtal2017 = FALSE,
    mc.cores = 1L
)

```

## Arguments

<code>X</code>	An object containing the data in some form. A <code>curve_set</code> object, or an envelope object of the <b>spatstat</b> package, as the <code>curve_sets</code> argument of <code>global_envelope_test</code> (need to provide <code>X.ls</code> ), or a fitted point process model of <b>spatstat</b> (e.g. object of class <code>ppm</code> or <code>kppm</code> ), or a point pattern object of class <code>ppp</code> of <b>spatstat</b> , or another data object (need to provide <code>simfun</code> , <code>fitfun</code> , <code>calcfun</code> ).
<code>X.ls</code>	A list of objects as <code>curve_sets</code> argument of <code>global_envelope_test</code> , giving the second stage simulations, see details.
<code>nsim</code>	The number of simulations to be generated in the primary test. Ignored if <code>X.ls</code> provided.
<code>nsimsub</code>	Number of simulations in each basic test. There will be <code>nsim</code> repetitions of the basic test, each involving <code>nsimsub</code> simulated realisations. Total number of simulations will be <code>nsim * (nsimsub + 1)</code> .
<code>simfun</code>	A function for generating simulations from the null model. If given, this function is called by <code>replicate(n=nsim, simfun(simfun.arg), simplify=FALSE)</code> to make <code>nsim</code> simulations. Here <code>simfun.arg</code> is obtained by <code>fitfun(X)</code> .
<code>fitfun</code>	A function for estimating the parameters of the null model. The function should return the fitted model in the form that it can be directly passed to <code>simfun</code> as its argument.
<code>calcfun</code>	A function for calculating a summary function from a simulation of the model. The default is the identity function, i.e. the simulations from the model are functions themselves. The use of <code>calcfun</code> is still experimental. Preferably provide <code>X</code> and <code>X.ls</code> instead, if <code>X</code> is not a point pattern or fitted point process model object of <b>spatstat</b> .
<code>testfuns</code>	A list of lists of parameters to be passed to the envelope function of <b>spatstat</b> if <code>X</code> is a point pattern or a fitted point process model of <b>spatstat</b> . A list of parameters should be provided for each test function that is to be used in the combined test.
<code>...</code>	Additional parameters to the envelope function of <b>spatstat</b> in the case where only one test function is used. In that case, this is an alternative to providing the parameters in the argument <code>testfuns</code> . If <code>envelope</code> is also used to generate simulations under the null hypothesis (if <code>simfun</code> not provided), then also recall to specify how to generate the simulations.
<code>type</code>	The type of the global envelope with current options for <code>'rank'</code> , <code>'erl'</code> , <code>'cont'</code> , <code>'area'</code> , <code>'qdir'</code> , <code>'st'</code> and <code>'unscaled'</code> . See details.
<code>alpha</code>	The significance level. The $100(1-\alpha)\%$ global envelope will be calculated under the <code>'fwer'</code> or <code>'fdr'</code> control. If a vector of values is provided, the global envelopes are calculated for each value.
<code>alternative</code>	A character string specifying the alternative hypothesis. Must be one of the following: <code>"two.sided"</code> (default), <code>"less"</code> or <code>"greater"</code> . The last two options only available for types <code>'rank'</code> , <code>'erl'</code> , <code>'cont'</code> and <code>'area'</code> .

<code>probs</code>	A two-element vector containing the lower and upper quantiles for the measure 'q' or 'qdir', in that order and on the interval [0, 1]. The default values are 0.025 and 0.975, suggested by Myllymäki et al. (2015, 2017).
<code>r_min</code>	The minimum argument value to include in the test.
<code>r_max</code>	The maximum argument value to include in the test. in calculating functions by the envelope function of <b>spatstat</b> .
<code>take_residual</code>	Logical. If TRUE (needed for visual reasons only) the mean of the simulated functions is reduced from the functions in each first and second stage test.
<code>save.cons.envelope</code>	Logical flag indicating whether to save the unadjusted envelope test results.
<code>savefuncs</code>	Logical flag indicating whether to save all the simulated function values. Similar to the same argument of the envelope function of <b>spatstat</b> .
<code>verbose</code>	Logical flag indicating whether to print progress reports during the simulations. Similar to the same argument of envelope function of <b>spatstat</b> .
<code>MrkvickaEtal2017</code>	Logical. If TRUE, type is "st" or "qdir" and several test functions are used, then the combined scaled MAD envelope presented in Mrkvicka et al. (2017) is calculated. Otherwise, the two-step procedure described in <a href="#">global_envelope_test</a> is used for combining the tests. Default to FALSE. The option is kept for historical reasons.
<code>mc.cores</code>	The number of cores to use, i.e. at most how many child processes will be run simultaneously. Must be at least one, and parallelization requires at least two cores. On a Windows computer mc.cores must be 1 (no parallelization). For details, see <a href="#">mclapply</a> , for which the argument is passed. Parallelization can be used in generating simulations and in calculating the second stage tests.

## Details

The specification of `X`, `X.ls`, `fitfun`, `simfun` is important:

- If `X.ls` is provided, then the global envelope test is calculated based on functions in these objects. `X` should be a [curve\\_set](#) object, or an envelope object of **spatstat**, including the observed function and simulations from the tested model. `X.ls` should be a list of `curve_set` or envelope (of R package **spatstat**) objects, where each component contains an "observed" function `f` that has been simulated from the model fitted to the data and the simulations that have been obtained from the same model that has been fitted to the "observed" `f`. The user has the responsibility that the functions have been generated correctly, the test is done based on these provided simulations. See the examples.
- Otherwise, if `simfun` and `fitfun` are provided, `X` can be general. The function `fitfun` is used for fitting the desired model `M` and the function `simfun` for simulating from a fitted model `M`. These functions should be coupled with each other such that the object returned by `fitfun` is directly accepted as the (single) argument in `simfun`. In the case, that the global envelope should not be calculated directly for `X` (`X` is not a function), `calcfun` can be used to specify how to calculate the function from `X` and from simulations generated by `simfun`. Special attention is needed in the correct specification of the functions, see examples.
- Otherwise, `X` should be either a fitted (point process) model object or a ppp object of the R package **spatstat**.

- If  $X$  is a fitted (point process) model object of the R package **spatstat**, then the simulations from this model are generated and summary functions for testing calculated by the envelope function of **spatstat**. Which summary function to use and how to calculate it, can be passed to `envelope` either in `...` or `testfuns`. Unless otherwise specified the default function of `envelope`, i.g. the K-function, is used. The argument `testfuns` should be used to specify the test functions in the case where one wants to base the test on several test functions.
- If  $X$  is a ppp object of **spatstat**, then the envelope function is used for simulations and model fitting and the complete spatial randomness (CSR) is tested (with intensity parameter).

For the rank envelope test, the global envelope test is the test described in Myllymäki et al. (2017) with the adjustment of Baddeley et al. (2017). For other test types, the test (also) uses the two-stage procedure of Dao and Genton (2014) with the adjustment of Baddeley et al. (2017) as described in Myllymäki and Mrkvička (2024).

See examples also in [saplings](#).

### Value

An object of class `global_envelope` or `combined_global_envelope`, which can be printed and plotted directly. See [global\\_envelope\\_test](#).

### References

- Baddeley, A., Hardegen, A., Lawrence, T., Milne, R. K., Nair, G. and Rakshit, S. (2017). On two-stage Monte Carlo tests of composite hypotheses. *Computational Statistics and Data Analysis* 114: 75-87. doi: <http://dx.doi.org/10.1016/j.csda.2017.04.003>
- Dao, N.A. and Genton, M. (2014). A Monte Carlo adjusted goodness-of-fit test for parametric models describing spatial point patterns. *Journal of Graphical and Computational Statistics* 23, 497-517.
- Mrkvička, T., Myllymäki, M. and Hahn, U. (2017) Multiple Monte Carlo testing, with applications in spatial point processes. *Statistics & Computing* 27(5): 1239-1255. doi: [10.1007/s11222-016-9683-9](https://doi.org/10.1007/s11222-016-9683-9)
- Myllymäki, M., Mrkvička, T., Grabarnik, P., Seijo, H. and Hahn, U. (2017). Global envelope tests for spatial point patterns. *Journal of the Royal Statistical Society: Series B (Statistical Methodology)*, 79: 381-404. doi: [10.1111/rssb.12172](https://doi.org/10.1111/rssb.12172)
- Myllymäki, M. and Mrkvička, T. (2024). GET: Global envelopes in R. *Journal of Statistical Software* 111(3), 1-40. doi: [10.18637/jss.v111.i03](https://doi.org/10.18637/jss.v111.i03)

### See Also

[global\\_envelope\\_test](#), [plot.global\\_envelope](#), [saplings](#)

### Examples

```
# Graphical normality test (Myllymaki and Mrkvicka, 2020, Section 3.3.)
#=====
if(require("fda.usc", quietly=TRUE)) {
```



```

data("poblenou")
dat <- poblenou[["nox"]][["data"]][, 'H10']
n <- length(dat)

# The number of simulations
nsim <- nsimsub <- 199

set.seed(200127)
# General setup
#=====
# 1. Fit the model
mu <- mean(dat)
sigma <- sd(dat)
# 2. Simulate a sample from the fitted null model and
#   compute the test vectors for data (obs) and each simulation (sim)
r <- seq(min(dat), max(dat), length=100)
obs <- stats::ecdf(dat)(r)
sim <- sapply(1:nsimsub, function(i) {
  x <- rnorm(n, mean = mu, sd = sigma)
  stats::ecdf(x)(r)
})
cset <- create_curve_set(list(r = r, obs = obs, sim_m = sim))

# 3. Simulate another sample from the fitted null model.
# 4. Fit the null model to each of the patterns,
#   simulate a sample from the null model,
#   and compute the test vectors for all.
cset.ls <- vector("list", nsim)
for(rep in 1:nsim) {
  x <- rnorm(n, mean = mu, sd = sigma)
  mu2 <- mean(x)
  sigma2 <- sd(x)
  obs2 <- stats::ecdf(x)(r)
  sim2 <- sapply(1:nsimsub, function(i) {
    x2 <- rnorm(n, mean = mu2, sd = sigma2)
    stats::ecdf(x2)(r)
  })
  cset.ls[[rep]] <- create_curve_set(list(r = r, obs = obs2, sim_m = sim2))
}
# Perform the adjusted test
res <- GET.composite(X = cset, X.ls = cset.ls, type = 'erl')
plot(res) + ggplot2::labs(x = "NOx", y = "Ecdf")
}

# Example of a point pattern data
#=====
# Test the fit of a Matern cluster process.

if(require("spatstat.model", quietly=TRUE)) {
  data("saplings")
  saplings <- as.ppp(saplings, W = square(75))
}

```

```

# First choose the r-distances
rmin <- 0.3; rmax <- 10; rstep <- (rmax-rmin)/500
r <- seq(0, rmax, by = rstep)

# Number of simulations
nsim <- 19 # Increase nsim for serious analysis!

# Option 1: Give the fitted model object to GET.composite
#-----
# This can be done and is preferable when the model is
# a point process model of spatstat.
# 1. Fit the Matern cluster process to the pattern
# (using minimum contrast estimation with the K-function)
M1 <- kppm(saplings~1, clusters = "MatClust", statistic = "K")
summary(M1)
# 2. Make the adjusted global area rank envelope test using the L(r)-r function
adjenvL <- GET.composite(X = M1, nsim = nsim,
                        testfuns = list(L = list(fun="Lest", correction="translate",
                                                  transform=expression(.-r), r=r)), # passed to envelope
                        type = "area", r_min = rmin, r_max = rmax)
# Plot the test result
plot(adjenvL)

# Option 2: Generate the simulations "by yourself"
#-----
# and provide them as curve_set or envelope objects
# Preferable when you want to have a control
# on the simulations yourself.
# 1. Fit the model
M1 <- kppm(saplings~1, clusters = "MatClust", statistic = "K")
# 2. Generate nsim simulations by the given function using the fitted model
X <- envelope(M1, nsim = nsim, savefuns = TRUE,
              fun = "Lest", correction = "translate",
              transform = expression(.-r), r = r)
plot(X)
# 3. Create another set of simulations to be used to estimate
# the second-state p-value (as proposed by Baddeley et al., 2017).
simpatterns2 <- simulate(M1, nsim = nsim)
# 4. Calculate the functions for each pattern
simf <- function(rep) {
  # Fit the model to the simulated pattern Xsims[[rep]]
  sim_fit <- kppm(simpatterns2[[rep]], clusters = "MatClust", statistic = "K")
  # Generate nsim simulations from the fitted model
  envelope(sim_fit, nsim = nsim, savefuns = TRUE,
           fun = "Lest", correction = "translate",
           transform = expression(.-r), r = r)
}
X.ls <- parallel::mclapply(X = 1:nsim, FUN = simf, mc.cores = 1) # list of envelope objects
# 5. Perform the adjusted test
res <- GET.composite(X = X, X.ls = X.ls, type = "area",
                    r_min = rmin, r_max = rmax)
plot(res)
}

```

---

GET.distrequa	<i>Graphical n sample test of correspondence of distribution functions</i>
---------------	--

---

## Description

Compare the distributions of two (or more) samples.

## Usage

```
GET.distrequa(
  x,
  stat = "ECDF",
  nsim,
  r = seq(min(unlist((lapply(x, min))))), max(unlist((lapply(x, max))))), length = 100),
  tau = seq(0.1, 0.9, length = 100),
  contrasts = FALSE,
  GET.args = NULL,
  density.args = NULL,
  approxfun.args = NULL,
  rq.args = NULL,
  savefuns = FALSE,
  ...
)
```

## Arguments

<b>x</b>	A list of numeric vectors, one for each sample.
<b>stat</b>	Character string indicating which test statistic to be used. See details.
<b>nsim</b>	The number of random permutations.
<b>r</b>	The sequence of argument values at which the test functions are to be compared. The default is 100 equally spaced values between the minimum and maximum over all groups.
<b>tau</b>	The sequence of argument values for the QR test statistic. The default values are 100 equally spaced values between 0.1 and 0.9.
<b>contrasts</b>	Logical. FALSE and TRUE specify the two test functions as described in description part of this help file.
<b>GET.args</b>	A named list of additional arguments to be passed to <a href="#">global_envelope_test</a> , e.g. typeone specifies the type of multiple testing control, FWER or FDR. See <a href="#">global_envelope_test</a> for the defaults and available options.
<b>density.args</b>	A named list of additional arguments to be passed for the estimation of the test statistic "DEN". For more details see <a href="#">density</a> .
<b>approxfun.args</b>	A named list of additional arguments to be passed for the estimation of the the test statistic "QQ". For more details see <a href="#">approxfun</a> .
<b>rq.args</b>	A named list of additional arguments to be passed for the estimation of the test statistic "QR". For more details see the function rq of <a href="#">quantreq</a> .

savefuns	Logical. If TRUE, then the functions from permutations are saved to the attribute simfuns.
...	Additional parameters to be passed to <a href="#">global_envelope_test</a> . For example, the type of multiple testing control, FWER or FDR must be set by typeone. And, if typeone = "fwer", the type of the global envelope can be chosen by specifying the argument type. See <a href="#">global_envelope_test</a> for the defaults and available options. (The test here uses alternative="two.sided" and nstep=1 (when relevant), but all the other specifications are to be given in ...)

## Details

A global envelope test can be performed to investigate whether the  $n$  distribution functions differ from each other and how do they differ. This test is a generalization of the two-sample Kolmogorov-Smirnov test with a graphical interpretation. We assume that the observations in the sample  $i$  are an i.i.d. sample from the distribution  $F_i(r)$ ,  $i = 1, \dots, n$ , and we want to test the hypothesis

$$F_1(r) = \dots = F_n(r).$$

If contrasts = FALSE (default), then the default test statistic ("ECDF") is taken to be

$$\mathbf{T} = (\hat{F}_1(r), \dots, \hat{F}_n(r))$$

where  $\hat{F}_i(r) = (\hat{F}_i(r_1), \dots, \hat{F}_i(r_k))$  is the ecdf of the  $i$ th sample evaluated at argument values  $r = (r_1, \dots, r_k)$ .

Another possibility is given by stat = "DIFF", and then the test statistic is still based on the ECDFs and constructed from all pairwise differences,

$$\mathbf{T} = (\hat{F}_1(r) - \hat{F}_2(r), \hat{F}_1(r) - \hat{F}_3(r), \dots, \hat{F}_{n-1}(r) - \hat{F}_n(r))$$

The choices contrasts = TRUE and stat = "ECDF" lead to the same test statistic. For other (or multiple) values of stat, the argument contrasts is ignored.

All the options as the test statistics are the following:

1. "ECDF": The ECDFs of the  $n$ -samples, as specified above
2. "DIFF": The pairwise differences between the ECDFs, as specified above
3. "DEN": The kernel estimated density functions of the  $n$ -samples as the test statistic
4. "QQ": The pairwise comparisons of empirical quantiles
5. "SHIFT" The de-trended QQ-plot (shift plot)
6. "QR": The quantile regression process, i.e. the  $\beta$ -coefficients of the quantile regression. By default, the reference category of this test statistic is the first sample.

The test statistics are described in detail in Konstantinou et al. (2024).

The simulations under the null hypothesis that the distributions are the same are obtained by permuting the individuals of the groups. The default number of permutation, if nsim is not specified, is  $n \cdot 1000 - 1$  for the case contrasts = FALSE and  $(n \cdot (n - 1)/2) \cdot 1000 - 1$  for the case contrasts = TRUE, where  $n$  is the length of  $x$ .

## References

Konstantinou K., Mrkvička T. and Myllymäki M. (2024) Graphical n-sample tests of correspondence of distributions. arXiv:2403.01838 [stat.ME] <https://doi.org/10.48550/arXiv.2403.01838>

## Examples

```
if(require("fda", quietly=TRUE)) {
  # Heights of boys and girls at age 10
  f.a <- growth$hgtf["10",] # girls at age 10
  m.a <- growth$hgtm["10",] # boys at age 10
  # Empirical cumulative distribution functions
  plot(ecdf(f.a))
  plot(ecdf(m.a), col='grey70', add=TRUE)
  # Create a list of the data
  fm.list <- list(Girls=f.a, Boys=m.a)

  res <- GET.distrequa(fm.list)
  plot(res)
  # If you want to change the labels:
  plot(res, scales = "free", labels = c("Girls", "Boys"))
  # If you want to change the x-label (y-label similarly):
  require("ggplot2")
  myxlab <- substitute(paste(italic(i), " (", j, ")", sep = ""),
                       list(i = "x", j = "Height in cm"))
  plot(res, scales = "free") + xlab(myxlab)
  # Use instead the test statistics QQ and DEN
  res <- GET.distrequa(fm.list, stat = c("QQ", "DEN"))
  plot(res, scales = "free")

  # Heights of boys and girls at age 14
  f.a <- growth$hgtf["14",] # girls at age 14
  m.a <- growth$hgtm["14",] # boys at age 14
  # Empirical cumulative distribution functions
  plot(ecdf(f.a))
  plot(ecdf(m.a), col='grey70', add=TRUE)
  # Create a list of the data
  fm.list <- list(Girls=f.a, Boys=m.a)

  res <- GET.distrequa(fm.list)
  plot(res) + xlab(myxlab)
  res <- GET.distrequa(fm.list, stat = c("QQ", "DEN"))
  plot(res, scales = "free") + xlab(myxlab)
}

if(require("datasets", quietly=TRUE)) {
  data("iris")
  virginica <- subset(iris, Species == "virginica")
  setosa <- subset(iris, Species == "setosa")
  versicolor <- subset(iris, Species == "versicolor")
}
```

```

res <- GET.distrequal(x = list(virginica = virginica$Sepal.Length,
                              setosa = setosa$Sepal.Length,
                              versicolor = versicolor$Sepal.Length),
                    stat = c("QQ", "DEN"))
plot(res, scales = "free", ncol = 3)

}

```

---

GET.distrindep

*Test of independence of two general distributions*


---

### Description

Permutation-based test of independence in a bivariate vector using as the test statistic either 1) the empirical joint cumulative distribution function, 2) the matrix of observed counts of a 2D contingency table, or 3) the smoothed Q-Q plot.

### Usage

```

GET.distrindep(
  X,
  nsim = 999,
  statistic = c("cdf", "contingency", "qq"),
  ngrid,
  seq.x,
  seq.y,
  sigma,
  atoms.x,
  atoms.y,
  ...
)

```

### Arguments

X	A matrix with n rows and 2 columns. Each row contains one bivariate observation.
nsim	The number of random permutations used.
statistic	Either "cdf", "contingency" or "qq" corresponding to the three test functions.
ngrid	Vector with two elements, giving the number of grid points to be used in the test statistic for each of the two marginals. The default is 20 in each marginal for "cdf" and 64 for "qq". (This is not relevant for "contingency".)
seq.x	For the first marginal, the values at which the empirical cumulative distribution function will be evaluated. If NULL (the default), sequence of quantiles will be used, equidistant in terms of probability. seq.x and seq.y only relevant for "cdf".

<code>seq.y</code>	For the second marginal, the values at which the empirical cumulative distribution function will be evaluated. If NULL (the default), sequence of quantiles will be used, equidistant in terms of probability. <code>seq.x</code> and <code>seq.y</code> only relevant for "cdf".
<code>sigma</code>	Standard deviation of the smoothing kernel to be used for smoothing the Q-Q plot when computing the test statistic. If NULL, sensible default value is used based on the number of observations.
<code>atoms.x</code>	Vector specifying atomic values in the first marginal. Only relevant for "qq". See Examples.
<code>atoms.y</code>	Vector specifying atomic values in the second marginal. Only relevant for "qq". See Examples.
<code>...</code>	Additional parameters to be passed to <a href="#">global_envelope_test</a> . In particular, <code>alpha</code> specifies the nominal significance level of the test, and <code>type</code> the type of the global envelope test.

## Details

The function performs permutation-based test of independence in a bivariate sample based on three different test statistics chosen by the argument `statistic`.

If the observed data are the pairs  $\{(X_1, Y_1), \dots, (X_n, Y_n)\}$ , the permutations are obtained by randomly permuting the values in the second marginal, i.e.  $\{(X_1, Y_{\pi(1)}), \dots, (X_n, Y_{\pi(n)})\}$ .

The first alternative `statistic = "cdf"` is the empirical joint cumulative distribution function computed on a grid of `ngrid[1]` times `ngrid[2]` arguments. The grid points are chosen according to the quantiles of the marginal distributions. The second alternative `statistic = "contingency"` is to test of independence in a 2D contingency table, using the matrix of observed counts as the test statistic. The third alternative `statistic = "qq"` is based on Q-Q representation and estimate of the intensity function computed on a regular grid of `ngrid[1]` times `ngrid[2]` points.

The test itself is in each case performed using the global envelope test of the chosen version, see the argument `type` of [global\\_envelope\\_test](#).

In the case of a 2D contingency table, instead of plotting, text output can be printed in the console by typing the object name. The cells in which the observed value exceeds the upper envelope printed in red, and cells in which the observed value is lower than the lower envelope printed in cyan. Standard output of the global envelope test is also returned and can be plotted or analyzed accordingly.

## References

Dvořák, J. and Mrkvička, T. (2022). Graphical tests of independence for general distributions. *Computational Statistics* 37, 671–699.

## Examples

```
#- Example of cdf
#-----
# Generate sample data
data <- matrix(rnorm(n=200), ncol=2) %*% matrix(c(1,1,0,1), ncol=2)
plot(data)
```

```

# Compute the CDF test and plot the significant regions
res <- GET.distrindep(data, statistic="cdf", ngrid=c(20,15), nsim=1999)

plot(res) + ggplot2::scale_radius(range = 2 * c(1, 6))

# Extract the p-value
attr(res,"p")

#- Example of a 2D contingency table
#-----
# Generate sample data:
data <- matrix(c(sample(4, size=100, replace=TRUE), sample(2, size=100, replace=TRUE)), ncol=2)
data[,2] <- data[,2] + data[,1]

# Observed contingency table (with row names and column names)
table(data[,1], data[,2])

# Permutation-based envelope test
res <- GET.distrindep(data, statistic="contingency", nsim=999)

res
plot(res) + ggplot2::scale_radius(range = 5 * c(1, 6))

# Extract the p-value
attr(res,"p")

# Example of QQ
#-----
# Generate sample data
data <- matrix(rnorm(n=200), ncol=2) %*% matrix(c(1,1,0,1), ncol=2)

plot(data)

# Compute the QQ test and plot the significant regions
res <- GET.distrindep(data, statistic="qq", ngrid=c(30,20), nsim=999)

plot(res)
# Extract the p-value
attr(res,"p")

# With atoms, independent
data <- cbind(rnorm(n=100), sample(4, size=100, replace=TRUE))
plot(data)
res <- GET.distrindep(data, statistic="qq", nsim=999, atoms.y=c(1,2,3,4))

plot(res)

# With atoms, dependent
data <- cbind(sort(rnorm(n=100)), sort(sample(4, size=100, replace=TRUE)))
plot(data)
res <- GET.distrindep(data, statistic="qq", nsim=999, atoms.y=c(1,2,3,4))
plot(res, sign.type="col", what=c("obs", "lo", "hi", "lo.sign", "hi.sign"))

```



```
# Atoms in both variables
data <- cbind(rnorm(n=100), rnorm(n=100)) %% matrix(c(1,1,0,1), ncol=2)
data[,1][data[,1]<=-1] <- -1
data[,2][data[,2]<=-0.5] <- -0.5
plot(data)

# Perform the test
res <- GET.distrindep(data, statistic="qq", nsim=999, atoms.x=c(-1), atoms.y=c(-0.5))

plot(res, sign.type="col", what=c("obs", "lo", "hi", "lo.sign", "hi.sign"))
```

---

GET.localcor

*The test of local correlations*


---

## Description

The test of local correlations using Vilodomat et al. (2014) procedure for resamples and the FDR envelope of Mrkvička and Myllymäki (2023).

## Usage

```
GET.localcor(
  data,
  Delta,
  nsim = 1000,
  ...,
  varying.bandwidth = FALSE,
  bandwidth.nn = 0.1,
  bandwidth.h = 5.281,
  maxk = 300,
  savefuns = FALSE,
  N_s = 1000,
  mc.cores = 1L,
  mc.args = NULL,
  cl = NULL,
  notest = FALSE
)
```

## Arguments

data	A data.frame where the first two columns correspond to the values of the two random fields, whose correlations are to be studied, and the third and fourth columns correspond to the x- and y-coordinates where these random fields have been observed. In addition, the width and height of the pixels at each (x,y) can be given in the fifth and sixth column. Warning: no checks for the data input.
------	--

Delta	A smoothing parameter of the local correlation. According to Vilodomat et al. (2014): Delta is a set of values for the proportion of neighbors to consider for the smoothing step. No default. The user may have to experiment with different values to find one suitable for their data.
nsim	The number of resamples.
...	Additional parameters to be passed to <a href="#">global_envelope_test</a> . Note that for testing local correlations, it may often be preferable to use FDR control. This can be specified by setting <code>typeone = "fdr"</code> , while the default is FWER control. See <a href="#">global_envelope_test</a> for defaults and available options.
varying.bandwidth	Logical, whether to use a varying bandwidth to calculate the local correlations or not. See Vilodomat et al. (2014).
bandwidth.nn	Nearest neighbor component of the smoothing parameter for varying bandwidth to be passed to the argument <code>nn</code> of the function <code>lp</code> of the <b>locfit</b> package. The user may have to experiment with different values to find one suitable for their data. Default set to 0.1 according to Vilodamat et al. (2014, supporting information).
bandwidth.h	Non-varying bandwidth, to be passed to the argument <code>h</code> of the function <code>lp</code> of <b>locfit</b> . The user may have to experiment with different values to find one suitable for their data. Default to 5.281 according to Vilodamat et al. (2014, supporting information).
maxk	See <code>locfit</code> and <code>locfit.raw</code> of <b>locfit</b> . Default here to 300 following Vilodomat et al. (2014).
savefuns	Logical. If TRUE, then the functions from permutations are saved to the attribute <code>simfuns</code> .
N_s	If the number of observations is bigger than <code>N_s</code> , following Vilodomat et al. (2014) a subsample of size <code>N_s</code> is taken every time when a variogram is calculated.
mc.cores	The number of cores to use, i.e. at most how many child processes will be run simultaneously. Must be at least one, and parallelization requires at least two cores. On a Windows computer <code>mc.cores</code> must be 1 (no parallelization). For details, see <a href="#">mclapply</a> , for which the argument is passed. Parallelization can be used in generating simulations and in calculating the second stage tests.
mc.args	A named list of additional arguments to be passed to <a href="#">mclapply</a> . Only relevant if <code>mc.cores</code> is more than 1.
c1	Allows parallelization through the use of <a href="#">parLapply</a> (works also in Windows), see the argument <code>c1</code> there, and examples.
notest	Logical. FALSE means that the test is done. TRUE allows to calculate only local correlation for the data, which can be beneficial for choosing the bandwidths before running the test. If TRUE, then only the observed local correlations will be returned.

## Details

The code is a modification of the supporting information code of Vilodomat et al. (2014) available at <https://doi.org/10.1111/biom.12139>. The modification includes the FDR or FWER envelopes (as

specified by the argument `typeone` in `...`, passed to `global_envelope_test`) for the test of local correlations, i.e. multiple testing correction and graphical illustration of the test results.

Variograms are calculated using the package **geoR** and the local correlations using the R package **loclfit**. These packages should be installed to use `GET.localcor`.

Currently the data is provided in the format of Viladomat et al. (2014, Supporting information). Additionally width and height of area represented by a data point can be provided, see the argument `data`. This information is used for plotting purposes when plotting the output by `plot()`.

Examples are provided in the vignette 'FDRenvelopes', see e.g. <https://cran.r-project.org/package=GET>.

## Value

A global envelope object (with possible additional classes), see description of main components in `global_envelope` (Value). Additional attributes: `p_global` contains the Monte Carlo p-value for the global test of correlation. `cor_global` and `cor_global_sim` contain the value of the correlation for data and permuted data, respectively. If `savefuns = TRUE`, then `permutations` contain the permuted values of the first random field according to Viladomat et al. (2014) procedure, and `cset` contains all the local correlations for the data and permuted data in a `curve_set` object (see `create_curve_set`).

## References

Viladomat, J., Mazumder, R., McInturff, A., McCauley, D.J. and Hastie, T. (2014). Assessing the significance of global and local correlations under spatial autocorrelation: A nonparametric approach. *Biometrics* 70, 409-418. doi: 10.1111/biom.12139

Mrkvička, T., Myllymäki, M. (2023) False discovery rate envelopes. *Statistics and Computing* 33, 109. <https://doi.org/10.1007/s11222-023-10275-7>

---

GET.spatialF

*Testing global and local dependence of point patterns on covariates*

---

## Description

Compute the spatial F- and S-statistics and perform the one-stage global envelope tests proposed by Myllymäki et al. (2020).

## Usage

```
GET.spatialF(
  X,
  formula.full,
  formula.reduced,
  fitfun,
  covariates,
  nsim,
  bw = spatstat.explore::bw.scott(X),
  bw.S = bw,
```

```

    dimyx = NULL,
    ...
)

```

### Arguments

<code>X</code>	A ppp object of <b>spatstat</b> representing the observed point pattern.
<code>formula.full</code>	A formula for the trend of the full model.
<code>formula.reduced</code>	A formula for the trend of the reduced model that is a submodel of the full model.
<code>fitfun</code>	A function of a point pattern, model formula and covariates, giving a fitted model object that can be used with <a href="#">simulate</a> .
<code>covariates</code>	A list of covariates.
<code>nsim</code>	The number of simulations.
<code>bw</code>	The bandwidth for smoothed residuals.
<code>bw.S</code>	The radius for the local S(u)-statistic.
<code>dimyx</code>	Pixel array dimensions for smoothed residuals. See as.mask of <b>spatstat</b> .
<code>...</code>	Additional arguments to be passed to <a href="#">global_envelope_test</a> .

### Value

list with three components

- `F` = the global envelope test based on the  $F(u)$  statistic
- `S` = the global envelope test based on the  $S(u)$  statistic
- `coef` = the coefficients of the full model given by `fitfun`

### References

Myllymäki, M., Kuronen, M. and Mrkvička, T. (2020). Testing global and local dependence of point patterns on covariates in parametric models. *Spatial Statistics* 42, 100436. doi: 10.1016/j.spasta.2020.100436

### Examples

```

if(require("spatstat.model", quietly=TRUE)) {
  # Example of tropical rain forest trees
  data("bei")

  fullmodel <- ~ grad
  reducedmodel <- ~ 1
  fitppm <- function(X, model, covariates) {
    ppm(X, model, covariates=covariates)
  }

  nsim <- 19 # Increase nsim for serious analysis!
  res <- GET.spatialF(bei, fullmodel, reducedmodel, fitppm, bei.extra, nsim)
}

```

```

plot(res$F)
plot(res$S)

# Example of forest fires
data("clmfires")
# Choose the locations of the lightnings in years 2004-2007:
pp.lightning <- unmark(subset(clmfires, cause == "lightning" &
                             date >= "2004-01-01" & date < "2008-01-01"))

covariates <- clmfires.extra$clmcov100
covariates$forest <- covariates$landuse == "conifer" | covariates$landuse == "denseforest" |
                    covariates$landuse == "mixedforest"

fullmodel <- ~ elevation + landuse
reducedmodel <- ~ landuse
nsim <- 19 # Increase nsim for serious analysis!
res <- GET.spatialF(pp.lightning, fullmodel, reducedmodel, fitppm, covariates, nsim)
plot(res$F)
plot(res$S)

# Examples of the fitfun functions for clustered and regular processes
# fitfun for the log Gaussian Cox Process with exponential covariance function
fitLGCPexp <- function(X, model, covariates) {
  kppm(X, model, clusters="LGCP", model="exponential", covariates=covariates)
}
# fitfun for the hardcore process with hardcore radius 0.01
fitHardcore <- function(X, model, covariates) {
  ppm(X, model, interaction=Hardcore(0.01), covariates=covariates)
}
}

```

---

GET.variogram

---

Variogram and residual variogram with global envelopes

---

## Description

The function accompanies the function `variogram` with global envelopes that are based on permutations of the variable(s) or residuals for which the variogram is calculated. Therefore, one can inspect the hypothesis of "no spatial autocorrelation" of the variable or the residuals of the fitted model.

## Usage

```

GET.variogram(
  object,
  nsim = 999,
  data = NULL,

```

```
...,
GET.args = NULL,
savefuns = TRUE
)
```

## Arguments

object	An object of class <code>gstat</code> or a <code>variogram.formula</code> . In the first case, direct (residual) variograms are calculated for the variable defined in object. Only one variable allowed. In the second case, a formula defining the response vector and (possible) regressors, in case of absence of regressors, use e.g. <code>z~1</code> . See <a href="#">variogram</a> .
nsim	The number of permutations.
data	A data frame where the names in formula are to be found. If <code>NULL</code> , the data are assumed to be found in the object.
...	Additional parameters to be passed to <a href="#">variogram</a> .
GET.args	A named list of additional arguments to be passed to <a href="#">global_envelope_test</a> .
savefuns	Logical. If <code>TRUE</code> , then the functions from permutations are saved to the attribute <code>simfuns</code> .

## Examples

```
if(require("sp", quietly=TRUE) & require("gstat", quietly=TRUE)) {
  # Examples from gstat complemented with global envelopes
  #-----
  data("meuse")
  coordinates(meuse) <- ~x+y
  # topsoil zinc concentration, mg kg-1 soil ("ppm")
  bubble(meuse, "zinc",
         col=c("#00ff0088", "#00ff0088"), main="zinc concentrations (ppm)")
  # Variogram can be calculated as follows by the function variogram of the gstat package.
  # The function variogram takes a formula as its first argument:
  # log(zinc)~1 means that we assume a constant trend for the variable log(zinc).
  lzn.vgm <- variogram(object=log(zinc)~1, data=meuse)
  plot(lzn.vgm)
  # Variogram with global envelopes is as easy:
  lzn.vgm.GET <- GET.variogram(object=log(zinc)~1, data=meuse)

  plot(lzn.vgm.GET)

  # Instead of the constant mean, denoted by ~1, a mean function can
  # be specified, e.g. using ~sqrt(dist) as a predictor variable:
  lznr.vgm <- variogram(log(zinc)~sqrt(dist), meuse)
  # In this case, the variogram of residuals with respect
  # to a fitted mean function are shown.
  plot(lznr.vgm)
  # The variogram with global envelopes (obtained by permuting the residuals):
  lznr.vgm.GET <- GET.variogram(object=log(zinc)~sqrt(dist), data=meuse)

  plot(lznr.vgm.GET)
```

```

# Directional variograms
lzn.dir <- variogram(object=log(zinc)~1, data=meuse, alpha=c(0, 45, 90, 135))
plot(lzn.dir)
# with global envelopes
lzn.dir.GET <- GET.variogram(object=log(zinc)~1, data=meuse, alpha=c(0, 45, 90, 135))

plot(lzn.dir.GET)

# Use instead gstat objects
g <- gstat(id="ln.zinc", formula=log(zinc)~1, data=meuse)
# or: g <- gstat(id="ln.zinc", formula=log(zinc)~sqrt(dist), data=meuse)
# The variogram
plot(variogram(g))
# The variogram with global envelopes:
g.GET <- GET.variogram(object=g)

plot(g.GET)
}

```

---

global\_envelope\_test    *Global envelope test*

---

## Description

Global envelope test, global envelopes and p-values

## Usage

```

global_envelope_test(
  curve_sets,
  typeone = c("fwer", "fdr"),
  alpha = 0.05,
  alternative = c("two.sided", "less", "greater"),
  type = "erl",
  algorithm = c("IATSE", "ATSE"),
  ties = "erl",
  probs = c(0.025, 0.975),
  quantile.type = 7,
  central = "mean",
  nstep = 2,
  ...,
  lower = NULL,
  upper = NULL
)

```

**Arguments**

curve_sets	A <a href="#">curve_set</a> object or a list of <a href="#">curve_set</a> objects containing a data function and simulated functions from which the envelope is to be constructed. Also envelope objects of <b>spatstat</b> are accepted instead of curve_set objects. If an envelope object is given, it must contain the summary functions from simulated patterns which can be achieved by setting savefuncs = TRUE when calling the envelope function.
typeone	Character string indicating which type I error rate to control, either the family-wise error rate ('fwer') or false discovery rate ('fdr').
alpha	The significance level. The 100(1-alpha)% global envelope will be calculated under the 'fwer' or 'fdr' control. If a vector of values is provided, the global envelopes are calculated for each value.
alternative	A character string specifying the alternative hypothesis. Must be one of the following: "two.sided" (default), "less" or "greater". The last two options only available for types 'rank', 'erl', 'cont' and 'area'.
type	The type of the global envelope with current options for 'rank', 'erl', 'cont', 'area', 'qdir', 'st' and 'unscaled'. See details.
algorithm	The algorithm for the computation of the FDR envelope. Either "IATSE" or "ATSE" standing for the iteratively adaptive two-stage envelope and the adaptive two-stage envelope, respectively, see Mrkvička and Myllymäki (2023).
ties	The method to obtain a unique p-value when typeone = 'fwer' and type = 'rank'; otherwise ignored. Possible values are 'midrank', 'random', 'conservative', 'liberal' and 'erl'. For 'conservative' the resulting p-value will be the highest possible. For 'liberal' the p-value will be the lowest possible. For 'random' the rank of the obs within the tied values is uniformly sampled so that the resulting p-value is at most the conservative option and at least the liberal option. For 'midrank' the mid-rank within the tied values is taken. For 'erl' the extreme rank length p-value is calculated. The default is 'erl'.
probs	A two-element vector containing the lower and upper quantiles for the measure 'q' or 'qdir', in that order and on the interval [0, 1]. The default values are 0.025 and 0.975, suggested by Myllymäki et al. (2015, 2017).
quantile.type	As type argument of <a href="#">quantile</a> , how to calculate quantiles for 'q' or 'qdir'.
central	Either "mean" or "median". If the curve sets do not contain the component theo for the theoretical central function, then the central function (used for plotting only) is calculated either as the mean or median of functions provided in the curve sets. For 'qdir', 'st' and 'unscaled' only the mean is allowed as an option, due to their definition.
nstep	1 or 2 for how to construct a combined global envelope if list of curve sets is provided. 2 (default) for a two-step combining procedure, 1 for one-step.
...	Additional parameters to be passed to <a href="#">central_region</a> for the computation of the 'fwer' envelope.
lower	A single number (or a vector of suitable length) giving a lower bound for the functions. Used only for the extension of the FDR envelope.
upper	A single number (or a vector of suitable length) giving an upper bound for the functions. Used only for the extension of the FDR envelope.



## Details

Given a `curve_set` object, or an envelope object of `spatstat`, which contains both the data curve (or function or vector)  $T_1(r)$  (in the component `obs`) and the simulated curves  $T_2(r), \dots, T_{s+1}(r)$  (in the component `sim_m`), the function `global_envelope_test` performs a global envelope test, either under the control of family-wise error rate (FWER) or false discovery rate (FDR), as specified by the argument `typeone`. The function `global_envelope_test` is the main function for global envelope tests (for simple hypotheses).

The case `typeone = "fdr"` corresponds to the FDR envelopes proposed by Mrkvička and Myllymäki (2023). See details in `fdr_envelope` and in the vignette `vignette("FDRenvelopes")`. Note there also the arguments that are the relevant ones for the FDR envelope specification. The descriptions below concern the FWER envelopes.

If `typeone = "fwer"`, the functionality of the function is rather similar to the function `central_region`, but in addition to ordering the functions from the most extreme one to the least extreme one using different measures and providing the global envelopes with intrinsic graphical interpretation, p-values are calculated for the test. Thus, while `central_region` can be used to construct global envelopes in a general setting, the function `global_envelope_test` is devoted to testing as its name suggests.

Different type of global envelope tests under the control of FWER can be computed. We use such ordering of the functions for which we are able to construct global envelopes with intrinsic graphical interpretation (IGI, see Myllymäki and Mrkvička, 2023).

- `'rank'`: the completely non-parametric rank envelope test (Myllymäki et al., 2017) based on minimum of pointwise ranks
- `'erl'`: the completely non-parametric rank envelope test based on extreme rank lengths (Myllymäki et al., 2017; Mrkvička et al., 2018) based on number of minimal pointwise ranks
- `'cont'`: the completely non-parametric rank envelope test based on continuous rank (Hahn, 2015; Mrkvička et al., 2022) based on minimum of continuous pointwise ranks
- `'area'`: the completely non-parametric rank envelope test based on area rank (Mrkvička et al., 2022) based on area between continuous pointwise ranks and minimum pointwise ranks for those argument (`r`) values for which pointwise ranks achieve the minimum (it is a combination of `erl` and `cont`)
- `"qdir"`, the directional quantile envelope test, protected against unequal variance and asymmetry of  $T(r)$  for different distances  $r$  (Myllymäki et al., 2015, 2017)
- `"st"`, the studentised envelope test, protected against unequal variance of  $T(r)$  for different distances  $r$  (Myllymäki et al., 2015, 2017)
- `"unscaled"`, the unscaled envelope (providing a baseline) that has a constant width and that corresponds to the classical maximum deviation test (Ripley, 1981).

The first four types are global rank envelopes. The `'rank'` envelope test is a completely non-parametric test, which provides the  $100(1-\alpha)\%$  global envelope for the chosen test function  $T(r)$  on the chosen interval of distances and associated p-values. The other three are modifications of `'rank'` to treat the ties in the extreme rank ordering on which the `'rank'` test is based on. The last three envelopes are global scaled maximum absolute difference (MAD) envelope tests. The unscaled envelope test leads to envelopes with constant width over the distances  $r$ . Thus, it suffers from unequal variance of  $T(r)$  over the distances  $r$  and from the asymmetry of distribution of  $T(r)$ .

We recommend to use the other global envelope tests available. The unscaled envelope is provided as a reference.

See Myllymäki and Mrkvička (2023, Appendix.), i.e. `vignette("GET")`, for more detailed description of the measures and the corresponding envelopes.

See `vignette("pointpatterns")` for examples of point pattern analyses. See `vignette("FDRenvelopes")` for examples of FDR envelopes obtained by `typeone = "fdr"`.

## Value

Either an object of class `"global_envelope"` or `"combined_global_envelope"`, similarly as the objects returned by `central_region`. Further, if `typeone = "fdr"`, the objects have the further class `"fdr_envelope"` or `"combined_fdr_envelope"`.

The `global_envelope` is essentially a data frame containing columns

- the values of the argument `r` at which the test was made, copied from the argument `curve_sets` with the corresponding names
- `obs` = values of the data function, copied from the argument `curve_sets` (unlike for central regions, `obs` always exists for a global envelope test)
- `lo` = the lower envelope; in case of a vector of `alpha` values, several `'lo'` exist with names `paste0("lo.", 100*(1-alpha))`
- `hi` = the upper envelope; in case of a vector of `alpha` values, several `'lo'` exist with names `paste0("hi.", 100*(1-alpha))`
- `central` = a central curve as specified in the argument `central`.

Moreover, the returned object has the same attributes as the `global_envelope` object returned by `central_region` and in addition

- `p` = the p-value of the test

and in the case that `type = 'rank'` also

- `p_interval` = The p-value interval  $[p_{liberal}, p_{conservative}]$ .
- `ties` = As the argument `ties`.

The `combined_global_envelope` is a list of `global_envelope` objects containing the above mentioned columns and which all together form the global envelope. It has the same attributes as described in `central_region`, and in addition also the p-value `p`. The 2d classes are attached as described in `central_region`.

## Procedure

1) First the curves are ranked from the most extreme one to the least extreme one by a measure that is specified by the argument `type`. The options are

- `'rank'`: extreme ranks (Myllymäki et al., 2017)
- `'erl'`: extreme rank lengths (Myllymäki et al., 2017; Mrkvička et al., 2018)
- `'cont'`: continuous ranks (Hahn, 2015; Mrkvička et al., 2019)
- `'area'`: area ranks (Mrkvička et al., 2019)

- 'qdir': the directional quantile maximum absolute deviation (MAD) measure (Myllymäki et al., 2015, 2017)
- 'st': the studentized MAD measure (Myllymäki et al., 2015, 2017)
- 'unscaled': the unscaled MAD measure (Ripley, 1981)

2) Based on the measures used to rank the functions, the  $100(1-\alpha)\%$  global envelope is provided. It corresponds to the  $100*\text{coverage}\%$  central region.

3) P-values: In the case `type="rank"`, based on the extreme ranks  $k_i, i = 1, \dots, s+1$ , the p-interval is calculated. Because the extreme ranks contain ties, there is not just one p-value. The p-interval is given by the most liberal and the most conservative p-value estimate. Also a single p-value is calculated. By default this single p-value is the extreme rank length p-value ("erl") as specified by the argument `ties`. If the case of other measures, a (single) p-value based on the given ordering of the functions is calculated and returned in the attribute `p`.

### Number of simulations

For the global "rank" envelope test, Myllymäki et al. (2017) recommended to use at least 2500 simulations for testing at the significance level  $\alpha = 0.05$  for single function tests, based on experiments with summary functions for point processes evaluated approximately at 500 argument values. In this case, the width of the p-interval associated with the extreme rank measure tended to be smaller than 0.01. The tests 'erl', 'cont' and 'area', similarly as the MAD deviation/envelope tests 'qdir', 'st' and 'unscaled', allow in principle a lower number of simulations to be used than the test based on extreme ranks ('rank'), because no ties occur for these measures. If affordable, we recommend in any case some thousands of simulations for all the measures to achieve a good power and repeatability of the test. If the dimension of the test functions is higher, also the number of simulations should preferably be higher.

### Tests based on several functions

If a list of (suitable) objects are provided in the argument `curve_sets`, then by default (`nstep = 2`) the two-step combining procedure is used to perform the combined global test as described in Myllymäki and Mrkvička (2024). If `nstep = 1` and the lengths of the multivariate vectors in each component of the list are equal, then the one-step combining procedure is used where the functions are concatenated together into a one long vector.

### References

- Mrkvička, T., Myllymäki, M. and Hahn, U. (2017). Multiple Monte Carlo testing, with applications in spatial point processes. *Statistics & Computing* 27(5), 1239-1255. doi: 10.1007/s11222-016-9683-9
- Mrkvička, T., Myllymäki, M., Jilek, M. and Hahn, U. (2020) A one-way ANOVA test for functional data with graphical interpretation. *Kybernetika* 56(3), 432-458. doi: 10.14736/kyb-2020-3-0432
- Mrkvička, T., Myllymäki, M., Kuronen, M. and Narisetty, N. N. (2022) New methods for multiple testing in permutation inference for the general linear model. *Statistics in Medicine* 41(2), 276-297. doi: 10.1002/sim.9236
- Mrkvička and Myllymäki (2023). False discovery rate envelopes. *Statistics and Computing* 33, 109. <https://doi.org/10.1007/s11222-023-10275-7>

- Myllymäki, M., Grabarnik, P., Seijo, H. and Stoyan, D. (2015). Deviation test construction and power comparison for marked spatial point patterns. *Spatial Statistics* 11, 19-34. doi: 10.1016/j.spasta.2014.11.004
- Myllymäki, M., Mrkvička, T., Grabarnik, P., Seijo, H. and Hahn, U. (2017). Global envelope tests for spatial point patterns. *Journal of the Royal Statistical Society: Series B (Statistical Methodology)* 79, 381–404. doi: 10.1111/rssb.12172
- Myllymäki, M. and Mrkvička, T. (2024). GET: Global envelopes in R. *Journal of Statistical Software* 111(3), 1-40. doi: 10.18637/jss.v111.i03
- Ripley, B.D. (1981). *Spatial statistics*. Wiley, New Jersey.

### See Also

[plot.global\\_envelope](#), [central\\_region](#), [GET.composite](#)

### Examples

```
# Goodness-of-fit testing for simple hypothesis
if(require("spatstat.explore", quietly=TRUE)) {
  # Testing complete spatial randomness (CSR)
  #=====
  X <- unmark(spruces)

  nsim <- 1999 # Number of simulations

  # Illustration of general workflow for simple hypotheses
  #=====
  # First illustrate the general workflow for the test by this example
  # of CSR test for a point pattern X using the empirical L-function.
  # Define the argument values at which the functions are evaluated
  obs.L <- Lest(X, correction="translate")
  r <- obs.L[['r']]
  # The test function for the data
  obs <- obs.L[['trans']] - r
  # Prepare simulations and calculate test functions for them at same r as 'obs'
  sim <- matrix(nrow=length(r), ncol=nsim)
  for(i in 1:nsim) {
    sim.X <- runifpoint(ex=X) # simulation under CSR
    sim[, i] <- Lest(sim.X, correction="translate", r=r[['trans']] - r)
  }
  # Create a curve_set containing argument values, observed and simulated functions
  cset <- curve_set(r=r, obs=obs, sim=sim)
  # Perform the test
  res <- global_envelope_test(cset, type="erl")
  plot(res) + ggplot2::ylab(expression(italic(hat(L)(r)-r)))

  # Simple hypothesis for a point pattern utilizing the spatstat package
  #=====
  # Generate nsim simulations under CSR, calculate L-function for the data and simulations
  env <- envelope(X, fun="Lest", nsim=nsim,
    savefuns=TRUE, # save the functions
    correction="translate", # edge correction for L
```

```

        transform=expression(.-r), # centering
        simulate=expression(runifpoint(ex=X))) # Simulate CSR
# The rank envelope test (ERL)
res <- global_envelope_test(env, type="erl")
# Plot the result
plot(res)

## Advanced use:
# Choose the interval of distances [r_min, r_max] (at the same time create a curve_set from 'env')
cset <- crop_curves(env, r_min=1, r_max=7)
# Do the rank envelope test (erl)
res <- global_envelope_test(cset, type="erl")
plot(res) + ggplot2::ylab(expression(italic(L(r)-r)))

# A combined global envelope test
#=====
# As an example test CSR of the saplings point pattern by means of
# L, F, G and J functions.
data(saplings)
X <- as.ppp(saplings, W=square(75))

nsim <- 499 # Number of simulations

# Specify distances for different test functions
n <- 500 # the number of r-values
rmin <- 0; rmax <- 20; rstep <- (rmax-rmin)/n
rminJ <- 0; rmaxJ <- 8; rstepJ <- (rmaxJ-rminJ)/n
r <- seq(0, rmax, by=rstep) # r-distances for Lest
rJ <- seq(0, rmaxJ, by=rstepJ) # r-distances for Fest, Gest, Jest

# Perform simulations of CSR and calculate the L-functions
env_L <- envelope(X, nsim=nsim,
  simulate=expression(runifpoint(ex=X)),
  fun="Lest", correction="translate",
  transform=expression(.-r), # Take the L(r)-r function instead of L(r)
  r=r, # Specify the distance vector
  savefuns=TRUE, # Save the estimated functions
  savepatterns=TRUE) # Save the simulated patterns
# Take the simulations from the returned object
simulations <- attr(env_L, "simpatterns")
# Then calculate the other test functions F, G, J for each simulated pattern
env_F <- envelope(X, nsim=nsim, simulate=simulations,
  fun="Fest", correction="Kaplan", r=rJ,
  savefuns=TRUE)
env_G <- envelope(X, nsim=nsim, simulate=simulations,
  fun="Gest", correction="km", r=rJ,
  savefuns=TRUE)
env_J <- envelope(X, nsim=nsim, simulate=simulations,
  fun="Jest", correction="none", r=rJ,
  savefuns=TRUE)

# Crop the curves to the desired r-interval I

```

```

curve_set_L <- crop_curves(env_L, r_min=rmin, r_max=rmax)
curve_set_F <- crop_curves(env_F, r_min=rminJ, r_max=rmaxJ)
curve_set_G <- crop_curves(env_G, r_min=rminJ, r_max=rmaxJ)
curve_set_J <- crop_curves(env_J, r_min=rminJ, r_max=rmaxJ)

res <- global_envelope_test(curve_sets=list(L=curve_set_L, F=curve_set_F,
                                           G=curve_set_G, J=curve_set_J))

plot(res)
plot(res, labels=c("L(r)-r", "F(r)", "G(r)", "J(r)"))
}

# A test based on a low dimensional random vector
#=====
# Let us generate some example data.
X <- matrix(c(-1.6,1.6),1,2) # data pattern X=(X_1,X_2)
if(requireNamespace("mvtnorm", quietly=TRUE)) {
  Y <- mvtnorm::rmvnorm(200,c(0,0),matrix(c(1,0.5,0.5,1),2,2)) # simulations
  plot(Y, xlim=c(min(X[,1],Y[,1]), max(X[,1],Y[,1])), ylim=c(min(X[,2],Y[,2]), max(X[,2],Y[,2])))
  points(X, col=2)

  # Test the null hypothesis is that X is from the distribution of Y's (or if it is an outlier).

  # Case 1. The test vector is (X_1, X_2)
  cset1 <- curve_set(r=1:2, obs=as.vector(X), sim=t(Y))
  res1 <- global_envelope_test(cset1)
  plot(res1)

  # Case 2. The test vector is (X_1, X_2, (X_1-mean(Y_1))*(X_2-mean(Y_2))).
  t3 <- function(x, y) { (x[,1]-mean(y[,1]))*(x[,2]-mean(y[,2])) }
  cset2 <- curve_set(r=1:3, obs=c(X[,1],X[,2],t3(X,Y)), sim=rbind(t(Y), t3(Y,Y)))
  res2 <- global_envelope_test(cset2)
  plot(res2)
}

```

---

global\_rq

*Global quantile regression*


---

## Description

Global tests of significance for the effect of covariates in quantile regression

## Usage

```

global_rq(
  nsim,
  formula.full,
  formula.reduced,
  taus,
  data,
  contrasts = NULL,

```

```

permutationstrategy = c("Freedman-Lane", "Freedman-Lane+remove zeros",
  "within nuisance", "remove location", "remove location scale", "remove quantile"),
savefuns = FALSE,
rq.args = NULL,
lm.args = NULL,
GET.args = NULL,
mc.cores = 1L,
mc.args = NULL,
cl = NULL
)

```

### Arguments

<code>nsim</code>	The number of random permutations.
<code>formula.full</code>	The formula specifying the general linear model, see formula in <a href="#">lm</a> .
<code>formula.reduced</code>	The formula of the reduced model with nuisance factors only. This model should be nested within the full model.
<code>taus</code>	The quantiles to be used.
<code>data</code>	data.frame where to look for variables.
<code>contrasts</code>	Passed directly to <code>rq</code> .
<code>permutationstrategy</code>	The permutation strategy to be used. See details.
<code>savefuns</code>	Logical. If TRUE, then the functions from permutations are saved to the attribute <code>simfuns</code> .
<code>rq.args</code>	Additional arguments passed to <code>rq</code> .
<code>lm.args</code>	A named list of additional arguments to be passed to <a href="#">lm</a> . See details.
<code>GET.args</code>	A named list of additional arguments to be passed to <a href="#">global_envelope_test</a> , e.g. <code>typeone</code> specifies the type of multiple testing control, FWER or FDR. See <a href="#">global_envelope_test</a> for the defaults and available options.
<code>mc.cores</code>	The number of cores to use, i.e. at most how many child processes will be run simultaneously. Must be at least one, and parallelization requires at least two cores. On a Windows computer <code>mc.cores</code> must be 1 (no parallelization). For details, see <a href="#">mclapply</a> , for which the argument is passed. Parallelization can be used in generating simulations and in calculating the second stage tests.
<code>mc.args</code>	A named list of additional arguments to be passed to <a href="#">mclapply</a> . Only relevant if <code>mc.cores</code> is more than 1.
<code>cl</code>	Allows parallelization through the use of <a href="#">parLapply</a> (works also in Windows), see the argument <code>cl</code> there, and examples.

### Details

The possible permutation strategies include "Freedman-Lane" (FL), "Freedman-Lane+remove zeros" (FL+), "within nuisance" (WN), "remove location" (RL), "remove location scale" (RLS), "remove quantile" (RQ), which correspond to those in Mrkvička et al. (Section 4.1-4.6 and Table 1).

**Value**

A `global_envelope` or `combined_global_envelope` object, which can be printed and plotted directly.

**References**

Mrkvička, T., Konstantinou, K., Kuronen, M. and Myllymäki, M. (2023) Global quantile regression. arXiv:2309.04746 [stat.ME]. <https://doi.org/10.48550/arXiv.2309.04746>

Myllymäki, M. and Mrkvička, T. (2024). GET: Global envelopes in R. *Journal of Statistical Software* 111(3), 1-40. doi: 10.18637/jss.v111.i03

Freedman, D., & Lane, D. (1983) A nonstochastic interpretation of reported significance levels. *Journal of Business & Economic Statistics*, 1(4), 292-298. doi:10.2307/1391660

**Examples**

```
if(require("quantreg", quietly=TRUE)) {
  data("stackloss")
  res <- global_rq(nsim = 19, # Increase nsim for serious analysis!
    formula.full = stack.loss ~ Air.Flow + Water.Temp + Acid.Conc.,
    formula.reduced = stack.loss ~ Water.Temp,
    taus = seq(0.1, 0.9, length=10), permutationstrategy = "remove quantile",
    data = stackloss, GET.args = list(typeone = "fwer"))
  plot(res)
}
```

---

graph.fanova

---

One-way graphical functional ANOVA

---

**Description**

One-way ANOVA tests for functional data with graphical interpretation

**Usage**

```
graph.fanova(
  nsim,
  curve_set,
  groups,
  variances = "equal",
  contrasts = FALSE,
  n.aver = 1L,
  mirror = FALSE,
  savefuns = FALSE,
  test.equality = c("mean", "var", "cov"),
  cov.lag = 1,
  ...
)
```



**Arguments**

<code>nsim</code>	The number of random permutations.
<code>curve_set</code>	The original data (an array of functions) provided as a <a href="#">curve_set</a> object or a <code>fdata</code> object of <b>fda.usc</b> . The curve set should include the argument values for the functions in the component <code>r</code> , and the observed functions in the component <code>obs</code> .
<code>groups</code>	The original groups (a factor vector representing the assignment to groups).
<code>variances</code>	Either "equal" or "unequal". If "unequal", then correction for unequal variances as explained in details will be done. Only relevant for the case <code>test.equality = "means"</code> (default).
<code>contrasts</code>	Logical. FALSE and TRUE specify the two test functions as described in description part of this help file.
<code>n.aver</code>	If <code>variances = "unequal"</code> , there is a possibility to use variances smoothed by applying moving average to the estimated sample variances. <code>n.aver</code> determines how many values on each side do contribute (incl. value itself).
<code>mirror</code>	The complement of the argument circular of <a href="#">filter</a> . Another parameter for the moving average to estimate sample variances (see <code>n.aver</code> ).
<code>savefuns</code>	Logical. If TRUE, then the functions from permutations are saved to the attribute <code>simfuns</code> .
<code>test.equality</code>	A character with possible values mean (default), var and cov. If mean, the functional ANOVA is performed to compare the means in the groups. If var, then the equality of variances of the curves in the groups is tested by performing the graphical functional ANOVA test on the functions

$$Z_{ij}(r) = T_{ij}(r) - \bar{T}_j(r).$$

If cov, then the equality of lag cov. lag covariance is tested by performing the fANOVA with

$$W_{ij}(r) = \sqrt{|V_{ij}(r)| \cdot \text{sign}(V_{ij}(r))},$$

where

$$V_{ij}(r) = (T_{ij}(r) - \bar{T}_j(r))((T_{ij}(r + s) - \bar{T}_j(r + s))).$$

See Mrkvicka et al. (2020) for more details.

<code>cov.lag</code>	The lag of the covariance for testing the equality of covariances, see <code>test.equality</code> .
<code>...</code>	Additional parameters to be passed to <a href="#">global_envelope_test</a> . For example, the type of multiple testing control, FWER or FDR must be set by <code>typeone</code> . And, if <code>typeone = "fwer"</code> , the type of the global envelope can be chosen by specifying the argument <code>type</code> . See <a href="#">global_envelope_test</a> for the defaults and available options. (The test here uses <code>alternative="two.sided"</code> and <code>nstep=1</code> (when relevant), but all the other specifications are to be given in ...)

**Details**

This function can be used to perform one-way graphical functional ANOVA tests described in Mrkvicka et al. (2020). Both 1d and 2d functions are allowed in curve sets.

The tests assume that there are  $J$  groups which contain  $n_1, \dots, n_J$  functions  $T_{ij}, i = \dots, J, j = 1, \dots, n_j$ . The functions should be given in the argument `curve_set`, and the groups in the argument `groups`. The tests assume that  $T_{ij}, i = 1, \dots, n_j$  is an iid sample from a stochastic process with mean function  $\mu_j$  and covariance function  $\gamma_j(s, t)$  for  $s, t$  in  $\mathbb{R}$  and  $j = 1, \dots, J$ .

To test the hypothesis

$$H_0 : \mu_1(r) \equiv \mu_2(r) \equiv \dots \equiv \mu_J(r),$$

you can use the test function

$$\mathbf{T} = (\bar{T}_1(\mathbf{r}), \bar{T}_2(\mathbf{r}), \dots, \bar{T}_J(\mathbf{r}))$$

where  $\bar{T}_i(\mathbf{r})$  is a vector of mean values of functions in the group  $j$ . This test function is used when `contrasts = FALSE` (default).

The hypothesis can equivalently be written as

$$H_0 : \mu_i(r) - \mu_j(r) = 0, i = 1, \dots, J-1, j = 1, \dots, J.$$

and, alternatively, one can use the test function (vector) taken to consist of the differences of the group averages,

$$\mathbf{T}' = (\bar{T}_1(\mathbf{r}) - \bar{T}_2(\mathbf{r}), \bar{T}_1(\mathbf{r}) - \bar{T}_3(\mathbf{r}), \dots, \bar{T}_{J-1}(\mathbf{r}) - \bar{T}_J(\mathbf{r})).$$

The choice is available with the option `contrasts = TRUE`. This test corresponds to the post-hoc test done usually after an ANOVA test is significant, but it can be directed tested by means of the combined rank test (Mrkvička et al., 2017) with this test vector.

The test as such assumes that the variances are equal across the groups of functions. To deal with unequal variances, the differences are rescaled as the first step as follows

$$S_{ij}(r) = \frac{T_{ij}(r) - \bar{T}(r)}{\sqrt{\text{Var}(T_j(r))}} \sqrt{\text{Var}(T(r))} + \bar{T}(r)$$

where  $\bar{T}(\mathbf{r})$  is the overall sample mean and  $\sqrt{\text{Var}(T(r))}$  is the overall sample standard deviation. This scaling of the test functions can be obtained by giving the argument `variances = "unequal"`.

## References

- Mrkvička, T., Myllymäki, M., Jilek, M. and Hahn, U. (2020) A one-way ANOVA test for functional data with graphical interpretation. *Kybernetika* 56 (3), 432-458. doi: 10.14736/kyb-2020-3-0432
- Mrkvička, T., Myllymäki, M., and Hahn, U. (2017). Multiple Monte Carlo testing, with applications in spatial point processes. *Statistics and Computing* 27 (5): 1239-1255. doi:10.1007/s11222-016-9683-9
- Myllymäki, M. and Mrkvička, T. (2024). GET: Global envelopes in R. *Journal of Statistical Software* 111(3), 1-40. doi: 10.18637/jss.v111.i03

## See Also

[frank.fanova](#)

## Examples

```

#-- NOx levels example (see for details Myllymaki and Mrkvicka, 2020)
if(require("fda.usc", quietly=TRUE)) {
  # Prepare data
  data("poblenou")
  fest <- poblenou$df$day.festive; week <- as.integer(poblenou$df$day.week)
  Type <- vector(length=length(fest))
  Type[fest == 1 | week >= 6] <- "Free"
  Type[fest == 0 & week %in% 1:4] <- "MonThu"
  Type[fest == 0 & week == 5] <- "Fri"
  Type <- factor(Type, levels = c("MonThu", "Fri", "Free"))

  # (log) Data as a curve_set
  cset <- curve_set(r = 0:23,
    obs = t(log(poblenou[['nox']][['data']]))))
  # Graphical functional ANOVA
  nsim <- 2999

  res.c <- graph.fanova(nsim = nsim, curve_set = cset,
    groups = Type, variances = "unequal",
    contrasts = TRUE)
  plot(res.c) + ggplot2::labs(x = "Hour", y = "Diff.")
}

#-- Centred government expenditure centralization ratios example
# This is an example analysis of the centred GEC in Mrkvicka et al.
data("cgec")

# Number of simulations
nsim <- 2499 # increase to reduce Monte Carlo error

# Test for unequal lag 1 covariances
res.cov1 <- graph.fanova(nsim = nsim, curve_set = cgec$cgec,
  groups = cgec$group,
  test.equality = "cov", cov.lag = 1)
plot(res.cov1)
# Add labels
plot(res.cov1, labels = paste("Group ", 1:3, sep="")) +
  ggplot2::xlab(substitute(paste(italic(i), "(", j, ")", sep=""), list(i="r", j="Year")))
# Test for equality of variances among groups
res.var <- graph.fanova(nsim = nsim, curve_set = cgec$cgec,
  groups = cgec$group,
  test.equality = "var")
plot(res.var)

# Test for equality of means assuming equality of variances
# a) using 'means'
res <- graph.fanova(nsim = nsim, curve_set = cgec$cgec,
  groups = cgec$group,
  variances = "equal", contrasts = FALSE)
plot(res)

```

```
# b) using 'contrasts'
res2 <- graph.fanova(nsim = nsim, curve_set = cgec$cgec,
                    groups = cgec$group,
                    variances = "equal", contrasts = TRUE)

plot(res2)

# Image set examples
data("imageset3")

res <- graph.fanova(nsim = 19, # Increase nsim for serious analysis!
                  curve_set = imageset3$image_set,
                  groups = imageset3$Group)
plot(res, what = c("obs", "lo.sign", "hi.sign"), sign.type = "col")
```

graph.flm

*Graphical functional GLM***Description**

Non-parametric graphical tests of significance in functional general linear model (GLM)

**Usage**

```
graph.flm(
  nsim,
  formula.full,
  formula.reduced,
  curve_sets,
  factors = NULL,
  contrasts = FALSE,
  lm.args = NULL,
  GET.args = NULL,
  mc.cores = 1L,
  mc.args = NULL,
  cl = NULL,
  savefuns = FALSE,
  fast = TRUE
)
```

**Arguments**

nsim	The number of random permutations.
formula.full	The formula specifying the general linear model, see formula in <a href="#">lm</a> .
formula.reduced	The formula of the reduced model with nuisance factors only. This model should be nested within the full model.

curve_sets	A named list of sets of curves giving the dependent variable (Y), and possibly additionally factors whose values vary across the argument values of the functions. The dimensions of the elements should match with each other. Note that factors that are fixed across the functions can be given in the argument factors. Also <a href="#">fdata</a> objects allowed.
factors	A data frame of factors. An alternative way to specify factors when they are constant for all argument values of the functions. The number of rows of the data frame should be equal to the number of curves. Each column should specify the values of a factor.
contrasts	Logical or NULL. FALSE, TRUE and NULL specify the three test functions as described in description part of this help file.
lm.args	A named list of additional arguments to be passed to <a href="#">lm</a> . See details.
GET.args	A named list of additional arguments to be passed to <a href="#">global_envelope_test</a> , e.g. <code>typeone</code> specifies the type of multiple testing control, FWER or FDR. See <a href="#">global_envelope_test</a> for the defaults and available options.
mc.cores	The number of cores to use, i.e. at most how many child processes will be run simultaneously. Must be at least one, and parallelization requires at least two cores. On a Windows computer <code>mc.cores</code> must be 1 (no parallelization). For details, see <a href="#">mclapply</a> , for which the argument is passed. Parallelization can be used in generating simulations and in calculating the second stage tests.
mc.args	A named list of additional arguments to be passed to <a href="#">mclapply</a> . Only relevant if <code>mc.cores</code> is more than 1.
c1	Allows parallelization through the use of <a href="#">parLapply</a> (works also in Windows), see the argument <code>c1</code> there, and examples.
savefuns	Logical. If TRUE, then the functions from permutations are saved to the attribute <code>simfuns</code> .
fast	Logical. See details.

## Details

The function `graph.flm` performs the graphical functional GLM of Mrkvička et al. (2021), described also in Section 3.6 of Myllymäki and Mrkvička (2024) (type `vignette("GET")` in R). This is a nonparametric graphical test of significance of a covariate in functional GLM. The test is able to find not only if the factor of interest is significant, but also which functional domain is responsible for the potential rejection. In the case of functional multi-way main effect ANOVA or functional main effect ANCOVA models, the test is able to find which groups differ (and where they differ). In the case of functional factorial ANOVA or functional factorial ANCOVA models, the test is able to find which combination of levels (which interactions) differ (and where they differ). The described tests are global envelope tests applied in the context of GLMs. The Freedman-Lane algorithm (Freedman and Lane, 1983) is applied to permute the functions (to obtain the simulations under the null hypothesis of "no effects"); consequently, the test approximately achieves the desired significance level.

The specification of the full and reduced formulas is important. The reduced model should be nested within the full model. The full model should include in addition to the reduced model the interesting factors whose effects are under investigation. The implementation to find the coefficients of the interesting factors is based on [dummy.coef](#) and the restrictions there apply.

The regression coefficients serve as test functions in the graphical functional GLM. For a continuous interesting factor, the test function is its regression coefficient across the functional domain. For a discrete factor, there are three possibilities that are controlled by the arguments `contrasts`. If `contrasts = FALSE`, then the test statistic is the function/long vector where the coefficients related to all levels of the factor are joined together. If `contrasts = TRUE`, then the differences between the same coefficients are considered instead. Given the coefficients in a specific order that is obtained through the use of `lm` and `dummy.coef`, the differences are taken for couples  $i$  and  $j$  where  $i < j$  and reducing  $j$  from  $i$  (e.g. for three groups 1,2,3, the contrasts are 1-2, 1-3, 2-3). If `contrasts = NULL` the coefficients given by `lm` are used directly.

There are different versions of the implementation depending on the application. Given that the argument `fast` is `TRUE`, then

- If all the covariates are continuous or `contrasts = NULL` and `lm.args = NULL` the regression coefficients are computed using the normal equation approach instead of using `lm`.
- Otherwise, if all the covariates are constant across the functions, i.e. they can be provided in the argument `factors`, then a linear model is fitted separately by least-squares estimation to the data at each argument value of the functions fitting a multiple linear model by `lm`. The possible extra arguments passed in `lm.args` to `lm` must be of the form that `lm` accepts for fitting a multiple linear model. In the basic case, no extra arguments are needed.
- Otherwise, if some of the covariates vary across the space and there are user specified extra arguments given in `lm.args`, then the implementation fits a linear model at each argument value of the functions using `lm`, which can be rather slow. The arguments `lm.args` are passed to `lm` for fitting each linear model.

By setting `fast = FALSE`, it is possible to use the slow version (third option) for any case. Usually this is not desired.

## Value

A `global_envelope` or `combined_global_envelope` object, which can be printed and plotted directly.

## References

- Mrkvička, T., Roskovec, T. and Rost, M. (2021) A nonparametric graphical tests of significance in functional GLM. *Methodology and Computing in Applied Probability* 23, 593-612. doi: 10.1007/s11009-019-09756-y
- Myllymäki, M. and Mrkvička, T. (2024). GET: Global envelopes in R. *Journal of Statistical Software* 111(3), 1-40. doi: 10.18637/jss.v111.i03
- Freedman, D., & Lane, D. (1983) A nonstochastic interpretation of reported significance levels. *Journal of Business & Economic Statistics*, 1(4), 292-298. doi:10.2307/1391660

## Examples

```
data("rimov")
res <- graph.flm(nsim=19, # Increase the number of simulations for serious analysis!
  formula.full = Y~Year,
  formula.reduced = Y~1,
  curve_sets = list(Y=rimov), factors = data.frame(Year = 1979:2014))
```

```

plot(res)

# Test if there is a change in the slope in 1994,
# i.e. the full model is  $T = a + b \cdot \text{year} + c \cdot \text{year} : \text{Interv}$ ,
# where Interv is a dummy variable indicating the pre-intervention
# period (coded 0) or the post-intervention period (coded 1)
Year <- 1979:2014
res <- graph.flm(nsim = 19, # Increase the number of simulations for serious analysis!
  formula.full = Y ~ Year + Year:Interv,
  formula.reduced = Y ~ Year,
  curve_sets = list(Y=rimov),
  factors = data.frame(Year = Year,
    Interv = factor(c(rep(0,times=1994-1979+1), rep(1,times=2014-1994))),
    levels=0:1)),
  contrasts = NULL)
plot(res)

# An example of testing the joint effect of a discrete and a continuous variable
nsim <- 999

data("GDPtax")
factors.df <- data.frame(Group = GDPtax$Group, Tax = GDPtax$Profittax)
res.tax_within_group <- graph.flm(nsim = nsim,
  formula.full = Y~Group+Tax+Group:Tax,
  formula.reduced = Y~Group+Tax,
  curve_sets = list(Y=GDPtax$GDP),
  factors = factors.df)
plot(res.tax_within_group)

# Image data examples

data("abide_9002_23")
iset <- abide_9002_23$curve_set

# Testing the discrete factor 'group' with contrasts
# (Use contrasts = FALSE for 'means'; and for continuous factors)
res <- graph.flm(nsim = 19, # Increase nsim for serious analysis!
  formula.full = Y ~ Group + Sex + Age,
  formula.reduced = Y ~ Sex + Age,
  curve_sets = list(Y = iset),
  factors = abide_9002_23[['factors']],
  contrasts = TRUE,
  GET.args = list(type = "area"))
plot(res)

# Examples of modifying 2d plots
plot(res, sign.col="white") + ggplot2::scale_fill_viridis_c(option="magma")
plot(res, sign.col="white") + ggplot2::scale_fill_viridis_c(option="magma") +
  ggplot2::scale_radius(range = 2*c(1, 6))
plot(res, what=c("obs", "lo", "hi", "lo.sign", "hi.sign"))
plot(res, what=c("obs", "lo", "hi", "lo.sign", "hi.sign"), sign.type="col")

```

---

hotspots.MatClustlpp    *Hot spots detection for Matern point process - parallel version*

---

## Description

See the hotspots vignette available by starting R, typing `library("GET")` and `vignette("GET")`.

## Usage

```
hotspots.MatClustlpp(
  PP,
  formula,
  R,
  alpha,
  data,
  sigma = 250,
  nsim = 10000,
  ncores = 1L,
  ...
)
```

## Arguments

PP	Input, a point pattern object (ppp) of spatstat.
formula	An R formula to estimate the first order model. This formula can contain objects of full size. PP should be on the left side of the formula.
R	Cluster radius parameter of the Matern cluster process.
alpha	Parameter related to mean number of points per cluster.
data	Data from where the formula takes objects. Must be acceptable by the function <code>lppm</code> of <code>spatstat.linnet</code> .
sigma	To be passed to <code>density.lpp</code> .
nsim	Number of simulations to be performed.
ncores	Number of cores used for computations. Default to 1. If NULL, all available cores are used.
...	Additional parameters to be passed to <code>fdr_envelope</code> .

## References

Mrkvička et al. (2023). Hotspot detection on a linear network in the presence of covariates: A case study on road crash data. DOI: 10.2139/ssrn.4627591



---

hotspots.poislpp*Hot spots detection for Poisson point process - parallel version*

---

## Description

See the hotspots vignette available by starting R, typing `library("GET")` and `vignette("GET")`.

## Usage

```
hotspots.poislpp(  
  PP,  
  formula,  
  data,  
  sigma = 250,  
  nsim = 10000,  
  ncores = 1L,  
  ...  
)
```

## Arguments

PP	The point pattern living in a network.
formula	A formula for the intensity.
data	Data from where the formula takes objects. Must be acceptable by the function <code>lppm</code> of <code>spatstat.linnet</code> .
sigma	To be passed to <code>density.lpp</code> .
nsim	Number of simulations to be performed.
ncores	Number of cores used for computations. Default to 1. If NULL, all available cores are used.
...	Additional parameters to be passed to <a href="#">fdr_envelope</a> .

## References

Mrkvička et al. (2023). Hotspot detection on a linear network in the presence of covariates: A case study on road crash data. DOI: 10.2139/ssrn.4627591

imageset3

*A simulated set of images***Description**

A simulated set of images with a categorical factor (with three levels)

**Usage**

```
data("imageset3")
```

**Format**

A list of the `image_set` containing the simulated images, and the discrete group factor in the list component `Group`.

**Details**

We considered a categorical factor `Group` obtaining the values 0, 1 or 2 according to the group to which the image belongs to (10 images in each of the three groups). The images were simulated in the square window  $[-1,1]^2$  from the general linear model (GLM)

$$Y(r) = \exp(-10 \cdot \|r\|) \cdot (1 + \mathbf{1}(g = 2)) + \epsilon(r),$$

where  $\|r\|$  denotes the Euclidean distance of the pixel to the origin,  $g$  is the group and the error stems from an inhomogeneous distribution over  $\mathbb{R}^2$  with the normal and bimodal errors in the middle and periphery of the image:

$$\epsilon(r) = \mathbf{1}(\|r\| \leq 0.5)G(r) + \mathbf{1}(\|r\| > 0.5)\frac{1}{2}G(r)^{1/5},$$

where  $G(r)$  is a Gaussian random field with the exponential correlation structure with scale parameter 0.15 and standard deviation 0.2. Consequently, the first two groups (0,1) have the same mean, while a bigger bump appears in the third group (2) in the middle of the image.

**References**

Mrkvička, T., Myllymäki, M., Kuronen, M. and Narisetty, N. N. (2022) New methods for multiple testing in permutation inference for the general linear model. *Statistics in Medicine* 41(2), 276-297. doi: 10.1002/sim.9236

**See Also**

[graph.fanova](#), [frank.fanova](#)

**Examples**

```
data("imageset3")
plot(imageset3$image_set, idx=c(1:5, 11:15, 21:25), ncol=5)

# Colors can be changed as follows:
plot(imageset3$image_set, idx=c(1:5, 11:15, 21:25), ncol=5) +
  ggplot2::scale_fill_gradient(low="black", high="white")
```

---

is.curve_set	<i>Check class.</i>
--------------	---------------------

---

**Description**

Check class.

**Usage**

```
is.curve_set(x)
```

**Arguments**

x                      An object to be checked.

---

MatClust.lppm	<i>Fit a Matern cluster point process to a point pattern dataset on a linear network</i>
---------------	--

---

**Description**

Fit a Matern cluster point process to a point pattern dataset on a linear network. This function is provided in **GET** to support [hotspots.MatClustlpp](#). See the hotspots vignette available by starting R, typing `library("GET")` and `vignette("GET")`.

**Usage**

```
MatClust.lppm(
  PP,
  formula,
  subwin = NULL,
  valpha,
  vR,
  data,
  nsim = 10,
  ncores = 1L
)
```

Arguments

PP	Input, a point pattern object (ppp) of spatstat.
formula	An R formula to estimate the first order model. This formula can contain objects of full size. PP should be on the left side of the formula.
subwin	A part of the observation window of PP to be used for estimating the second order structure. NULL means that the full point pattern is used. Typically this is feasible (not too time consuming).
valpha	A vector of parameter values for the parameter alpha of the Matern cluster process.
vR	A vector of parameter values for the parameter R of the Matern cluster process.
data	Data from where the formula takes objects. Must be acceptable by the function lppm of spatstat.linnet.
nsim	The number of simulated Matern cluster point patterns for evaluating the K-function for any alpha and R values.
ncores	Number of cores used for computations. Default to 1. If NULL, all available cores are used.

Details

The function `MatClust.lppm`, can be used to estimate the Matern cluster point pattern with inhomogeneous cluster centers on linear network. This function provides the same outputs as the `pois.lppm` and further estimated parameters alpha and R. The `secondorder` provides again the diagnostics for checking if the clustered model is appropriate. The sample K-function must be close to the K-function of the estimated model (green line). If it is not the case the searching grid for parameters alpha and R that is input in the function must be manipulated to get the a closer result. If the estimated model is adequate one can proceed to the hotspot detection with the use of the function `hotspots.MatClustlpp`. Remark here, that for the estimation of the second order structure a smaller data can be used than for the estimation of the first order structure in order to save the computation time, since the second order is a local characteristics. This smaller window can be specified by `subwin`. Then the full point pattern will be used for estimation of first order intensity and the pattern in subwindow will be used for estimating second order characteristic. The input parameters are the same as in `pois.lppm`. Furthermore, `valpha`, i.e., vector of proposed alphas which should be considered in the optimization, `vR`, i.e., vector of proposed values for R which should be considered in the optimization, must be provided. The user can also specify how many cores should be used in the computation by parameter `ncores`.

---

naturalness	<i>Simulated data set</i>
-------------	---------------------------

---

Description

Simulated data set mimicing stand age distributions in natural, near-natural non-natural forests in the setup of Mrkvička et al. (2024).

**Usage**

```
data("naturalness")
```

**Format**

A data.frame containing the dominant species (three categories: Broadleaf, Conifer, Mixed), naturalness (Natural, Near-natural and Non-natural) and simulated stand age.

**Details**

The numbers of observations in the naturalness groups and in the dominant species groups correspond to the numbers of plots in on rich mineral soils in Finnish Lapland in the data set analysed in Myllymäki et al. (2023). The stand age values are simulated values from the quantile regression model of Mrkvička et al. (2024).

**References**

Myllymäki, M., Tuominen, S., Kuronen, M., Packalen, P. and Kangas, A. (2023) The relationship between forest structure and naturalness in the Finnish national forest inventory. *Forestry: An International Journal of Forest Research*, cpad053. DOI: <https://doi.org/10.1093/forestry/cpad053>

Mrkvička, T., Konstantinou, K., Kuronen, M. and Myllymäki, M. (2024) Global quantile regression. *arXiv:2309.04746 [stat.ME]* DOI: <https://doi.org/10.48550/arXiv.2309.04746>

**See Also**

[global\\_rq](#)

---

partial_forder	<i>Functional ordering in parts</i>
----------------	-------------------------------------

---

**Description**

If the functional data doesn't comfortably fit in memory it is possible to compute functional ordering by splitting the domain of the data (voxels in a brain image), using `partial_forder` on each part and finally combining the results with `combine_forder`.

**Usage**

```
partial_forder(  
  curve_set,  
  measure = c("erl", "rank", "cont", "area"),  
  alternative = c("two.sided", "less", "greater")  
)  
  
combine_forder(ls)
```

**Arguments**

curve_set	A curve_set object, usually a part of a larger curve_set. (No missing or infinite values allowed.)
measure	The measure to use to order the functions from the most extreme to the least extreme one. Must be one of the following: 'rank', 'erl', 'cont', 'area', 'max', 'int', 'int2'. Default is 'erl'.
alternative	A character string specifying the alternative hypothesis. Must be one of the following: "two.sided" (default), "less" or "greater". The last two options only available for types 'rank', 'erl', 'cont' and 'area'.
ls	List of objects returned by partial_forder

**Value**

See [forder](#)

**See Also**

[forder](#)

**Examples**

```
data("abide_9002_23")
res <- lapply(list(1:100, 101:200, 201:261), function(part) {
  set.seed(123) # When using partial_forder, all parts must use the same seed.
  fset <- frank.flm(nsim=99, formula.full = Y ~ Group + Sex + Age,
    formula.reduced = Y ~ Group + Sex,
    curve_sets = list(Y = abide_9002_23$curve_set[part,]),
    factors = abide_9002_23$factors, savefuns = "return")
  partial_forder(fset, measure="erl")
})
combine_forder(res)
```

---

plot.combined\_fboxplot

*Plot method for the class 'combined\_fboxplot'*

---

**Description**

Plot method for the class 'combined\_fboxplot'

**Usage**

```
## S3 method for class 'combined_fboxplot'
plot(
  x,
  labels,
  scales = "free",
```

```

    ncol = 2 + 1 * (length(x) == 3),
    digits = 3,
    outliers = TRUE,
    ...
)

```

### Arguments

x	an 'combined_fboxplot' object
labels	A character vector of suitable length. If dotplot = TRUE (for the level 2 test), then labels for the tests at x-axis; only valid/used when all components of x have the same dimension. Otherwise labels for the separate plots.
scales	See <a href="#">facet_wrap</a> . Use scales = "free" when the scales of the functions in the global envelope vary. scales = "fixed" is a good choice, when you want the same y-axis for all components. A sensible default based on r-values exists.
ncol	The maximum number of columns for the figures. Default 2 or 3, if the length of x equals 3. (Relates to the number of curve_sets that have been combined.)
digits	The number of digits used for printing the p-value or p-interval in the default main.
outliers	Logical. If TRUE, then the functions outside the functional boxplot are drawn.
...	Ignored.

---

```
plot.combined_global_envelope
```

*Plot method for the class 'combined\_global\_envelope'*

---

### Description

This function provides plots for combined global envelopes.

### Usage

```

## S3 method for class 'combined_global_envelope'
plot(
  x,
  labels,
  scales,
  sign.col = "red",
  dotplot = length(x[[1]]$obs) < 5,
  ncol = 2 + 1 * (length(x) == 3),
  digits = 3,
  level = 1,
  ...
)

```

**Arguments**

x	An 'combined_global_envelope' object
labels	A character vector of suitable length. If dotplot = TRUE (for the level 2 test), then labels for the tests at x-axis; only valid/used when all components of x have the same dimension. Otherwise labels for the separate plots.
scales	See <a href="#">facet_wrap</a> . Use scales = "free" when the scales of the functions in the global envelope vary. scales = "fixed" is a good choice, when you want the same y-axis for all components. A sensible default based on r-values exists.
sign.col	The color for the observed curve when outside the global envelope (significant regions). Default to "red". Setting the color to NULL corresponds to no coloring. If the object contains several envelopes, the coloring is done for the widest one.
dotplot	Logical. If TRUE, then instead of envelopes a dot plot is done. Suitable for low dimensional test vectors.
ncol	The maximum number of columns for the figures. Default 2 or 3, if the length of x equals 3. (Relates to the number of curve_sets that have been combined.)
digits	The number of digits used for printing the p-value or p-interval in the default main.
level	1 or 2. In the case of two-step combined tests (with several test functions), two different plots are available: 1 for plotting the combined global envelopes (default and most often wanted) or 2 for plotting the second level test result.
...	Ignored in most cases. If dotplot = TRUE, then parameters can be passed to <a href="#">arrow</a> , e.g. length = unit(0.25, "cm").

**Details**

Plotting method for the class 'combined\_global\_envelope', i.e. combined envelopes for 1d functions.

**See Also**

[central\\_region](#)

---

plot.combined\_global\_envelope2d

*Plotting function for combined 2d global envelopes*

---

**Description**

If fixedscales is FALSE (or 0) all images will have separate scale. If fixedscales is TRUE (or 1) each x[[i]] will have a common scale. If fixedscales is 2 all images will have common scale.

If more than one envelope has been calculated (corresponding to several coverage/alpha), only the largest one is plotted.



**Usage**

```
## S3 method for class 'combined_global_envelope2d'
plot(
  x,
  fixedscales = 2,
  labels,
  what = c("obs.sign", "obs", "lo", "hi", "lo.sign", "hi.sign"),
  sign.type = c("circles", "contour", "col"),
  sign.col = c("blue", "red"),
  transparency = 155/255,
  digits = 3,
  ...
)
```

**Arguments**

<code>x</code>	A 'global_envelope' object for two-dimensional functions
<code>fixedscales</code>	0, 1 or 2. See details.
<code>labels</code>	A character vector of suitable length giving the labels for the separate plots. Default exists. This parameter allows replacing the default.
<code>what</code>	Character vector specifying what information should be plotted for 2d functions. A combination of: Observed ("obs"), upper envelope ("hi"), lower envelope ("lo"), observed with significantly higher values highlighted ("hi.sign"), observed with significantly lower values highlighted ("lo.sign"), observed with significantly (lower and higher) values highlighted ("obs.sign"). Default to the last one. Combination <code>c("obs", "lo", "hi", "lo.sign", "hi.sign")</code> can also be of interest (earlier default).
<code>sign.type</code>	Either "col" for color showing the significant region, or "contour" for colored contour showing the significant region, or "circles" for plotting circles at locations where the observed function exceeds the envelope: diameters proportional to $(\text{obs}-\text{hi})/(\text{hi}-\text{lo})$ for values above the envelope and $(\text{lo}-\text{obs})/(\text{hi}-\text{lo})$ for values below the envelope. In the one-sided (testing) case, the divisors are instead $(\text{hi}-\text{central})$ (case 'greater') and $(\text{central}-\text{lo})$ (case 'less'). Default to "circles".
<code>sign.col</code>	A vector of length two giving the colors for significant parts below the envelope (first value) and above the envelope (second value).
<code>transparency</code>	A number between 0 and 1 (default 155/255, 60 Similar to alpha of <a href="#">rgb</a> . Used in plotting the significant regions for 2d functions.
<code>digits</code>	The number of digits used for printing the p-value or p-interval in the default main.
<code>...</code>	Ignored.

**Examples**

```
data("abide_9002_23")
iset <- subset(abide_9002_23[['curve_set']], 1:50)
factors <- abide_9002_23[['factors']][1:50,]
```

```

res <- graph.flm(nsim = 19, # Increase nsim for serious analysis!
  formula.full = Y ~ Group + Sex + Age,
  formula.reduced = Y ~ Sex + Age,
  curve_sets = list(Y=iset), factors = factors,
  contrasts = FALSE, GET.args = list(type="area"))
plot(res)

plot(res, what=c("obs", "hi"))

plot(res, what=c("hi", "lo"), fixed scales=1)

plot(res, what=c("obs", "lo", "hi"), fixed scales=FALSE)

if(requireNamespace("gridExtra", quietly=TRUE)) {
  # Edit style of "fixed scales = 2" plots
  plot(res, what=c("obs", "hi")) + ggplot2::theme_minimal()
  plot(res, what=c("obs", "hi")) + ggplot2::theme_bw()

  # Edit style (e.g. theme) of "fixed scales = 1 or 0" plots
  gs <- lapply(res, function(x, what) { plot(x, what=what) +
    ggplot2::ggtitle("") }, what=c("obs", "hi"))
  gridExtra::grid.arrange(grobs=gs, ncol=1, top="My main")

  gs <- outer(res, c("obs", "hi"), FUN=Vectorize(function(res, what)
    list(plot(res, what=what) + ggplot2::ggtitle("") +
      ggplot2::theme(axis.ticks=ggplot2::element_blank(),
        axis.text=ggplot2::element_blank(), axis.title=ggplot2::element_blank()))))
  gridExtra::grid.arrange(grobs=t(gs))
}

```

---

plot.curve\_set

*Plot method for the class 'curve\_set'*


---

## Description

Plot method for the class 'curve\_set'

## Usage

```

## S3 method for class 'curve_set'
plot(x, idx, col_idx, idx_name = "", col = "grey70", ...)

```

## Arguments

x                      An curve\_set object.

idx	Indices of functions to highlight with color col_idx. Default to the observed function, if there is just one. The legend of curves' colours is shown if indices are given or x contains one observed function. See examples to remove the legend if desired.
col_idx	A color for the curves to highlight, or a vector of the same length as idx containing the colors for the highlighted functions. Default exists.
idx_name	A variable name to be printed with the highlighted functions' idx. Default to empty.
col	The basic color for the curves (which are not highlighted).
...	Ignored.

**See Also**

[create\\_curve\\_set](#)

**Examples**

```
cset <- curve_set(r = 1:10, obs = matrix(runif(10*5), ncol=5))
plot(cset)
# Highlight some functions
plot(cset, idx=c(1,3))
plot(cset, idx=c(1,3), col_idx=c("black", "red"))
# Change legend
plot(cset, idx=c(1,3), col_idx=c("black", "red"), idx_name="Special functions")
plot(cset, idx=c(1,3)) + ggplot2::theme(legend.position="bottom")
# Add labels
plot(cset, idx=c(1,3)) + ggplot2::labs(x="x", y="Value")
# and title
plot(cset) + ggplot2::labs(title="Example curves", x="x", y="Value")
# A curve_set with one observed function (other simulated)
if(requireNamespace("mvtnorm", quietly=TRUE)) {
  cset <- curve_set(obs = c(-1.6, 1.6),
    sim = t(mvtnorm::rmvnorm(200, c(0,0), matrix(c(1,0.5,0.5,1), 2, 2))))
  plot(cset)
  # Remove legend
  plot(cset) + ggplot2::theme(legend.position="none")
}
```

---

plot.curve\_set2d      *Plot method for the class 'curve\_set2d'*

---

**Description**

Plot method for the class 'curve\_set2d', i.e. two-dimensional functions

**Usage**

```
## S3 method for class 'curve_set2d'
plot(x, idx = 1, ncol = 2 + 1 * (length(idx) == 3), ...)
```

**Arguments**

x	An curve_set2d object
idx	Indices of 2d functions to plot.
ncol	The maximum number of columns for the figures. Default 2 or 3, if the length of x equals 3. (Relates to the number of curve_sets that have been combined.)
...	Ignored.

**Examples**

```
data("abide_9002_23")
plot(abide_9002_23$curve_set, idx=c(1, 27))
```

---

plot.fboxplot	<i>Plot method for the class 'fboxplot'</i>
---------------	---

---

**Description**

Plot method for the class 'fboxplot'

**Usage**

```
## S3 method for class 'fboxplot'
plot(x, digits = 3, outliers = TRUE, ...)
```

**Arguments**

x	an 'fboxplot' object
digits	The number of digits used for printing the p-value or p-interval in the default main.
outliers	Logical. If TRUE, then the functions outside the functional boxplot are drawn.
...	Ignored.

**Examples**

```
if(requireNamespace("fda", quietly=TRUE)) {
  years <- paste(1:18)
  curves <- fda::growth[['hgtf']][years,]
  # Heights
  cset1 <- curve_set(r = as.numeric(years),
                    obs = curves)
  bp <- fBoxplot(cset1, coverage=0.50, type="area", factor=1)
  plot(bp)
  plot(bp) + ggplot2::theme(legend.position="bottom")
  plot(bp) + ggplot2::theme(legend.position="none")
  plot(bp, plot_outliers=FALSE)
}
```

---

plot.fclust	<i>Plot method for the class 'fclust'</i>
-------------	---

---

## Description

Plot method for the 'fclust' objects returned by [fclustering](#).

## Usage

```
## S3 method for class 'fclust'
plot(x, plotstyle = c("marginal", "joined"), coverage = 0.5, nstep, ncol, ...)
```

## Arguments

x	An 'fclust' object.
plotstyle	The resulting central regions of clusters can be plotted by sorting the appropriate curve_set only 'marginal' or by sorting the joined list of curve_set objects 'joined'. If 'joined' is used the shown central regions corresponds to the joined ordering used to cluster the functional data. If 'marginal' is used the shown central regions do not correspond to the joined ordering used to cluster the functional data, but better express the shape of cluster with respect to given curve_set.
coverage	The coverage of central regions to be used to show the clusters.
nstep	1 or 2 for how to construct a combined (joined) global envelope if there are more than one sets of curves. Default to 1, if the numbers of points where the curves are observed (r) are the same in each set, and 2 otherwise.
ncol	The number of columns in the graphical output, when there is just one set of curves that has been ordered. If not given, c(1, k+1) is used, which gives all plots in one row. For more sets of curves, the rows are fixed to correspond to the sets (one row for each set).
...	Ignored.

## Details

The clusters are shown respectively for each curve\_set. Thus for each curve\_set the panel with all the medoids is shown followed by all clusters represented by central region, medoid and all curves belonging to it.

For all sources, the function plots the deepest curves for all clusters and the deepest curve of each cluster together with the desired central region and all the curves of the group.

## References

Dai, W., Athanasiadis, S., Mrkvička, T. (2021) A new functional clustering method with combined dissimilarity sources and graphical interpretation. Intech open, London, UK. DOI: 10.5772/intechopen.100124

---

plot.global\_envelope    *Plot method for the class 'global\_envelope'*

---

## Description

Plot method for the class 'global\_envelope'

## Usage

```
## S3 method for class 'global_envelope'
plot(
  x,
  dotplot = length(x$r) < 10,
  sign.col = "red",
  labels = NULL,
  digits = 3,
  ...
)
```

## Arguments

x	An 'global_envelope' object
dotplot	Logical. If TRUE, then instead of envelopes a dot plot is done. Suitable for low dimensional test vectors. Default: TRUE if the dimension is less than 10, FALSE otherwise.
sign.col	The color for the observed curve when outside the global envelope (significant regions). Default to "red". Setting the color to NULL corresponds to no coloring. If the object contains several envelopes, the coloring is done for the widest one.
labels	A character vector of suitable length. If dotplot = TRUE, then labels for the tests at x-axis.
digits	The number of digits used for printing the p-value or p-interval in the default main.
...	Ignored.

## Details

If several envelopes have been computed, their are plotted in different grey scales so that the smallest envelope has the darkest color and the widest envelope consist of all grey scales with the lightest color in the outskirts.

## See Also

[central\\_region](#), [global\\_envelope\\_test](#)

**Examples**

```

if(require("spatstat.explore", quietly=TRUE)) {
  X <- unmark(spruces)
  nsim <- 1999 # Number of simulations

  env <- envelope(X, fun="Kest", nsim=nsim,
                 savefuns=TRUE, # save the functions
                 correction="translate", # edge correction for K
                 simulate=expression(runifpoint(ex=X))) # Simulate CSR
  res <- global_envelope_test(env, type="erl")

  # Default plot
  plot(res)
  # Plots can be edited, e.g.
  # Remove legend
  plot(res) + ggplot2::theme(legend.position="none")
  # Change its position
  plot(res) + ggplot2::theme(legend.position="right")
  # Change the outside color
  plot(res, sign.col="#5DC863FF")
  plot(res, sign.col=NULL)
  # Change default title and x- and y-labels
  plot(res) + ggplot2::labs(title="95% global envelope", x="x", y="f(x)")

  # Prior to the plot, you can set your preferred ggplot theme by theme_set
  old <- ggplot2::theme_set(ggplot2::theme_bw())
  plot(res)

  # Do other edits, e.g. turn off expansion with the default limits
  plot(res) + ggplot2::coord_cartesian(expand=FALSE)

  # Go back to the old theme
  ggplot2::theme_set(old)

  # If you are working with the R package spatstat and its envelope-function,
  # you can obtain global envelope plots in the style of spatstat using plot.fv:
  plot.fv(res)
}

```

---

plot.global\_envelope2d

*Plotting function for 2d global envelopes*


---

**Description**

If more than one envelope has been calculated (corresponding to several coverage/alpha), only the largest one is plotted.

**Usage**

```
## S3 method for class 'global_envelope2d'
plot(
  x,
  fixedscales = TRUE,
  what = c("obs.sign", "obs", "lo", "hi", "lo.sign", "hi.sign"),
  sign.type = c("circles", "contour", "col"),
  sign.col = c("blue", "red"),
  transparency = 155/255,
  digits = 3,
  ...
)
```

**Arguments**

<code>x</code>	A 'global_envelope' object for two-dimensional functions
<code>fixedscales</code>	Logical. TRUE for the same scales for all images.
<code>what</code>	Character vector specifying what information should be plotted for 2d functions. A combination of: Observed ("obs"), upper envelope ("hi"), lower envelope ("lo"), observed with significantly higher values highlighted ("hi.sign"), observed with significantly lower values highlighted ("lo.sign"), observed with significantly (lower and higher) values highlighted ("obs.sign"). Default to the last one. Combination c("obs", "lo", "hi", "lo.sign", "hi.sign") can also be of interest (earlier default).
<code>sign.type</code>	Either "col" for color showing the significant region, or "contour" for colored contour showing the significant region, or "circles" for plotting circles at locations where the observed function exceeds the envelope: diameters proportional to (obs-hi)/(hi-lo) for values above the envelope and (lo-obs)/(hi-lo) for values below the envelope. In the one-sided (testing) case, the divisors are instead (hi-central) (case 'greater') and (central-lo) (case 'less'). Default to "circles".
<code>sign.col</code>	A vector of length two giving the colors for significant parts below the envelope (first value) and above the envelope (second value).
<code>transparency</code>	A number between 0 and 1 (default 155/255, 60 Similar to alpha of <a href="#">rgb</a> ). Used in plotting the significant regions for 2d functions.
<code>digits</code>	The number of digits used for printing the p-value or p-interval in the default main.
<code>...</code>	Ignored.

**See Also**

[graph.flm](#)



---

pois.lppm	<i>Fit a Poisson point process model to a point pattern dataset on a linear network</i>
-----------	---

---

## Description

Fit a Poisson point process model to a point pattern dataset on a linear network. This function is provided in **GET** to support [hotspots.poislpp](#) and [hotspots.MatClustlpp](#). See the hotspots vignette available by starting R, typing `library("GET")` and `vignette("GET")`.

## Usage

```
pois.lppm(PP, formula, data, subwin = NULL, r_max = NULL)
```

## Arguments

PP	Input, a point pattern object (ppp) of spatstat.
formula	An R formula to estimate the first order model. This formula can contain objects of full size. PP should be on the left side of the formula.
data	Data from where the formula takes objects. Must be acceptable by the function lppm of spatstat.linnet.
subwin	A part of the observation window of PP to be used for estimating the second order structure. NULL means that the full point pattern is used. Typically this is feasible (not too time consuming).
r_max	The maximum distance on which the K-function is evaluated. Default is computed as $\sqrt{A}/10$ where $A$ is the area of the window of observation of $X$ .

## Details

The function `pois.lppm`, can be used to estimate the inhomogeneous Poisson point process model on linear network. This function provides the `firstordermodel`, i.e. the regression model of dependence of crashes on the spatial covariates, EIP, i.e. estimated inhomogeneous intensity from the data and `secondorder`, i.e. estimation of the inhomogeneous K-function. The plot of the `secondorder` provides diagnostics, if the model is adequate for the data. If the estimated  $K$ -function lies close to the theoretical line, the data does not report any clustering, and the function `hotspots.poislpp` can be used for final hotspots detection. If the estimated K-function does not lie close to the theoretical line, and it is above, the data report clustering, and the a clustered point pattern model must be fitted to the data and hotspots detected using this clustered model instead.

---

popgrowthmillion	<i>Population growth</i>
------------------	--------------------------

---

**Description**

Population growth

**Usage**

```
data("popgrowthmillion")
```

**Format**

A matrix, where each row corresponds to a year and each column to a country. Column names correspond to the countries, and row names to the years.

**Details**

This dataset includes population growth, i.e. population at the end of the year divided by population at the beginning of the year, in 134 countries in years from 1950 to 2015. The dataset includes only countries over million inhabitants in 1950. The data were extracted from the supplement of Nagy et al. (2017) distributed under the GPL-2 license.

**References**

Nagy, S., I. Gijbels, and D. Hlubinka (2017). Depth-based recognition of shape outlying functions. *Journal of Computational and Graphical Statistics* 26 (4), 883-893.

---

```
print.combined_fboxplot
```

*Print method for the class 'combined\_fboxplot'*

---

**Description**

Print method for the class 'combined\_fboxplot'

**Usage**

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

**Arguments**

x	an 'combined_fboxplot' object
...	Ignored.

---

```
print.combined_global_envelope
```

*Print method for the class 'combined\_global\_envelope'*

---

### Description

Print method for the class 'combined\_global\_envelope'

### Usage

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

### Arguments

x	A 'combined_global_envelope' object
...	Ignored.

---

```
print.curve_set
```

*Print method for the class 'curve\_set'*

---

### Description

Print method for the class 'curve\_set'

### Usage

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

### Arguments

x	an 'curve_set' object
...	Passed to <a href="#">str</a> .

---

<code>print.deviation_test</code>	<i>Print method for the class 'deviation_test'</i>
-----------------------------------	--

---

**Description**

Print method for the class 'deviation\_test'

**Usage**

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

**Arguments**

- `x` an 'deviation\_test' object
- `...` Ignored.

---

<code>print.fboxplot</code>	<i>Print method for the class 'fboxplot'</i>
-----------------------------	--

---

**Description**

Print method for the class 'fboxplot'

**Usage**

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

**Arguments**

- `x` an 'fboxplot' object
- `...` Ignored.

---

print.fclust	<i>Print method for the class 'fclust'</i>
--------------	--

---

**Description**

Print method for the 'fclust' objects returned by [fclustering](#).

**Usage**

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

**Arguments**

x	A object of class 'fclust'.
...	Ignored.

---

print.fdr_envelope	<i>Print method for the class 'fdr_envelope'</i>
--------------------	--

---

**Description**

Print method for the class 'fdr\_envelope'

**Usage**

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

**Arguments**

x	An 'fdr_envelope' object
...	Ignored.

---

print.GET\_contingency *Print method for the class 'GET\_contingency'*

---

### Description

Print method for the class 'GET\_contingency'

### Usage

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

### Arguments

x	A 'GET_contingency' object
...	Ignored.

---

print.global\_envelope *Print method for the class 'global\_envelope'*

---

### Description

Print method for the class 'global\_envelope'

### Usage

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

### Arguments

x	A 'global_envelope' object.
...	Ignored.

qdir\_envelope

*Global scaled maximum absolute difference (MAD) envelope tests*

## Description

Performs the global scaled MAD envelope tests, either directional quantile or studentised, or the unscaled MAD envelope test. These tests correspond to calling the function [global\\_envelope\\_test](#) with `type="qdir"`, `type="st"` and `type="unscaled"`, respectively. The functions `qdir_envelope`, `st_envelope` and `unscaled_envelope` have been kept for historical reasons; preferably use [global\\_envelope\\_test](#) with the suitable `type` argument.

## Usage

```
qdir_envelope(curve_set, ...)

st_envelope(curve_set, ...)

unscaled_envelope(curve_set, ...)
```

## Arguments

<code>curve_set</code>	A <a href="#">curve_set</a> object, or an envelope object of <b>spatstat</b> . If an envelope object is given, it must contain the summary functions from the simulated patterns which can be achieved by setting <code>savefuns = TRUE</code> when calling the function of <b>spatstat</b> .
<code>...</code>	Additional parameters to be passed to <a href="#">global_envelope_test</a> .

## Details

The directional quantile envelope test (Myllymäki et al., 2015, 2017) takes into account the unequal variances of the test function  $T(r)$  for different distances  $r$  and is also protected against asymmetry of  $T(r)$ .

The studentised envelope test (Myllymäki et al., 2015, 2017) takes into account the unequal variances of the test function  $T(r)$  for different distances  $r$ .

The unscaled envelope test (Ripley, 1981) corresponds to the classical maximum deviation test without scaling, and leads to envelopes with constant width over the distances  $r$ . Thus, it suffers from unequal variance of  $T(r)$  over the distances  $r$  and from the asymmetry of distribution of  $T(r)$ . We recommend to use the other global envelope tests available, see [global\\_envelope\\_test](#) for full list of alternatives.

## Value

An object of class `global_envelope` of `combined_global_envelope` which can be printed and plotted directly. See [global\\_envelope\\_test](#) for more details.

## References

- Myllymäki, M., Grabarnik, P., Seijo, H. and Stoyan, D. (2015). Deviation test construction and power comparison for marked spatial point patterns. *Spatial Statistics* 11: 19-34. doi: 10.1016/j.spasta.2014.11.004
- Myllymäki, M., Mrkvička, T., Grabarnik, P., Seijo, H. and Hahn, U. (2017). Global envelope tests for spatial point patterns. *Journal of the Royal Statistical Society: Series B (Statistical Methodology)*, 79: 381–404. doi: 10.1111/rssb.12172
- Ripley, B.D. (1981). *Spatial statistics*. Wiley, New Jersey.

## See Also

[global\\_envelope\\_test](#)

## Examples

```
# See more examples in ?global_envelope_test
## Testing complete spatial randomness (CSR)
#-----
if(require("spatstat.explore", quietly=TRUE)) {
  X <- spruces
  nsim <- 999 # Number of simulations

  ## Test for complete spatial randomness (CSR)
  # Generate nsim simulations under CSR, calculate centred L-function for the data and simulations
  env <- envelope(X, fun="Lest", nsim=nsim, savefuns=TRUE,
                 correction="translate", transform=expression(.-r),
                 simulate=expression(runifpoint(ex=X)))
  res_qdir <- qdir_envelope(env) # The directional quantile envelope test
  plot(res_qdir)

  ## Advanced use:
  # Create a curve set, choosing the interval of distances [r_min, r_max]
  curve_set <- crop_curves(env, r_min=1, r_max=8)
  # The directional quantile envelope test
  res_qdir <- qdir_envelope(curve_set); plot(res_qdir)
  # The studentised envelope test
  res_st <- st_envelope(curve_set); plot(res_st)
  # The unscaled envelope test
  res_unscaled <- unscaled_envelope(curve_set); plot(res_unscaled)
}
```

---

rank\_envelope

*The rank envelope test*


---

## Description

The rank envelope test, p-values and global envelopes. The test corresponds to the global envelope test that can be carried out by [global\\_envelope\\_test](#) by specifying the type for which the options "rank", "erl", "cont" and "area" are available. The last three are modifications of the first one



to treat the ties in the extreme rank ordering used in "rank". This function is kept for historical reasons.

### Usage

```
rank_envelope(curve_set, type = "rank", ...)
```

### Arguments

curve_set	A <a href="#">curve_set</a> object, or an envelope object of <b>spatstat</b> . If an envelope object is given, it must contain the summary functions from the simulated patterns which can be achieved by setting <code>savefuns = TRUE</code> when calling the function of <b>spatstat</b> .
type	The type of the global envelope with current options for "rank", "erl", "cont" and "area". If "rank", the global rank envelope accompanied by the p-interval is given (Myllymäki et al., 2017). If "erl", the global rank envelope based on extreme rank lengths accompanied by the extreme rank length p-value is given (Myllymäki et al., 2017, Mrkvička et al., 2018). See details and additional sections thereafter.
...	Additional parameters to be passed to <a href="#">global_envelope_test</a> .

### Details

The "rank" envelope test is a completely non-parametric test, which provides the  $100(1-\alpha)\%$  global envelope for the chosen test function  $T(r)$  on the chosen interval of distances and associated p-values. The other three types are solutions to break the ties in the extreme ranks on which the "rank" envelope test is based on.

Note: The method to break ties for the global type = "rank" envelope (Myllymäki et al., 2017) can be done by the argument `ties` with default to `ties = "erl"` corresponding to the extreme rank length breaking of ties. In this case the global envelope corresponds to the extreme rank measure. If instead choosing type to be "erl", "cont" or "area", then the global envelope corresponds to these measures.

### Value

An object of class `global_envelope` of `combined_global_envelope` which can be printed and plotted directly. See [global\\_envelope\\_test](#) for more details.

### Number of simulations

The global "erl", "cont", "area" envelope tests allow in principle a lower number of simulations to be used than the global "rank" test based on extreme ranks. However, if feasible, we recommend some thousands of simulations in any case to achieve a good power and repeatability of the test. For the global "rank" envelope test, Myllymäki et al. (2017) recommended to use at least 2500 simulations for testing at the significance level  $\alpha = 0.05$  for single function tests, experimented with summary functions for point processes.

## References

Myllymäki, M., Mrkvička, T., Grabarnik, P., Seijo, H. and Hahn, U. (2017). Global envelope tests for spatial point patterns. *Journal of the Royal Statistical Society: Series B (Statistical Methodology)*, 79: 381–404. doi: 10.1111/rssb.12172

Mrkvička, T., Myllymäki, M. and Hahn, U. (2017). Multiple Monte Carlo testing, with applications in spatial point processes. *Statistics & Computing* 27 (5): 1239-1255. doi: 10.1007/s11222-016-9683-9

Mrkvička, T., Myllymäki, M., Jilek, M. and Hahn, U. (2020) A one-way ANOVA test for functional data with graphical interpretation. *Kybernetika* 56 (3), 432-458. doi: 10.14736/kyb-2020-3-0432

## See Also

[global\\_envelope\\_test](#)

## Examples

```
# See ?global_envelope_test for more examples

## Testing complete spatial randomness (CSR)
#-----
if(require("spatstat.explore", quietly=TRUE)) {
  X <- unmark(spruces)
  nsim <- 2499 # Number of simulations

  # Generate nsim simulations under CSR, calculate centred L-function for the data and simulations
  env <- envelope(X, fun="Lest", nsim=nsim, savefuns=TRUE,
                 correction="translate", transform=expression(.-r),
                 simulate=expression(runifpoint(ex=X)))

  # The rank envelope test
  res <- rank_envelope(env)
  # Plot the result.
  plot(res)

  ## Advanced use:
  # Choose the interval of distances [r_min, r_max] (at the same time create a curve_set from 'env')
  curve_set <- crop_curves(env, r_min=1, r_max=7)
  # Do the rank envelope test
  res <- rank_envelope(curve_set); plot(res)
}
```

---

residual

*Residual form of the functions*

---

## Description

Subtract the theoretical function  $S_{H_0}$  or the mean of the functions in the curve set. If the curve\_set object contains already residuals  $T_i(r) - T_0(r)$ , use\_theo ignored and the same object returned.

**Usage**

```
residual(curve_set, use_theo = TRUE)
```

**Arguments**

curve_set	A <a href="#">curve_set</a> object, or an envelope object of <b>spatstat</b> . If an envelope object is given, it must contain the summary functions from the simulated patterns which can be achieved by setting <code>savefuns = TRUE</code> when calling the envelope function.
use_theo	Whether to use the theoretical summary function or the mean of the functions in the <code>curve_set</code> .

**Details**

The mean of the functions in the `curve_set` is the mean of all functions. If `use_theo = TRUE`, but the component `theo` does not exist in the `curve_set`, the mean of the functions is used silently.

**Value**

A curve set object containing residual summary functions. `theo` is no longer included.

---

rimov

---

*Year temperature curves*


---

**Description**

Year temperature curves

**Usage**

```
data("rimov")
```

**Format**

A `curve_set` object with water temperatures in 365 days of the 36 years. The component `curve_set[['r']]` is a vector of days (from 1 to 365), whereas `curve_set[['obs']]` contains the water temperatures such that each column gives year temperatures in a year.

**Details**

The water temperature data sampled at the water level of Rimov reservoir in Czech republic every day for the 36 years between 1979 and 2014.

**References**

Mrkvička, T., Myllymäki, M., Jilek, M. and Hahn, U. (2020) A one-way ANOVA test for functional data with graphical interpretation. *Kybernetika* 56 (3), 432-458. doi: 10.14736/kyb-2020-3-0432

**See Also**

graph.fanova

**Examples**

```
data("rimov")
groups <- factor(c(rep(1, times=12), rep(2, times=12), rep(3, times=12)))
for(i in 1:3)
  assign(paste0("p", i), plot(subset(rimov, groups==i)) +
    ggplot2::labs(title=paste("Group ", i, sep=""), y="Temperature"))
p3
if(require("patchwork", quietly=TRUE))
  p1 + p2 + p3
# See example analysis in ?graph.fanova
```

---

rMatClustlpp

---

*Simulating the Matern cluster process on a linear network*


---

**Description**

Simulating the Matern clusters with parameters  $\alpha$  and  $R$  given the (x,y)- coordinates of the parent points. This function is provided in **GET** to support [hotspots.MatClustlpp](#). See the hotspots vignette available by starting R, typing `library("GET")` and `vignette("GET")`.

**Usage**

```
rMatClustlpp(Centers, R, alpha, LL, check_vol = FALSE)
```

**Arguments**

Centers	The (x,y)-coordinates of parent points
R	Cluster radius parameter of the Matern cluster process.
alpha	Parameter related to mean number of points per cluster.
LL	The linear network on which the point pattern should be simulated.
check_vol	Logical. TRUE for checking if the ball producing the cluster has any intersection with linear network.

---

`roadcrash`*Road crashes*

---

**Description**

Road crashes

**Usage**`data("roadcrash")`**Format**

A list with

- `x` = x-coordinates of road accidents
- `y` = y-coordinates of road accidents
- `xrange` = x coordinate limits of enclosing box (-774936.86,-727048.86)
- `yrange` = y coordinate limits of enclosing box (-1201599.83,-1125679.83)
- `Vertices.x` = x-coordinates of vertices of the linear network
- `Vertices.y` = y-coordinates of vertices of the linear network
- `Edges` = a 2 column matrix giving the ID (index) of the origin and destination vertices (in `Vertices.x` and `Vertices.y`)
- `Traffic` = matrix of traffic volume
- `ForestDensity` = matrix of forest density
- `BuildingDensity` = matrix of building density

**Details**

Mrkvička et al. (2023) worked with the database of road crashes reported to the Police in the Czech Republic from 1 January 2016 to 31 December 2020. The data available here is a subpattern of this full data set, included here with the permission of the Police in the Czech Republic. The full data is published as open data, see <https://policyvisuals.eu/traffic-accidents-data-in-the-czech-republic/>. The subpattern 7700 crashes lying on a linear network with 269 vertices and 354 lines. Further average traffic volume (number of vehicles per 24 hours), forest density and building density in the cell are available in the region of the linear network.

**References**

Mrkvička, T., Kraft, S., Blažek, V. and Myllymäki, M. (2023) Hotspots detection on a linear network with presence of covariates: a case study on road crash data.

**Examples**

```

if(require("spatstat.geom", quietly = TRUE) & require("spatstat.linnet", quietly = TRUE)) {
  data("roadcrash")
  win <- owin(xrange = roadcrash$xrange,
             yrange = roadcrash$yrange)
  X <- ppp(x = roadcrash$x, y = roadcrash$y, window = win)
  Vertices.pp <- ppp(x = roadcrash$Vertices.x,
                    y = roadcrash$Vertices.y,
                    window=win)
  L <- linnet(vertices=Vertices.pp,
             edges = roadcrash$Edges)
  PP <- lpp(X, L)
  z1 <- im(roadcrash$Traffic,
          xrange = roadcrash$xrange,
          yrange = roadcrash$yrange)
  z2 <- im(roadcrash$ForestDensity,
          xrange = roadcrash$xrange,
          yrange = roadcrash$yrange)
  z3 <- im(roadcrash$BuildingDensity,
          xrange = roadcrash$xrange,
          yrange = roadcrash$yrange)
}

```

---

saplings

*Saplings data set*


---

**Description**

Saplings data set

**Usage**

```
data("saplings")
```

**Format**

A data.frame containing the locations (x- and y-coordinates) of 123 trees in an area of 75 m x 75 m.

**Details**

A pattern of small trees (height  $\leq 15$  m) originating from an uneven aged multi-species broadleaf nonmanaged forest in Kaluzhskie Zaseki, Russia.

The pattern is a sample part of data collected over 10 ha plot as a part of a research program headed by project leader Prof. O.V. Smirnova.

## References

- Grabarnik, P. and Chiu, S. N. (2002) Goodness-of-fit test for complete spatial randomness against mixtures of regular and clustered spatial point processes. *Biometrika*, **89**, 411–421.
- van Lieshout, M.-C. (2010) Spatial point process theory. In Handbook of Spatial Statistics (eds. A. E. Gelfand, P. J. Diggle, M. Fuentes and P. Guttorp), Handbooks of Modern Statistical Methods. Boca Raton: CRC Press.
- Myllymäki, M., Mrkvička, T., Grabarnik, P., Seijo, H. and Hahn, U. (2017). Global envelope tests for spatial point patterns. *Journal of the Royal Statistical Society: Series B (Statistical Methodology)*, 79: 381-404. doi: 10.1111/rssb.12172

## See Also

[adult\\_trees](#)

## Examples

```
# This is an example analysis of the saplings data set
#=====
# Example of Myllymaki et al. (2017, Supplement S4).
if(require("spatstat.explore", quietly=TRUE)) {
  data("saplings")
  saplings <- as.ppp(saplings, W=square(75))

  # First choose the r-distances for L (r) and J (rJ) functions, respectively.
  nr <- 500
  rmin <- 0.3; rminJ <- 0.3
  rmax <- 10; rmaxJ <- 6
  rstep <- (rmax-rmin)/nr; rstepJ <- (rmaxJ-rminJ)/nr
  r <- seq(0, rmax, by=rstep)
  rJ <- seq(0, rmaxJ, by=rstepJ)

  #-- CSR test --# (a simple hypothesis)
  #-----#
  # First, a CSR test using the L(r)-r function:
  # Note: CSR is simulated by fixing the number of points and generating nsim simulations
  # from the binomial process, i.e. we deal with a simple hypothesis.
  nsim <- 999 # Number of simulations

  env <- envelope(saplings, nsim=nsim,
    simulate=expression(runifpoint(ex=saplings)), # Simulate CSR
    fun="Lest", correction="translate", # T(r) = estimator of L with translational edge correction
    transform=expression(.-r),          # Take the L(r)-r function instead of L(r)
    r=r,                                # Specify the distance vector
    savefuns=TRUE)                      # Save the estimated functions
  # Crop the curves to the interval of distances [rmin, rmax]
  # (at the same time create a curve_set from 'env')
  curve_set <- crop_curves(env, r_min=rmin, r_max=rmax)
  # Perform a global envelope test
  res <- global_envelope_test(curve_set, type="erl") # type="rank" and larger nsim was used in S4.
  # Plot the result.
  plot(res) + ggplot2::ylab(expression(italic(hat(L)(r)-r)))
```

```

# -> The CSR hypothesis is clearly rejected and the rank envelope indicates clear
# clustering of saplings. Next we explore the Matern cluster process as a null model.
}

if(require("spatstat.model", quietly=TRUE)) {
  #-- Testing the Matern cluster process --# (a composite hypothesis)
  #-----#
  # Fit the Matern cluster process to the pattern (using minimum contrast estimation with the pair
  # correction function)
  fitted_model <- kppm(saplings~1, clusters="MatClust", statistic="pcf")
  summary(fitted_model)

  nsim <- 19 # 19 just for experimenting with the code!!
  #nsim <- 499 # 499 is ok for type = 'qdir' (takes > 1 h)

  # Make the adjusted directional quantile global envelope test using the L(r)-r function
  # (For the rank envelope test, choose type = "rank" instead and increase nsim.)
  adjenvL <- GET.composite(X=fitted_model,
    fun="Lest", correction="translate",
    transform=expression(.-r), r=r,
    type="qdir", nsim=nsim, nsimsub=nsim,
    r_min=rmin, r_max=rmax)

  # Plot the test result
  plot(adjenvL) + ggplot2::ylab(expression(italic(L(r)-r)))

  # From the test with the L(r)-r function, it appears that the Matern cluster model would be
  # a reasonable model for the saplings pattern.
  # To further explore the goodness-of-fit of the Matern cluster process, test the
  # model with the J function:
  # This takes quite some time if nsim is reasonably large.
  adjenvJ <- GET.composite(X=fitted_model,
    fun="Jest", correction="none", r=rJ,
    type="qdir", nsim=nsim, nsimsub=nsim,
    r_min=rminJ, r_max=rmaxJ)

  # Plot the test result
  plot(adjenvJ) + ggplot2::ylab(expression(italic(J(r))))
  # -> the Matern cluster process not adequate for the saplings data

  # Test with the two test functions jointly
  adjenvLJ <- GET.composite(X=fitted_model,
    testfuns=list(L=list(fun="Lest", correction="translate",
      transform=expression(.-r), r=r),
      J=list(fun="Jest", correction="none", r=rJ)),
    type="erl", nsim=nsim, nsimsub=nsim,
    r_min=c(rmin, rminJ), r_max=c(rmax, rmaxJ),
    save.cons.envelope=TRUE)

  plot(adjenvLJ)
}

```



**Description**

ggproto objects for central region plot. Not to be used directly.

**Usage**

StatCentralRegion

GeomCentralRegion

GeomCentralRegionMulti

**Format**

An object of class StatCentralRegion (inherits from Stat, ggproto, gg) of length 4.

An object of class GeomCentralRegion (inherits from Geom, ggproto, gg) of length 4.

An object of class GeomCentralRegionMulti (inherits from Geom, ggproto, gg) of length 4.

**See Also**

[geom\\_central\\_region](#)

---

subset.curve_set	<i>Subsetting curve sets</i>
------------------	------------------------------

---

**Description**

Return subsets of curve sets which meet conditions.

**Usage**

```
## S3 method for class 'curve_set'
subset(x, subset, ...)
```

**Arguments**

x	A curve_set object.
subset	A logical expression indicating curves to keep.
...	Ignored.

**Examples**

```
if(require("fda.usc", quietly=TRUE)) {  
  # Prepare data  
  data("poblenou")  
  Free <- poblenou$df$day.festive == 1 |  
    as.integer(poblenou$df$day.week) >= 6  
  MonThu <- poblenou$df$day.festive == 0 & poblenou$df$day.week %in% 1:4  
  Friday <- poblenou$df$day.festive == 0 & poblenou$df$day.week == 5  
  
  # Data as a curve_set  
  cset <- curve_set(r=0:23, obs=t(poblenou[['nox']][['data']]))  
  plot(subset(cset, MonThu))  
  plot(subset(cset, Friday))  
  plot(subset(cset, Free))  
}
```

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