

Package ‘briskaR’

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Description A spatio-temporal exposure-hazard model for assessing biological risk and impact. The model is based on stochastic geometry for describing the landscape and the exposed individuals, a dispersal kernel for the dissemination of contaminants, a set of tools to handle spatio-temporal dataframe and ecotoxicological equations.

Walker E, Leclerc M, Rey JF, Beaudouin R, Soubeyrand S, and Messean A, (2017),
A Spatio-Temporal Exposure-Hazard Model for Assessing Biological Risk and Impact,
Risk Analysis, <doi:10.1111/risa.12941>.

Leclerc M, Walker E, Messean A, Soubeyrand S (2018),
Spatial exposure-hazard and landscape models for assessing the impact of GM crops on non-target organisms,
Science of the Total Environment, 624, 470-479.

URL <https://gitlab.paca.inrae.fr/biosp/briskaR>

BugReports <https://gitlab.paca.inrae.fr/biosp/briskaR/-/issues>

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| <i>briskaR-package</i> | <i>Biological Risk Assessment</i> |
|------------------------|-----------------------------------|

Description

A spatio-temporal exposure-hazard model for assessing biological risk and impact. The model is based on stochastic geometry for describing the landscape and the exposed individuals, a dispersal kernel for the dissemination of contaminants and an ecotoxicological equation.

Details

The *briskaR* package contains functions and methods for quantifying spatio-temporal variation in contamination risk around known polygon sources of contaminants, and quantifies the impact of the contaminants on the surrounding population of individuals which are located in habitat areas and are susceptible to the contaminants.

The package implements an spatio-temporal exposure-hazard model based on (i) tools of stochastic geometry (marked polygon and point processes) for structuring the landscape and describing the location of exposed individuals, (ii) a method based on a dispersal kernel describing the dissemination of contaminant particles from polygon sources, and (iii) ecotoxicological equations describing how contaminants affect individuals of the exposed population.

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See Also

Useful links:

- <https://gitlab.paca.inrae.fr/biosp/briskaR>
- Report bugs at <https://gitlab.paca.inrae.fr/biosp/briskaR/-/issues>

```
briskaRLoadInternProjection
```

Load an internal working projection PROJ.4

Description

Will load a projection as internal package working projection

Usage

```
briskaRSetInternProjection(proj = LAMBERT_93)
```

Arguments

| | |
|------|---|
| proj | A character string of projection arguments, must be in the PROJ.4 documentation |
|------|---|

```
brk_addFD
```

Functional DATA

Description

Add time series to sf objects

Usage

```
brk_addFD(sf, key, FUN, ...)
```

```
brk_addFD2(sf, keyConstraint, key, FUN, ...)
```

Arguments

| | |
|---------------|---|
| sf | object of class sf . See sf for details. |
| key | the name of the new column, as strings or symbols |
| FUN | the function to be applied to each element of sf . In the case of functions like <code>+</code> , backquoted or quoted. See lapply functions for details. |
| ... | optional arguments to FUN. See lapply functions for details. |
| keyConstraint | character string. The reference of the column to be checked |

Value

A [sf](#) object with addition functional data feature (or feature dynamic).

| | |
|---------------|---|
| brk_cFilterFD | <i>Function used to filter functional data in sf objects.</i> |
|---------------|---|

Description

Function used to filter functional data in [sf](#) objects.

Usage

```
brk_cFilterFD(sf, key, index)
brk_cFilterFD_(sf, key, index)
brk_cFilterFD2(sf, key1, key2, index)
brk_cFilterFD3(sf, key1, key2, key3, index)
```

Arguments

| | |
|-------|--|
| sf | sf. An object of class sf |
| key | character string. The name of the column to select |
| index | integer (or vector). The index of the functional data. |
| key1 | character string. The name of the column to select |
| key2 | character string. The name of the column to select |
| key3 | character string. The name of the column to select |

| | |
|---------------|--------------------------|
| brk_dispersal | <i>Compute dispersal</i> |
|---------------|--------------------------|

Description

Simulate contaminants or individuals frequency over the landscape by two steps: dispersal of elements and local intensity/frequency of elements after dispersal.

Usage

```
brk_dispersal(
  object,
  size_raster,
  tolerance_square,
  kernel,
  kernel.options,
  nbr_cores,
  squared_frame
)
```

Arguments

| | |
|------------------|--|
| object | sf or patialPolygonsDataFrame. A simple feature of class sf or SpatialPolygons-DataFrame |
| size_raster | integer. Raster size (default = 2^10) |
| tolerance_square | numeric. Tolerance rate to test if an sf set is squared |
| kernel | string. Dispersion kernel, function name (default = NIG) |
| kernel.options | list. Parameters list for the kernel function |
| nbr_cores | integer. Parameters for parallel computing: the number of cores to use, i.e. at most how many child processes will be run simultaneously. Default is 1 (non parallel). |
| squared_frame | sf. Select the sf to be considered as frame to rasterized. Default is 'NULL', and 'object' is used. |

Details

The dispersal of contaminants or individuals is implemented by rastering the landscape and by computing the convolution between sources emissions and a dispersal kernel.

The dispersion kernel by default is Normal Inverse Gaussian kernel ("NIG" function). Currently, two others are implemented "geometric" (with parameter a) and "2Dt" kernels (with parameters a, b, c1, c2).

| | |
|--------------|--------------------------------------|
| brk_emission | <i>Wrapper function brk_emission</i> |
|--------------|--------------------------------------|

Description

'brk_emission' This function simulates emissions. Will simulate emissions shape in source fields of a landscape.

'brk_emission_landscape' Simulate sources emission intensity. With: - 'emission' is the quantity of emission per time unit per spatial unit. With other argument, we have: - 'emission = density x intensity = density x production x intensity_pmf'. - 'density' is the density of the source, so the quantity of source per spatial unit. - 'intensity' is the intensity of the emission for 1 source: that is the quantity of emission per time unit per source. - 'production' is the overall production, the total emission, for one source unit: quantity of emission for the period - 'intensity_pmf' is the distribution of emission along time. So we have 'intensity = production x intensity_pmf'.

Usage

```
brk_emission(sf, keyTime, key, FUN)
```

```
brk_emission_landscape(
  sf,
  timeline = 1:61,
```

```

    emission = NULL,
    density = NULL,
    intensity = NULL,
    intensity_pmf = NULL,
    production = NULL
  )

```

Arguments

| | |
|----------------------------|---|
| <code>sf</code> | A <code>sf</code> object |
| <code>keyTime</code> | character. Name of the column for timeline. Used to check length of vectors in column vector object |
| <code>key</code> | character. Name of the column which is going to be created |
| <code>FUN</code> | A function applied on the <code>sf</code> object. See <code>lapply</code> functions for further details assuming $X = 1:nrow(sf)$, that is <code>sf[[key]] <- lapply(1:nrow(sf), FUN, ...)</code> . |
| <code>timeline</code> | Vector of time units (e.g. days) covering all the function |
| <code>emission</code> | Vector or Matrix given the quantity of emission per time unit per spatial unit. Length of vector equal the length of the 'timeline' vector (time unit matching). Size of the matrix, 'n,m', is such as the number of row equal the number of sources in 'sf' object, and the number of column equals the length of the 'timeline' vector. |
| <code>density</code> | Scalar or Vector (with length equal to the number of sources in 'sf' object) given the density of the source(s) (e.g. number of plant by squared meter) |
| <code>intensity</code> | Vector or Matrix given the quantity of emission per time unit per source. Length of vector equal the length of the 'timeline' vector (time unit matching). Size of the matrix, 'n,m', is such as the number of row equal the number of sources in 'sf' object, and the number of column equals the length of the 'timeline'. |
| <code>intensity_pmf</code> | Vector or Matrix given distribution of emission along time (given a probability mass function with time) . Length of vector equal the length of the 'timeline' vector (time unit matching). Size of the matrix, 'n,m', is such as the number of row equal the number of sources in 'sf' object, and the number of column equals the length of the 'timeline'. |
| <code>production</code> | Scalar or Vector (with length equal to the number of sources in 'sf' object) total emission for one source (e.g. total number of grains by plant) |

Details

This function is a wrapper of `with` and `lapply` function and is like this: `sf[["key"]] <- with(sf, lapply(1:nrow(sf), FUN, ...))`

So, all column of `sf` can be called in `FUN`

Value

A matrix indexed by sources ID (in rows) and by time (in columns) whose rows give the values of intensity emission (number of grains) for every source.

| | |
|--------------|--|
| brk_exposure | <i>Compute exposure for a 'RasterStack' class object from package 'raster'</i> |
|--------------|--|

Description

Compute exposure for a 'RasterStack' class object from package 'raster'

Usage

```
brk_exposure(
  RasterStack_dispersal,
  sf,
  key,
  keyTime,
  loss,
  beta,
  nbr_cores,
  quiet,
  unit
)
```

Arguments

| | |
|-----------------------|---|
| RasterStack_dispersal | RasterStack. An object of class RasterStack (see package raster for details). |
| sf | sf. And object of class 'sf' on which exposure is computed from the previous list of raster by patch 'RasterStack_dispersal'. See sf for details. |
| key | name of the column in 'sf' object providing emission amount |
| keyTime | name of the column of sf for time |
| loss | numeric. scalar or vector (of the same length as the number the timeline include is argument sf) to apply a loss on exposure cells. |
| beta | numeric. toxic adherence parameter between 0 and 1 (default = 0.4). |
| nbr_cores | integer. Set the number of cores to used for parallel computing. |
| quiet | boolean. Set 'TRUE' to remove progress bar. |
| unit | default is meter "m". but should be more generic: "any". |

Details

Local intensity depends of beta and alpha parameters. Beta represents the toxic adherence between [0,1]. Alpha represents a list of parameters of the lost of toxic particules due to covariates (precipitation). There are two configurations to integrate the loss in the function : (i) simulating covariate (simulate=TRUE) or (ii) uploading covariate (simulate=FALSE). The covariate is linked to the loss by a linear regression with paramaters minalpha, maxalpha, covariate_threshold.

brk_exposureMatch *Add raster value to element of sf object*

Description

Add the raster value(s) of a Raster* object to element of sf object.

Usage

```
brk_exposureMatch(
  stackRaster_exposure,
  sf,
  stackTimeline,
  keyTime = "TIMELINE",
  key = "EXPOSURE"
)
```

Arguments

| | |
|----------------------|--|
| stackRaster_exposure | The Raster* object |
| sf | the sf object |
| stackTimeline | sequence with matching to elements of RasterStack (length of stackTimeline must be the same as length of a list of stackRaster_exposure) |
| keyTime | name of the column to match exposure timeline from stackRaster_exposure |
| key | name of the new column |

brk_FDtoDF *Convert list.column data.frame into scalar.column data.frame*

Description

Convert data.frame with 2 column.list into data.frame with only column.scalar

Usage

```
brk_FDtoDF_(sf, key1, key2, id = NULL)

brk_FDtoDF(sf, key1, key2, id = NULL, keep = NULL)

brk_FDtoDF_STICK(sf, key1, keep)
```

Arguments

| | |
|------|---|
| sf | sf. An object of class sf |
| key1 | character string. The name of the column to select |
| key2 | character string. The name of the column to select |
| id | name of the replicate for the id. As to be of the same length as the number of row of the sf object |
| keep | vector of column name to keep |

| | |
|-----------------|-------------------|
| brk_findIndexFD | <i>find index</i> |
|-----------------|-------------------|

Description

find index

Usage

```
brk_findIndexFD(sf, key, value)
```

Arguments

| | |
|-------|---|
| sf | An object of class sf |
| key | character string. name of the column to select |
| value | value of the element to return index from the column defined by key |

Value

vector if not all index are equal. scalar if all equal.

| | |
|---------------|--|
| brk_newPoints | <i>Simulate new points on a specific sf object</i> |
|---------------|--|

Description

Simulate new points on a specific [sf](#) object. See [st_sample](#) for details.

Usage

```
brk_newPoints(sf, size = 100)
```

Arguments

| | |
|------|--|
| sf | object of class sf or sfc |
| size | sample size(s) requested; either total size, or a numeric vector with sample sizes for each feature geometry. When sampling polygons, the returned sampling size may differ from the requested size, as the bounding box is sampled, and sampled points intersecting the polygon are returned. |

| | |
|-----------------|---|
| brk_rbindLStoDF | <i>Combine list of data.frame by Rows</i> |
|-----------------|---|

Description

Function used to filter functional data in [sf](#) objects.

Usage

```
brk_rbindLStoDF(ls, id = NULL)
```

Arguments

| | |
|----|--|
| ls | A list of data.frame |
| id | id to provide to each data.frame. Must be the length of the list |

Value

Return a data.frame

| | |
|--------------|---|
| brk_sampling | <i>Sampling based on distribution provided by Raster* object over a sf object</i> |
|--------------|---|

Description

Sampling based on distribution provided by Raster* object (does not have to be summed to one) over an sf object used as mask

Usage

```
brk_sampling(rasterStack, sf, sizeSite = 1)
```

Arguments

| | |
|-------------|---|
| rasterStack | Raster* object (RasterLayer or RasterStack) |
| sf | An sf object |
| sizeSite | scalar or vector of the number of individual to sample per site. Default is 1. If vector, as to be the length of Raster* object |

| | |
|--------------|---|
| brk_timeline | <i>Add sequence to element of a sf object</i> |
|--------------|---|

Description

Add column with sequence for each row to a [sf](#) object

Usage

```
brk_timeline(sf, key, from, to, by)
```

Arguments

| | |
|------|---|
| sf | object of class sf or sfc |
| key | a character string used as name of the new column |
| from | the starting value of the sequence. Of length 1 unless just from is supplied as an unnamed argument. |
| to | the end (maximal) value of the sequence. Of length 1 unless just from is supplied as an unnamed argument. |
| by | number: increment of the sequence. |

| | |
|------------|--|
| brk_toxFun | <i>Functions for Toxicokinetic-Toxidynamic (TKTD) models</i> |
|------------|--|

Description

Functions for Toxicokinetic-Toxidynamic (TKTD) models

ODE solver applied to IT-GUTS model

ODE solver applied to SD-GUTS model

Usage

```
brk_toxFun_damage1(exposure, kin, kout)
```

```
brk_toxFun_survival1(damage, alpha1, alpha2, alpha3)
```

```
brk_toxFun_survival2(damage, LC50, slope)
```

```
brk_survIT(time, Cw, listParameters)
```

```
brk_survSD(time, Cw, listParameters)
```

Arguments

| | |
|----------------|--|
| exposure | exposure level of individual |
| kin | parameter describing the intake rate of the element |
| kout | parameter describing the excretion rate of the element |
| damage | damage level on which the survival model is going to be applied |
| alpha1 | parameter describing the natural background death mortality rate |
| alpha2 | parameter describing the killing rate of the element on individual |
| alpha3 | exponential parameter describing the killing rate of the element on individual |
| LC50 | parameter describing the lethal concentration for 50% of the population |
| slope | parameter describing the slope of the curve |
| time | vector of time of the exposure profile |
| Cw | vector of concentration of the exposure profile |
| listParameters | A list of parameter for the SD model |

create_pollen_sources *Pollen sources emission simulation*

Description

Simulate pollen sources emission for maize crop. The proportion of plants emitting pollen per day (during 12 days) was observed by Frédérique Angevin (Angevin et al. 2008).

Usage

```
create_pollen_sources(
  sf,
  timeline = 1:61,
  density = runif(1, 7, 11),
  pollen = rgamma(1, shape = 1.6, scale = 1/(2 * 10^-7))
)

create.pollen.sources()
```

Arguments

| | |
|----------|---|
| sf | A sf object defining sources fields |
| timeline | Vector of interger. Time units (e.g. days) including the pollen emission (for simulation, default is time.max). Minimal value is a vector of 12 days. |
| density | Plant density (number of plant by squared meter) |
| pollen | Pollen production (number of grains by plant) |

Value

A matrix indexed by sources ID (in rows) and by time (in columns) whose rows give the values of pollen emission (number of grains) for every source.

| | |
|----------|---|
| data_brk | <i>Data set included in the package</i> |
|----------|---|

Description

Data set included in the package

Usage

```
data(maize_65)
data(sfMaize65)
data(maize.emitted_pollen)
data(maize.proportion_pollen)
data(Hofmann_2009)
data(Lang_2004)
data(Precipitation)
data(temperatureGermany)
data(df_precipitation)
```

Arguments

| | |
|-------------------------|--|
| maize_65 | A SpatialPolygonsDataFrame of class SpatialPolygonsDataFrame defining a patchy landscape |
| sfMaize65 | A set of MULTIPOLYGON of class sf defining a patchy landscape |
| maize.emitted_pollen | A data.frame of pollen emission |
| maize.proportion_pollen | A data.frame of pollen proportion |
| Hofmann_2009 | A data.frame of pollen emission |
| Lang_2004 | A data.frame of pollen emission |
| Precipitation | data.frame of daily Precipitation from 01/01/2003 to 31/12/2013 |
| temperatureGermany | A data.frame of temperature over a year in south and north of germany |
| df_precipitation | A data.frame of daily Precipitation from 01/01/2003 to 31/12/2013 |

| | |
|---------------------|---|
| GetInternProjection | <i>Get the internal working projection PROJ.4</i> |
|---------------------|---|

Description

Will print and return the internal projection of briskaR package

Usage

```
briskaRGetInternProjection()
```

| | |
|--------------|--|
| is_square_sf | <i>Test if an sf is a square</i> |
|--------------|--|

Description

Test if an sf is a square, with tolerance. Default is 5

Usage

```
is_square_sf(.sf, tolerance = 0.05)
```

Arguments

| | |
|-----------|--|
| .sf | and object of class sf |
| tolerance | tolerance rate between both square side length |

| | |
|------------|-------------------|
| LAMBERT_93 | <i>LAMBERT_93</i> |
|------------|-------------------|

Description

SIG projection Lambert_93 references "+init=epsg:2154" PROJ.4

Usage

```
LAMBERT_93
```

Format

An object of class character of length 1.

| | |
|-----------------|---|
| loadIndividuals | <i>Wrapper function loadIndividuals</i> |
|-----------------|---|

Description

Wrapper function to create an individuals object using [sf](#) or [SpatialPoints](#) and data.frame.

The [SpatialPoints](#) object and the data.frame have to contain the same number of coordinates and rows.

Usage

```
loadIndividuals(sf, data, timeline)
```

Arguments

| | |
|-----------------------|--|
| <code>sf</code> | a sf object |
| <code>data</code> | a data.frame containing individuals attributes. Rows numbers as individuals ID, columns names as dob (date of birth) life_duration toxic_threshold |
| <code>timeline</code> | Vector of the time line |

Value

an Individuals object

| | |
|---------------|---|
| loadLandscape | <i>Wrapper function : loadLandscape</i> |
|---------------|---|

Description

Wrapper function to create a Landscape object using [SpatialPolygons](#) and dataframe. The [SpatialPolygons](#) object and the data.frame have to contain the same number of polygons and row (row ID is polygons ID).

Usage

```
loadLandscape(sp, data)
```

Arguments

| | |
|-------------------|--|
| <code>sp</code> | a SpatialPolygons object designing the landscape |
| <code>data</code> | a data.frame containing fields (polygons) information. Row names as fields ID, column names as sources neutral receptors (for a given field, the value is 1 for the type of the field (source or neutral or receptor), otherwise 0). |

Value

A [SpatialPolygonsDataFrame](#) object

Examples

```
data(maize_65)
maize_data <- maize_65@data
maize_sp_only <- maize_65 ; maize_sp_only@data = data.frame(remove = rep(0,nrow(maize_65@data)))
load_landscape <- loadLandscape(maize_sp_only, maize_data)
```

loadLandscapeSIG *Create a Landscape object from SIG shapefile file*

Description

Create a Landscape object from SIG shapefile. Shapefile has to contain a [SpatialPolygonsDataFrame](#). Data in the data frame contain fields (polygons) information. Row names as fields ID, cols names as sources | neutral | receptors (for a given field, the value is 1 for the type of the field (source or neutral or receptor), otherwise 0).

Usage

```
loadLandscapeSIG(dsn, layer, format = TRUE)
```

Arguments

| | |
|--------|---|
| dsn | folder path to the source files |
| layer | file name without extension |
| format | only load data needed Landscapae-class (default TRUE) |

Value

A [SpatialPolygonsDataFrame](#) object

Examples

```
## Not run:
land <- loadLandscapeSIG("/path/to/directory/", "fileName")
plot(land)

## End(Not run)
```

| | |
|--------------------|---|
| loss_precipitation | <i>Loss precipitation function from a set of function implemented in earlier version of briskaR</i> |
|--------------------|---|

Description

Loss precipitation function from a set of function implemented in earlier version of briskaR

Usage

```
loss_precipitation(
  starttime = "16/07",
  endtime = "14/07",
  alpha = list(minalpha = 0.1, maxalpha = 0.95, covariate_threshold = 30, simulate =
    TRUE, covariate = NULL)
)
```

Arguments

| | |
|-----------|--------------------------------|
| starttime | data of begining of simulation |
| endtime | data of end of simulation |
| alpha | list of parameter |

| | |
|--------------|--|
| saveIntoFile | <i>Save Particles Dispersion 3D Array to tiff file</i> |
|--------------|--|

Description

Save into tiff file particles dispersion 3D array from toxicIntensity. The output is a RasterStack with a layer per time unit with projection set to CRS="+proj=longlat +datum=WGS84"

Usage

```
saveIntoFile(
  objectL,
  objectT,
  filename = "ParticlesDispersion.tif",
  format = "GTiff",
  overwrite = TRUE
)
```

Arguments

| | |
|-----------|---|
| objectL | a Landscape object |
| objectT | a 3D array particles dispersion indexed by time (output from toxicIntensity) |
| filename | output file name (default "ParticlesDispersion.tif") |
| format | output format (default=GTiff) |
| overwrite | if TRUE overwrite file (default TRUE) |

Value

a RasterStack object

Examples

```
## Not run:  
data(maize_65)  
ti <- toxicIntensity(maize.landscape,maize.emitted_pollen)  
saveIntoFile(maize.landscape,ti,filename="ParticlesDispersion.tiff",format="GTiff",overwrite=T)  
  
## End(Not run)
```

simul.precipitation *Simulate precipitation between two dates*

Description

Will evaluate parameters from data and simulate precipitation between the two dates.

Usage

```
simul.precipitation(starttime = "15/07", endtime = "15/09", data = NULL)
```

Arguments

| | |
|-----------|-------------------------------|
| starttime | *string*. Date shape: "mm/yy" |
| endtime | *string*. Date shape: "mm/yy" |
| data | *data.frame*. default is NULL |

Value

an array of length of the period between the two dates included

Author(s)

Jean-Francois Rey

simulateIndividuals *Wrapper function SimulateIndividuals*

Description

This function simulates individuals as an Individuals object.

Will simulate n individuals in receptors fields of a landscape.

Usage

```
simulateIndividuals(  
  sf,  
  size = 100,  
  timeline = 1:61,  
  dob,  
  life_duration,  
  toxic_threshold  
)
```

Arguments

| | |
|-----------------|--|
| sf | A sf object |
| size | Number of individuals to simulate |
| timeline | Vector of the time line |
| dob | A vector for the Date Of Birth of each individual between min and max of the timeline. |
| life_duration | A vector for the life duration of each individual |
| toxic_threshold | A vector for the internal toxic threshold value leading to death for each individual |

Details

The Individuals object output includes for each individual the coordinates, the date of birth, the life duration, the toxic threshold and

`simulateInitialPartition`*simulateInitialPartition Method*

Description

This function creates an object [SpatialPolygonsDataFrame](#) and simulates a landscape with neutral and source fields.

Usage

```
simulateInitialPartition(  
  n = 500,  
  prop = 0.4,  
  range = 10,  
  xmin = 0,  
  xmax = 5000,  
  ymin = 0,  
  ymax = 5000  
)
```

Arguments

| | |
|--------------------|---|
| <code>n</code> | Numeric, numbers of cells |
| <code>prop</code> | Numeric [0,1] toxic cells proportion |
| <code>range</code> | Aggregation parameter (range of the spatial Exponential covariance of Gaussian process) |
| <code>xmin</code> | x-axis left coordinates in space unit (see <code>projections_briskaR</code>) |
| <code>xmax</code> | x-axis right coordinates in space unit (see <code>projections_briskaR</code>) |
| <code>ymin</code> | y-axis bottom coordinates in space unit (see <code>projections_briskaR</code>) |
| <code>ymax</code> | y-axis top coordinates in space unit (see <code>projections_briskaR</code>) |

Details

In the function the first step is a binomial point process to simulate a [SpatialPointsDataFrame](#) with sources and neutral marks, which depends on the aggregation parameter. The second step of the function is the Voronoi tessellation from the simulated points and returns a [SpatialPolygonsDataFrame](#).

Value

An [SpatialPolygonsDataFrame](#) object with `n` fields, `prop` pourcentage of toxic fields of size `(xmin,xmax)` `(ymin,ymax)`

Examples

```
## Not run:
# Simulate a 5000m x 5000m landscape with 500 cells (e.g. fields)
# whose 40% (200 cells) are sources.
# The projection by default is Lambert93 projection.
land <- simulateInitialPartition(n=500,prop=0.4,range=10,xmin=0,xmax=5000,ymin=0,ymax=5000)
plot(land)

## End(Not run)
```

| | |
|-------------------|---------------------------------|
| simulateLandscape | <i>Simulate a new landscape</i> |
|-------------------|---------------------------------|

Description

Create an object of class [SpatialPolygonsDataFrame](#). Simulate a landscape with neutral and source fields and receptors margins.

Usage

```
simulateLandscape(
  n = 500,
  prop = 0.4,
  range = 10,
  xmin = 0,
  xmax = 5000,
  ymin = 0,
  ymax = 5000,
  border_size = 200,
  prob = runif(1, 0.1, 0.9),
  mean_thickness = runif(1, 2, 20),
  v_thickness = 50
)
```

Arguments

| | |
|-------|--|
| n | Numeric, numbers of fields |
| prop | Numeric [0,1] toxic fields proportion |
| range | aggregation parameter (range of the spatial exponential covariance of gaussian process) in meters. |
| xmin | x-axis left coordinates |
| xmax | x-axis right coordinates |
| ymin | y-axis bottom coordinates |
| ymax | y-axis top coordinates |

| | |
|----------------|---------------------------------------|
| border_size | A numeric, bbox margin |
| prob | Probability to inflate a filed margin |
| mean_thickness | Margin width expectation |
| v_thickness | Margin width variance |

Details

Execute both [simulateInitialPartition](#) and [simulateThickMargins](#) functions.

Value

A [SpatialPolygonsDataFrame](#) object with n fields, prop pourcentage of toxic fields.

See Also

[simulateInitialPartition](#) and [simulateThickMargins](#)

Examples

```
## Not run:  
land <- simulateLandscape(n=100, prop=0.4, range=10,  
xmin=0, xmax=1000, ymin=0, ymax=1000, border_size=100,  
prob=runif(1,0.1,0.9), mean_thickness=runif(1,2,20),  
v_thickness=50)  
plot(land)  
## End(Not run)
```

simulateThickMargins *Simulate thick margin to a landscape*

Description

Simulate thick margins as receptors in a landscape.

Usage

```
simulateThickMargins(  
  objectL,  
  border_size = 200,  
  prob = runif(1, 0.1, 0.9),  
  mean_thickness = runif(1, 2, 20),  
  v_thickness = 50  
)
```

Arguments

| | |
|----------------|---|
| objectL | sf, sp or Landscape (earlier version of briskaR). |
| border_size | A numeric, bbox margin |
| prob | Probability to inflate a margin |
| mean_thickness | Margin width expectation in meter |
| v_thickness | Margin width variance in meter |

Details

Margin width use a Gamma distribution with shape and scale parameters based on thickness mean and variance.

Value

A [SpatialPolygonsDataFrame](#) object

See Also

[simulateInitialPartition](#) and [simulateLandscape](#)

Examples

```
## Not run:  
data(maize_65)  
plot(maize.landscape)  
landscape.margin <- simulateThickMargins(maize.landscape)  
plot(landscape.margin)  
## End(Not run)
```

st_multibuffer

Simulate thick margin to a landscape

Description

Add buffer around each objects of a [sf](#) file

Usage

```
st_multibuffer(  
  sf,  
  dist = 50,  
  nQuadSegs = 30,  
  endCapStyle = "ROUND",  
  joinStyle = "ROUND",  
  mitreLimit = 1  
)
```


Arguments

| | |
|--------------------------|---|
| <code>sf</code> | object of class <code>sfg</code> , <code>sfg</code> or <code>sf</code> |
| <code>dist</code> | numeric; buffer distance for all, or for each of the elements in <code>x</code> ; in case <code>dist</code> is a units object, it should be convertible to <code>arc_degree</code> if <code>x</code> has geographic coordinates, and to <code>st_crs(x)\$units</code> otherwise |
| <code>nQuadSegs</code> | integer; number of segments per quadrant (fourth of a circle), for all or per-feature |
| <code>endCapStyle</code> | character; style of line ends, one of 'ROUND', 'FLAT', 'SQUARE' |
| <code>joinStyle</code> | character; style of line joins, one of 'ROUND', 'MITRE', 'BEVEL' |
| <code>mitreLimit</code> | numeric; limit of extension for a join if <code>joinStyle</code> 'MITRE' is used (default 1.0, minimum 0.0) |

Details

see package [st_buffer](#) for details

`st_squared_geometry` *Add squared frame polygon*

Description

Return a square frame surrounding a list of `sf`

Usage

```
st_squared_geometry(list_sf, buffer_dist = NULL)
```

Arguments

| | |
|--------------------------|---|
| <code>list_sf</code> | A list of objects of class <code>sf</code> |
| <code>buffer_dist</code> | numeric; buffer distance for all, or for each of the elements in <code>x</code> ; in case <code>dist</code> is a units object, it should be convertible to <code>arc_degree</code> if <code>x</code> has geographic coordinates, and to <code>st_crs(x)\$units</code> otherwise. See function 'sf_buffer' from package 'sf' for details |

toxicIntensity *toxicIntensity function wrapping dispersal and exposure*

Description

toxicIntensity function wrapping dispersal and exposure

Usage

```
toxicIntensity(
  object,
  sf,
  size_raster = 2^10,
  tolerance_square = 0.1,
  kernel = "NIG",
  kernel.options = list(a1 = 0.2073, a2 = 0.2073, b1 = 0.3971, b2 = 0.3971, b3 =
    0.0649, theta = 0),
  loss = NULL,
  beta = 0.4,
  nbr_cores = 1,
  squared_frame = NULL,
  quiet = FALSE
)
```

Arguments

| | |
|------------------|--|
| object | sf or SpatialPolygonsDataFrame. A simple feature of class sf or SpatialPolygonsDataFrame |
| sf | sf. And object of class 'sf' on which exposure is computed from the previous list of raster by patch 'RasterStack_dispersal'. See sf for details. |
| size_raster | integer. Raster size (default = 2^10) |
| tolerance_square | numeric. Tolerance rate to test if an sf set is squared |
| kernel | string. Dispersion kernel, function name (default = NIG) |
| kernel.options | list. Parameters list for the kernel function |
| loss | numeric. Numeric vector to applied a loss on exposure cells. |
| beta | numeric. toxic adherence parameter between 0 and 1 (default = 0.4). |
| nbr_cores | integer. Parameters for parallel computing: the number of cores to use, i.e. at most how many child processes will be run simultaneously. Default is 1 (non parallel). |
| squared_frame | sf. Select the sf to be considered as frame to rasterized. Default is 'NULL', and 'object' is used. |
| quiet | boolean. Set 'TRUE' to remove progress bar. |

Details

The dispersal of contaminants is implemented by rastering the landscape and by computing the convolution between sources emissions and a dispersal kernel.

The dispersion kernel by default is Normal Inverse Gaussian kernel ("NIG" function). Currently, two others are implemented "geometric" (with parameter a) and "2Dt" kernels (with parameters a , b , c_1 , c_2).

Local intensity depends of beta and alpha parameters. Beta represents the toxic adherence between $[0,1]$. Alpha represents a list of parameters of the loss of toxic particules due to covariates (precipitation). There are two configurations to integrate the loss in the function : (i) simulating covariate (simulate=TRUE) or (ii) uploading covariate (simulate=FALSE). The covariate is linked to the loss by a linear regression with paramaters `minalpha`, `maxalpha`, `covariate_threshold`.

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