

# Package ‘gear’

April 10, 2020

**Type** Package

**Title** Geostatistical Analysis in R

**Version** 0.3.4

**Date** 2020-04-06

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**Description** Implements common geostatistical methods in a clean, straightforward, efficient manner. The methods are discussed in Schabenberger and Gotway (2004, <ISBN:9781584883227>) and Waller and Gotway (2004, <ISBN:9780471387718>).

**License** GPL (>= 2)

**Imports** autoimage, stats, optimx, Rcpp

**Suggests** sp, sf, testthat, Matrix, geoR, gstat, spam, ggplot2, lattice

**Encoding** UTF-8

**NeedsCompilation** yes

**RoxygenNote** 7.1.0

**LinkingTo** Rcpp

**Repository** CRAN

**Date/Publication** 2020-04-10 21:20:02 UTC

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---

angle2d	<i>Determine angle</i>
---------	------------------------

---

### Description

angle2d determines the angle between pairs of coordinates in degrees or radians. The coordinates are assumed to be in two-dimensional space.

### Usage

```
angle2d(coords1, coords2, radians = FALSE, invert = TRUE)
```

### Arguments

coords1	An $N \times 2$ matrix of spatial coordinates.
coords2	An $N \times 2$ matrix of spatial coordinates.
radians	A logical value indicating whether the angles returned should be in degrees or radians. The default is FALSE, indicating that the returned angles are in degrees.
invert	A logical value indicating whether the axes of the coordinates should be inverted (i.e., the x- and y-axis are switched). The default is TRUE to mimic results from other geostatistical R packages like <code>gstat</code> , <code>geoR</code> , and other software like <code>GSLIB</code> and <code>GeoEAS</code> . Set to FALSE to use the typical x- and y-axes.

**Details**

Note that the angle is between the actual pairs of points, not the angle between the vectors extending from the origin to the points. e.g., the angle between (0, 1) and (1, 1) is 90 degrees, not 45. The sign of the direction not accounted for, e.g., a -135 degree angle is rotated by 180 degrees to become a 45 degree angle. All angles returned are in the interval [0, 180].

**Value**

Returns a vector of angles.

**Author(s)**

Joshua French

**Examples**

```
coords1 = matrix(0, nrow = 8, ncol = 2)
coords2 = cbind(c(2, 2, 0, -2, -2, -2, 0, 2), c(0, 2, 2, 2, 0, -2, -2, -2))
angle2d(coords1, coords2)
angle2d(coords1, coords2, radians = TRUE)
```

---

autoplot.evgram      *Plot evgram object*

---

**Description**

Plot an evgram object produced by the [evgram](#) function. The plotting function internally calls the [autoplot](#) function. Note: the `ggplot2` package must be loaded (i.e., `library(autoplot)` or `ggplot2::autoplot` must be specifically called for this function to work. See Examples.

**Usage**

```
autoplot.evgram(object, ..., split = FALSE)
```

**Arguments**

<code>object</code>	An evgram object produced by the <a href="#">evgram</a> function.
<code>...</code>	Not used
<code>split</code>	A logical value indicating whether, for a directional semivariogram, the directional semivariograms should be displayed in a single or split panels. Default is FALSE, for a single panel.

**Author(s)**

Joshua French

## Examples

```
data(co)
v = evgram(A1 ~ 1, co, ~ easting + northing)
if (requireNamespace("ggplot2")) {
  ggplot2::autoplot(v)
}
v2 = evgram(A1 ~ 1, co, ~ easting + northing, angle = 22.5, ndir = 4)
# ggplot2 must manually be loaded for this to work
if (requireNamespace("ggplot2")) {
  ggplot2::autoplot(v2)
  ggplot2::autoplot(v2, split = TRUE)
}
```

---

cmod.man

*Standard covariance models for geostatistical data.*

---

## Description

cmod.man manually creates a covariance matrix object (cmodMan) for geostatistical data. This function will be deprecated in the future. Please update your code to use the [cmod\\_man](#) function.

## Usage

```
cmod.man(v, evar = 0)
```

## Arguments

v	The covariance matrix of the observed data, including any errors. The matrix should be square, symmetric, and positive definite, though that latter two conditions are not checked.
evar	The variance of the errors. Must be non-negative number. The default is 0.

## Details

Note that  $v$  includes the error variance, i.e.,  $v = v_z + v_e$ , where  $v_z$  is the covariance matrix of the filtered process and the variance matrix of the errors is  $v_e = \text{diag}(\text{evar}/\text{weights})$ , where the weights come from the `geo1m` object the `cmodMan` object is associated with.

## Value

Returns a `cmodMan` object.

## Author(s)

Joshua French

**Examples**

```

coords = matrix(runif(20), ncol = 2)
d = as.matrix(dist(coords))
cmod.man(v = exp(-d), evar = 1)

```

cmod.std

*Standard covariance models for geostatistical data.***Description**

Creates a standard covariance model (cmodStd) object for geostatistical data. This function will be deprecated in the future. Please update your code to use the `cmod_std` function, which also allows the user to specify geometric anisotropy.

**Usage**

```
cmod.std(model, psill, r, evar = 0, fvar = 0, par3 = 0.5)
```

**Arguments**

model	A covariance model (e.g., "exponential"). See Details for the complete list of choices.
psill	The partial sill of the model. Must be a positive number.
r	The range parameter r. Must be a positive number.
evar	The variance of the errors. Must be non-negative number. The default is 0.
fvar	The finescale variance (microscale error). Must be a non-negative number. The default is 0.
par3	The value of the third parameter for 3 parameter models. Must be a positive number. The default is 0.5.

**Details**

The general form of the specified covariance function is  $psill * \rho(d; r) + (evar + fvar) * (d == 0)$ , where  $\rho$  is the covariance function of the parametric models.

For the exponential model,  $\rho(d; r)$  is  $\exp(-d/r)$ .

For the gaussian model,  $\rho(d; r)$  is  $\exp(-d^2/r^2)$ .

For the matern model,  $\rho(d; r)$  is  $2^{(1-par3)}/\gamma(par3)*sd^{par3}*besselK(sd, nu = par3)$ , where  $sd = d/r$ .

For the amatern (alternative Matern) model,  $\rho(d; r)$  is  $2^{(1-par3)}/\gamma(par3)*sd^{par3}*besselK(sd, nu = par3)$ , where  $sd = 2 * \sqrt{par3} * d/r$ .

For the spherical model,  $\rho(d; r)$  is  $1 - 1.5*sd + 0.5*(sd)^3$  if  $d < r$ , and 0 otherwise, with  $sd = d/r$ .

For the wendland1 model,  $\rho(d; r)$  is  $(1 - sd)^4 * (4*sd + 1)$  if  $d < r$ , and 0 otherwise, with  $sd = d/r$ .

For the wendland2 model,  $\rho(d; r)$  is  $(1 - sd)^6 * (35*sd^2 + 18*sd + 3)/3$  if  $d < r$ , and 0 otherwise, with  $sd = d/r$ .

For the wu1 model,  $\rho(d; r)$  is  $(1 - sd)^3 * (1 + 3*sd + sd^2)$  if  $d < r$ , and 0 otherwise, with  $sd = d/r$ .

For the wu2 model,  $\rho(d; r)$  is  $(1 - sd)^4 * (4 + 16*sd + 12*sd^2 + 3*sd^3)/4$  if  $d < r$ , and 0 otherwise, with  $sd = d/r$ .

For the wu3 model,  $\rho(d; r)$  is  $(1 - sd)^6 * (1 + 6*sd + 41/3*sd^2 + 12*sd^3 + 5*sd^4 + 5/6*sd^5)$  if  $d < r$ , and 0 otherwise, with  $sd = d/r$ .

### Value

Returns a cmodStd object.

### Author(s)

Joshua French

### References

Waller, L. A., & Gotway, C. A. (2004). Applied Spatial Statistics for Public Health Data. John Wiley & Sons.

### Examples

```
cmod.std(model = "exponential", psill = 1, r = 1)
```

---

cmod\_man

*Manual covariance models for geostatistical data.*

---

### Description

cmod\_man manually creates a covariance model object (cmodMan) for geostatistical data.

### Usage

```
cmod_man(v, evar = 0)
```

### Arguments

- |      |   |
|------|---|
| v    | The covariance matrix of the observed data, including any errors. The matrix should be square, symmetric, and positive definite, though that latter two conditions are not checked. |
| evar | The variance of the errors. Must be non-negative number. The default is 0.  |

**Details**

Note that  $v$  includes the error variance, i.e.,  $v = v_z + v_e$ , where  $v_z$  is the covariance matrix of the filtered process and the variance matrix of the errors is  $v_e = \text{diag}(\text{evar}/\text{weights})$ , where the weights come from the `geolm` object the `cmodMan` object is associated with.

**Value**

Returns a `cmodMan` object.

**Author(s)**

Joshua French

**Examples**

```
coords = matrix(runif(6), ncol = 2)
d = as.matrix(dist(coords))
cmod_man(v = exp(-d), evar = 1)
```

---

cmod\_std

*Standard covariance models for geostatistical data.*

---

**Description**

Creates a standard covariance model (`cmodStd`) object for geostatistical data.

**Usage**

```
cmod_std(
  model,
  psill,
  r,
  evar = 0,
  fvar = 0,
  par3 = 0.5,
  longlat = FALSE,
  angle = 0,
  ratio = 1,
  radians = FALSE,
  invert = TRUE
)
```

**Arguments**

model	A covariance model (e.g., "exponential"). See Details for the complete list of choices.
psill	The partial sill of the model. Must be a positive number.
r	The range parameter $r$ . Must be a positive number.
evar	The variance of the errors. Must be non-negative number. The default is 0.
fvar	The finescale variance (microscale error). Must be a non-negative number. The default is 0.
par3	The value of the third parameter for 3 parameter models. Must be a positive number. The default is 0.5.
longlat	A logical value indicating whether great circle distance should be used. The default is FALSE.
angle	The major axis of geometric anisotropy (the direction of strongest spatial dependence). Must be between $[0, 180)$ if <code>radians = FALSE</code> , otherwise it must be between $[0, \pi)$ .
ratio	The ratio of the minor axis range over the major axis range. The value must be between $(0, 1]$ .
radians	A logical value indicating whether the angles returned should be in degrees or radians. The default is FALSE, indicating that the returned angles are in degrees.
invert	A logical value indicating whether the axes of the coordinates should be inverted (i.e., the x- and y-axis are switched). The default is TRUE to mimic results from other geostatistical R packages like <code>gstat</code> , <code>geoR</code> , and other software like <code>GSLIB</code> and <code>GeoEAS</code> . Set to FALSE to use the typical x- and y-axes.

**Details**

The general, isotropic form of the specified covariance function is  $\text{psill} * \rho(d; r) + (\text{evar} + \text{fvar}) * (d == 0)$ , where  $\rho$  is the correlation function of the parametric models and  $d$  is the distance between the relevant coordinates.

For the `exponential` model,  $\rho(d; r)$  is  $\exp(-d/r)$ .

For the `gaussian` model,  $\rho(d; r)$  is  $\exp(-d^2/r^2)$ .

For the `matern` model,  $\rho(d; r)$  is  $2^{(1-\text{par3})}/\text{gamma}(\text{par3}) * \text{sd}^{\text{par3}} * \text{besselK}(\text{sd}, \text{nu} = \text{par3})$ , where  $\text{sd} = d/r$ .

For the `amatern` (alternative Matern) model,  $\rho(d; r)$  is  $2^{(1-\text{par3})}/\text{gamma}(\text{par3}) * \text{sd}^{\text{par3}} * \text{besselK}(\text{sd}, \text{nu} = \text{par3})$ , where  $\text{sd} = 2 * \text{sqrt}(\text{par3}) * d/r$ .

For the `spherical` model,  $\rho(d; r)$  is  $1 - 1.5 * \text{sd} + 0.5 * (\text{sd})^3$  if  $d < r$ , and 0 otherwise, with  $\text{sd} = d/r$ .

For the `wendland1` model,  $\rho(d; r)$  is  $(1 - \text{sd})^4 * (4 * \text{sd} + 1)$  if  $d < r$ , and 0 otherwise, with  $\text{sd} = d/r$ .

For the `wendland2` model,  $\rho(d; r)$  is  $(1 - \text{sd})^6 * (35 * \text{sd}^2 + 18 * \text{sd} + 3) / 3$  if  $d < r$ , and 0 otherwise, with  $\text{sd} = d/r$ .

For the `wu1` model,  $\rho(d; r)$  is  $(1 - \text{sd})^3 * (1 + 3 * \text{sd} + \text{sd}^2)$  if  $d < r$ , and 0 otherwise, with  $\text{sd} = d/r$ .



For the wu2 model,  $\rho(d; r)$  is  $(1 - sd)^4 * (4 + 16 * sd + 12 * sd^2 + 3 * sd^3) / 4$  if  $d < r$ , and 0 otherwise, with  $sd = d/r$ .

For the wu3 model,  $\rho(d; r)$  is  $(1 - sd)^6 * (1 + 6 * sd + 41/3 * sd^2 + 12 * sd^3 + 5 * sd^4 + 5/6 * sd^5)$  if  $d < r$ , and 0 otherwise, with  $sd = d/r$ .

### Value

Returns a `cmodStd` object.

### Author(s)

Joshua French

### References

Waller, L. A., & Gotway, C. A. (2004). *Applied Spatial Statistics for Public Health Data*. John Wiley & Sons.

### Examples

```
cmod_std(model = "exponential", psill = 1, r = 1)
```

---

co

*Geochemical measurements for 960 sites in Colorado.*

---

### Description

These data were collected by the United States Geological Survey (USGS). Their description is as follows: In 2006, soil samples were collected at 960 sites (1 site per 280 square kilometers) throughout the state of Colorado. These samples were collected from a depth of 0 to 15 centimeters and, following a near-total multi-acid digestion, were analyzed for a suite of more than 40 major and trace elements. The resulting data set provides a baseline for the natural variation in soil geochemistry for Colorado and forms the basis for detecting changes in soil composition that might result from natural processes or anthropogenic activities.

Latitude and Longitude determined by hand-held GPS instrument using WGS84 datum. The longitude and latitude coordinates were converted to UTM coordinates using the following commands:

```
library(sp)
lonlat = co[,c("longitude", "latitude")]
coordinates(lonlat) = c("longitude", "latitude")
proj4string(lonlat) = CRS("+proj=longlat +datum=WGS84")
xy = spTransform(lonlat, CRS("+proj=utm +zone=13 ellps=WGS84"))
```

### Usage

```
data(co)
```

**Format**

A data frame with 960 rows and 31 columns:

**easting** Easting (m)

**northing** Northing (m)

**latitude** The latitude of the site.

**longitude** The longitude of the site.

**Al** aluminum (percent)

**Ca** calcium (percent)

**Fe** iron (percent)

**K** potassium (percent)

**Mg** magnesium (percent)

**Na** sodium (percent)

**Ti** titanium (percent)

**Be** beryllium (mg/kg)

**Ce** cerium (mg/kg)

**Co** cobalt (mg/kg)

**Cr** chromium (mg/kg)

**Cu** copper (mg/kg)

**Ga** gallium (mg/kg)

**La** lanthanum (mg/kg)

**Li** lithium (mg/kg)

**Mo** manganese (mg/kg)

**Nb** molybdenum (mg/kg)

**Ni** niobium (mg/kg)

**Rb** rubidium (mg/kg)

**Sc** scandium (mg/kg)

**Sn** tin (mg/kg)

**Th** thorium (mg/kg)

**Tl** thallium (mg/kg)

**U** uranium (mg/kg)

**V** vanadium (mg/kg)

**W** tungsten (mg/kg)

**Y** yttrium (mg/kg)

**Source**

U.S. Geological Survey, Data Series 520, 9 p. <http://pubs.usgs.gov/ds/520/>.

**References**

Smith, D.B., Ellefsen, K.J., and Kilburn, J.E., 2010, Geochemical data for Colorado soils – Results from the 2006 state-scale geochemical survey: U.S. Geological Survey, Data Series 520, 9p.

---

decomp_cov	<i>Decompose covariance matrix</i>
------------	------------------------------------

---

**Description**

decomp\_cov decomposes a covariance matrix  $v$ . If  $A = \text{decomp\_cov}(v)$ , then  $\text{tcrossprod}(A, A) == v$ .

**Usage**

```
decomp_cov(v, method = "eigen")
```

```
decomp.cov(v, method = "eigen")
```

**Arguments**

$v$  An  $N \times N$  covariance matrix.  
 method The method used to decompose  $v$ . Valid options are "chol", "eigen", or "svd".

**Details**

The "chol" method is the fastest but least stable method. The "eigen" method is slower, but more stable. The "svd" method is the slowest method, but should be the most stable.

**Value**

Returns an  $N \times N$  matrix.

**Author(s)**

Joshua French

**Examples**

```
# generate data
n = 100
coords = matrix(runif(n*2), nrow = n, ncol = 2)
d = as.matrix(dist(coords))
# create covariance matrix
v = 3 * exp(-d/2) + 0.1 * diag(n)

# decompose v using the three methods
d1 = decomp_cov(v, "chol")
d2 = decomp_cov(v, "eigen")
d3 = decomp_cov(v, "svd")

# verify accuracy of decompositions
all.equal(v, tcrossprod(d1))
all.equal(v, tcrossprod(d2), check.attributes = FALSE)
all.equal(v, tcrossprod(d3), check.attributes = FALSE)
```

---

estimate	<i>Estimate model parameters</i>
----------	----------------------------------

---

### Description

estimate estimates the parameters of the specified model. The function is written to automatically adapt based on the class of object. Currently, the estimate function performs maximum likelihood estimation for objects produced by the `geolm` function.

### Usage

```
estimate(object, ...)
```

### Arguments

object	A model object produced by the <code>geolm</code> function.
...	Currently unimplemented

### Author(s)

Joshua French

### See Also

`estimate.geolm_cmodStd`,

### Examples

```
set.seed(10)
n = 100
```

---

estimate.geolm_cmodStd	<i>Determine MLEs of model parameters for a geostatistical model</i>
------------------------	--

---

### Description

estimate estimates the parameters of a geostatistical linear model of class `geolm_cmodStd` using maximum likelihood estimation.

**Usage**

```
## S3 method for class 'geolm_cmodStd'
estimate(
  object,
  reml = FALSE,
  noise_type = "e",
  lower = NULL,
  upper = NULL,
  method = "nlminb",
  itnmax = NULL,
  control = list(),
  est_nugget = TRUE,
  est_par3 = TRUE,
  est_angle = FALSE,
  est_ratio = FALSE,
  verbose = FALSE,
  ...
)
```

**Arguments**

object	A geostatistical linear model object produced by the <code>geolm</code> function.
reml	A logical value indicating whether standard maximum likelihood estimation should be performed ( <code>reml = FALSE</code> ). If <code>reml = TRUE</code> , then restricted maximum likelihood estimation is performed. The default is <code>FALSE</code> .
noise_type	A character vector indicating the type of noise (nugget) variance to estimate. The default is <code>est = "e"</code> , indicating the error variance should be estimated. Alternatively, the user can specify <code>est = "f"</code> , indicating that the finescale (microscale) variance should be estimated. The other type of noise variance is set to 0, otherwise the model is not identifiable. See Details.
lower	A named list with the names of the parameters you wish to set lower bounds for and the associated value. See Details.
upper	A named list with the names of the parameters you wish to set upper bounds for and the associated value.
method	The optimization method. The default is <code>"nlminb"</code> . <code>"L-BFGS-B"</code> is another acceptable choice. See <code>optimx</code> for further choices.
itnmax	An integer indicating the maximum number of iterations to allow for the optimization procedure.
control	A list of control parameters passed internally to <code>optimx</code> .
est_nugget	A logical value indicating whether the nugget variance ( <code>evar</code> or <code>fvar</code> ) should be estimated. The default is <code>TRUE</code> , indicating that the nugget should be estimated.
est_par3	A logical value indicating whether <code>par3</code> should be estimated (for an appropriate covariance model such as <code>"matern"</code> or <code>"amatern"</code> ). The default is <code>TRUE</code> , indicating that this parameter should be estimated.

est_angle	A logical value indicating whether the geometric anisotropy angle should be estimated. The default is FALSE, indicating that this parameter should not be estimated.
est_ratio	A logical value indicating whether the geometric anisotropy ratio of minor axis length to major axis length should be estimated. The default is FALSE, indicating that this parameter should not be estimated.
verbose	A logical value indicating whether potentially informative messages should be printed. The default is FALSE.
...	Currently unimplemented

### Details

The `optimx` function is used to find the MLEs. The control argument of `optimx` has a parameter `kkt` related to checking optimality conditions. This is internally set to FALSE. See `optimx` for Details.

Only the sum of `evar` and `fvar` is identifiable. Depending on the choice of `noise_type`, the covariance model is internally updated to estimate only one type of noise. e.g., if `noise_type = "e"`, then internally we update `evar` so that `evar = evar + fvar` and `fvar = 0`. Estimation is then performed on `evar` alone. Alternatively, the analogous estimated would be made for `fvar` if `noise_type = "fvar"`.

When `est_nugget` is true, the likelihood is profiled to simplify the optimization problem. In that case a parameter  $\lambda = (\text{evar} + \text{fvar})/\text{psill}$  is optimized. The optimal `psill` and noise variance are then determined.

The lower argument should be a named list with the names of the parameters you wish to set lower bounds for. If not specified, an attempt is made to specify reasonable lower bounds. The current choices are `r = 0.001`, `psill = 0.001`, `lambda = 0`, `angle = 0`, `ratio = 0.001`, `par3 = 0.001`.

The upper argument should be a named list with the names of the parameters you wish to set upper bounds for. If not specified, an attempt is made to specify reasonable upper bounds. The current choices are `r = 5 * maximum intercentroid distance`, `psill = 5 * var(object$y)`, `lambda = 5`, `angle = 179.99`, `ratio = 1`, `par3 = 3`.

### Author(s)

Joshua French

### See Also

[cmod\\_std](#), [optimx](#)

### Examples

```
data(toydata, package = "gear")
# setup standard covariance model
mod_std = cmod_std("exponential", psill = 1, r = 1, evvar = 0.1)
# setup dataframe with data
# fit Std geolm
object = geolm(y ~ x1 + x2, data = toydata, mod = mod_std,
               coordnames = c("x1", "x2"))
```

```
est_object = estimate(object, control = list(trace = 1),
                      verbose = TRUE,
                      lower = list(r = 0.05, lambda = 0.05))
```

---

`eval.cmod`*Evaluate covariance or semivariance model.*

---

## Description

`eval.cmod` evaluates the covariance of a model based on the provided arguments. This function will be deprecated in the future. Please use the [evaluate](#) function.

## Usage

```
eval.cmod(mod, d, coords = NULL)
```

## Arguments

<code>mod</code>	A covariance or semivariance model.
<code>d</code>	An $n \times m$ matrix of distances.
<code>coords</code>	Not used.

## Value

Returns the evaluated model with necessary components needed for [estimate](#) and [predict](#).

## Author(s)

Joshua French

## Examples

```
n = 10
coords = matrix(runif(2*n), nrow = n, ncol = 2)
d = as.matrix(dist(coords))
cmod = cmod_std(model = "exponential", psill = 1, r = 1)
eval.cmod(cmod, d)
```

---

<code>evaluate.cmodStd</code>	<i>Evaluate spatial dependence model</i>
-------------------------------	--

---

### Description

`evaluate` evaluates the spatial dependence model based on the provided arguments.

### Usage

```
## S3 method for class 'cmodStd'
evaluate(mod, d, e = TRUE, f = TRUE)

evaluate(mod, d, e = TRUE, f = TRUE)
```

### Arguments

<code>mod</code>	A covariance or semivariogram model.
<code>d</code>	An $n \times m$ matrix of distances. If <code>mod\$ratio != 1</code> , i.e., if geometric anisotropy has been specified, then <code>d</code> must be produced by the <a href="#">ganiso_d</a> function.
<code>e</code>	A single logical value indicating whether the error variance should be added to the returned covariance matrix. Default is TRUE.
<code>f</code>	A single logical value indicating whether the finescale/microscale variance should be added to the returned covariance matrix. Default is TRUE.

### Details

If `mod` is of class `cmodStd` (from the `cmod_std` function), then the function returns an  $n \times m$  matrix with the evaluated standard covariance function.

### Value

Returns the evaluated model with necessary components needed for [estimate](#) and [predict](#).

### Author(s)

Joshua French

### Examples

```
n = 10
coords = matrix(runif(2*n), nrow = n, ncol = 2)
d = as.matrix(dist(coords))
cmod = cmod_std(model = "exponential", psill = 1, r = 1)
evaluate(cmod, d)
```



---

evgram	<i>Empirical (semi-)variogram</i>
--------	-----------------------------------

---

### Description

evgram computes the empirical semivariogram of data based on the specified formula indicating the response and trend. See Details. The variogram is twice the semivariogram. If a trend is specified, then the semivariogram is constructed using the residuals of `lm(formula, data)`.

### Usage

```
evgram(
  formula,
  data,
  coordnames = NULL,
  nbins = 10,
  maxd = NULL,
  angle = 0,
  ndir = 1,
  type = "standard",
  npmin = 2,
  longlat = FALSE,
  verbose = TRUE,
  invert = TRUE
)
```

### Arguments

formula	A formula describing the relationship between the response and any covariates of interest, e.g., <code>response ~ 1</code> . The variogram is computed for the residuals of the linear model <code>lm(formula, data)</code> .
data	A <code>data.frame</code> , <code>SpatialPointsDataFrame</code> , <code>SpatialPixelsDataFrame</code> , or <code>SpatialGridDataFrame</code> object.
coordnames	The columns of data containing the spatial coordinates, provided as a formula (e.g., <code>~ x + y</code> ), column numbers (e.g., <code>c(1, 2)</code> ), or column names (e.g., <code>c("x", "y")</code> ). The default is <code>NULL</code> .
nbins	The number of bins (tolerance regions) to use when estimating the empirical semivariogram.
maxd	The maximum distance used when calculating the semivariogram. Default is <code>NULL</code> , in which case half the maximum distance between coordinates is used.
angle	A single value (in degrees) indicating the starting direction for a directional variogram. The default is 0.
ndir	The number of directions for which to calculate a empirical semivariogram. The default is 1, meaning calculate an omnidirectional semivariogram.

type	The name of the estimator to use in the estimation process. The default is "standard", the typical method-of-moments estimator. Other options include "cressie" for the robust Cressie-Hawkins estimator, and "cloud" for a semi-variogram cloud based on the standard estimator. If "cloud" is specified, the nbins argument is ignored.
npmin	The minimum number of pairs of points to use in the semivariogram estimator. For any bins with fewer points, the estimate for that bin is dropped.
longlat	A logical indicating whether Euclidean (FALSE) or Great Circle distance (WGS84 ellipsoid) (longlat = TRUE) should be used. Default is FALSE.
verbose	Logical value indicating whether computation information should be printed. Default is TRUE.
invert	A logical value indicating whether the axes of the coordinates should be inverted (i.e., the x- and y-axis are switched). The default is TRUE to mimic results from other geostatistical R packages like gstat, geoR, and other software like GSLIB and GeoEAS. Set to FALSE to use the typical x- and y-axes.

### Details

Note that the directions may be different from other packages (e.g., gstat or geoR packages) because those packages calculate angles clockwise from the y-axis, which is a convention frequently seen in geostatistics (e.g., the GSLIB software library). If `invert = TRUE`, the directions should be the same.

Computing the empirical semivariogram for the residuals of `lm(response ~ 1)` will produce identical results to simply computing the empirical semivariogram from the original response. However, if a trend is specified (the righthand side of `~` has non-trivial covariates), then the empirical semivariogram of the residuals will differ from that of the original response. A trend should be specified when the mean is non-stationary over the spatial domain.

### Value

Returns an `evgram`.

### Author(s)

Joshua French

### Examples

```
data(co)
v = evgram(A1 ~ 1, co, ~ easting + northing)
plot(v)
v2 = evgram(A1 ~ 1, co, c("easting", "northing"), angle = 22.5, ndir = 4)
plot(v2)
```

---

fitted.geolm	<i>Extract fitted values from a geolm object</i>
--------------	--

---

**Description**

Extract the fitted values, i.e., the estimated mean values, for an object produced by the `geolm` function for a specified set of covariates, `x`. If `x` is `NULL`, then `x` is taken from object.

**Usage**

```
## S3 method for class 'geolm'
fitted(object, x, ...)
```

**Arguments**

<code>object</code>	An object produced by the <code>geolm</code> function.
<code>x</code>	A $m \times p$ matrix of covariates for the locations where fitted values are desired. If <code>NULL</code> , <code>object\$x</code> is used.
<code>...</code>	Not currently implemented.

**Details**

If the object has a known mean, i.e., `object$mu` is not `NULL`, then the function returns the vector `rep(object$mu, m)`. If object has estimated coefficients, then `x %*% object$coeff` is returned.

If `x` is missing, then `object$x` is used for `x`. Naturally, `ncol(x)` must equal `length(object$coeff)`. If `x` is `NULL` and `object$mu` is not `NULL`, then `m` is taken to be 1.

**Value**

The vector of fitted values.

**Author(s)**

Joshua French

**See Also**

[fitted](#)

**Examples**

```
data = data.frame(y = rnorm(10), x1 = runif(10),
                  x2 = runif(10))
d = as.matrix(dist(data[,c("x1", "x2")]))
mod = cmod_man(v = exp(-d), evar = 1)
gearmod = geolm(y ~ x1, data = data,
                coordnames = ~ x1 + x2, mod = mod)
# fitted values for original observations
```

```
fitted(gearmod)
# fitted values for new observations
fitted(gearmod, x = cbind(1, rnorm(20)))
```

---

ganiso_d	<i>Anisotropic distance-related characteristics</i>
----------	---

---

### Description

Computes necessary distance-related characteristics when there is geometric anisotropy. This is essentially an internal function to `evaluate` a `cmodStd` object produced by `cmod_std` when the anisotropy ratio differs from 1.

### Usage

```
ganiso_d(coords1, coords2, radians = TRUE, invert = TRUE)
```

### Arguments

<code>coords1</code>	An $N \times 2$ matrix of spatial coordinates.
<code>coords2</code>	An $M \times 2$ matrix of spatial coordinates. Is missing, then <code>coords2 = coords1</code> .
<code>radians</code>	A logical value indicating whether the angles returned should be in degrees or radians. The default is <code>FALSE</code> , indicating that the returned angles are in degrees.
<code>invert</code>	A logical value indicating whether the axes of the coordinates should be inverted (i.e., the x- and y-axis are switched). The default is <code>TRUE</code> to mimic results from other geostatistical R packages like <code>gstat</code> , <code>geoR</code> , and other software like <code>GSLIB</code> and <code>GeoEAS</code> . Set to <code>FALSE</code> to use the typical x- and y-axes.

### Value

A `ganisoD` object with components `d` and `angles`, which is the distance matrix between the coordinates and the angles between the coordinates. The angles are returned in radians.

### Examples

```
ganiso_d(cbind(0, 0), cbind(1, 1))
```

---

gear	<i>gear</i>
------	-------------

---

### Description

\*Ge\*ostatistical \*A\*nalysis in \*R\*.

---

geardf	<i>Construct a geardf</i>
--------	---------------------------

---

### Description

`geardf` constructs a `geardf`, which is simply a `data.frame` with an attribute indicating which columns refer to the coordinates. The function either combines `x` and `coords` if `coordnames` isn't provided, or identifies the columns of `x` to which `coordnames` refers. The `geardf` class is only provided to make plotting results of the `predict.geolm*` functions simple without depending on the `sp` or `sf` packages. See [plot.geardf](#) for easy plotting.

### Usage

```
geardf(data, coords, coordnames)
```

### Arguments

<code>data</code>	A <code>data.frame</code>
<code>coords</code>	A <code>data.frame</code> or matrix with column names
<code>coordnames</code>	The columns of <code>data</code> containing the spatial coordinates, provided as a formula (e.g., <code>~ x + y</code> ), column numbers (e.g., <code>c(1,2)</code> ), or column names (e.g., <code>c("x", "y")</code> ). The default is <code>NULL</code> .

### Value

A `geardf`

### See Also

[plot.geardf](#)

### Examples

```
dtf = data.frame(a = 1:2, b = 3:4)

# create geardf with matrix coords (note column names)
coords = matrix(rnorm(4), ncol = 2)
colnames(coords) = c("u", "v")
geardf(dtf, coords)

# create geardf with data.frame coords
coords = as.data.frame(coords)
geardf(dtf, coords)

# create geardf using coordnames
dtf2 = cbind(dtf, u = rnorm(2), v = rnorm(2))
# vector form of coordnames
geardf(dtf2, coordnames = c("u", "v"))
```

```

# formula form of coordnames
geardf(dtf2, coordnames = ~ u + v)
# column index forum of coordnames
geardf(dtf2, coordnames = 3:4)

gdf = geardf(dtf2, coordnames = 3:4)
# looks like a data.frame
gdf
# but slightly more complicated
class(gdf)
attr(gdf, "coordnames")

```

---

geodist

---

*Compute distance for geographic coordinates*


---

### Description

geodist computes the distance between the coordinates in coords1 and coords2. If coords2 isn't supplied, then the distances are computed between the coordinates in coords1 alone. Otherwise, the pairwise distances between then points in coords1 and coords2 is computed. If longlat = TRUE, then the great circle distance is computed.

### Usage

```
geodist(coords1, coords2, longlat = FALSE)
```

### Arguments

coords1	A two-dimensional matrix of coordinates.
coords2	A two-dimensional matrix of coordinates.
longlat	A logical value indicating whether Euclidean distance (longlat = FALSE) or great circle distance (longlat = TRUE) should be computed. The default is longlat = FALSE.

### Details

The algorithm used when longlat = TRUE is a C++ port of the C code written by Roger Bivand for `spDists`, which appears to be based on a special case of the Vincenty formula with a slight correction based on the WGS84 flattening constant. See [https://en.wikipedia.org/wiki/Great-circle\\_distance](https://en.wikipedia.org/wiki/Great-circle_distance).

### Value

A matrix of distances

### Examples

```

coords = matrix(runif(10), ncol = 2)
d = geodist(coords)
all.equal(d, as.matrix(dist(coords)), check.attributes = FALSE)

```

---

geolm *Linear model for geostatistical data.*

---

## Description

geolm creates a geostatistical linear model object of the appropriate class based on the arguments, especially the cmod arguments.

## Usage

```
geolm(  
  formula,  
  data,  
  coordnames,  
  mod,  
  weights = NULL,  
  mu = NULL,  
  longlat = NULL,  
  cmod = NULL  
)
```

## Arguments

formula	An object of class <a href="#">formula</a> providing a symbolic description of the model to be fitted. See Details of this function and <a href="#">lm</a> .
data	A data frame containing the response, covariates, and location coordinates.
coordnames	The columns of data containing the spatial coordinates, provided as a formula (e.g., $\sim x + y$ ), column numbers (e.g., <code>c(1,2)</code> ), or column names (e.g., <code>c("x", "y")</code> )
mod	A model object produced by one of the <code>cmod_*</code> functions, e.g., <a href="#">cmod_std</a> .
weights	An optional vector of weights for the errors to be used in the fitting process. A vector that is proportional to the reciprocal variances of the errors, i.e., errors are assumed to be uncorrelated with variances <code>evar/weights</code> . Default is <code>NULL</code> , meaning that the weights are uniformly 1.
mu	A single numeric value indicating the constant mean of the spatial process if simple kriging is desired. Default is <code>NULL</code> , meaning that ordinary or universal kriging should be used.
longlat	A logical indicating whether Euclidean ( <code>FALSE</code> ) or Great Circle distance (WGS84 ellipsoid) ( <code>longlat = TRUE</code> ) should be used. Default is <code>FALSE</code> .
cmod	Retained for backwards compatibility. A model object produced by one of the <code>cmod_*</code> functions, e.g., <a href="#">cmod_std</a> .

**Details**

Note: for the multiresolution Gaussian process model, if `cmod$est == "f"` (i.e., if the nugget is finescale instead of measurement error), then the `weights` argument is internally set to `rep(1, n)`, where `n` is the number of observations.

formula should be specified after the form  $y \sim x_1 + x_2$ , where `y` is the response variable and `x1` and `x2` are the covariates of interest. If `mu` is provided, the variables to the right of `~` are ignored.

**Value**

Returns a `geolm_*` object, where `*` depends on `mod`.

**Author(s)**

Joshua French

**Examples**

```
data = data.frame(y = rnorm(10), x1 = runif(10),
                 x2 = runif(10))
d = geodist(data[,c("x1", "x2")])
mod = cmod_man(v = exp(-d), evar = 1)
gearmod = geolm(y ~ x1, data = data,
               coordnames = ~ x1 + x2, mod = mod)
```

---

geolm\_fit

*Fit a geolm*

---

**Description**

`geolm_fit` fits a `geolm` based on the specified `mod`. This is effectively an internal function.

**Usage**

```
geolm_fit(
  mod,
  x,
  y,
  coords,
  mu,
  weights,
  formula,
  coordnames,
  n,
  call,
  coeff_names
)
```



**Arguments**

mod	A model object produced by one of the <code>cm<sub>od</sub>_*</code> functions, e.g., <code>cm<sub>od</sub>_std</code> .
x	The matrix of covariates.
y	The vector of observed responses.
coords	The coordinates of the observed data set.
mu	A single numeric value indicating the constant mean of the spatial process if simple kriging is desired. Default is <code>NULL</code> , meaning that ordinary or universal kriging should be used.
weights	A vector that is proportional to the reciprocal variances of the errors.
formula	An object of class <code>formula</code> providing a symbolic description of the model to be fitted. See Details of this function and <code>lm</code> .
coordnames	A vector of length 2 with the names of the columns in data containing the coordinates, e.g., <code>c("long", "lat")</code> .
n	The number of observations.
call	The <code>match.call</code> for the <code>ge<sub>o</sub>lm</code> .
coeff_names	A character string with the variable names associated with the coefficients.

**Value**

Returns a `geolm` with necessary components needed for `estimate` and `predict`.

**Author(s)**

Joshua French

**Examples**

```
# no examples since you shouldn't be using this function!
```

---

mle	<i>Finds maximum likelihood estimates of model parameters for a geostatistical model</i>
-----	--

---

**Description**

`mle` estimates the parameters of a geostatistical linear model.

`mle` estimates the parameters of a geostatistical linear model. The `mle` function will be deprecated in the future. Please update your code to use the `estimate` function.

**Usage**

```
mle(
  object,
  reml = FALSE,
  est = "e",
  lower = NULL,
  upper = NULL,
  method = "nlminb",
  itnmax = NULL,
  control = list(),
  ...
)
```

```
mle(
  object,
  reml = FALSE,
  est = "e",
  lower = NULL,
  upper = NULL,
  method = "nlminb",
  itnmax = NULL,
  control = list(),
  ...
)
```

**Arguments**

<code>object</code>	A geostatistical linear model object produced by the <code>geolm</code> function.
<code>reml</code>	A logical value indicating whether standard maximum likelihood estimation should be performed ( <code>reml = FALSE</code> ). If <code>reml = TRUE</code> , then restricted maximum likelihood is performed. Default is <code>FALSE</code> .
<code>est</code>	A character vector indicator whether the error variance ( <code>est="e"</code> ) or finescale variance ( <code>est = "f"</code> ) should be estimated. The other component of the nugget variance is held constant, and in the case of a <code>geolmStd</code> object, is set to 0.
<code>lower</code>	A vector of 2 or 3 specifying the lowerbound of parameter values. See Details.
<code>upper</code>	lower A vector of 2 or 3 specifying the lowerbound of parameter values. See Details.
<code>method</code>	The optimization method. Default is <code>"nlminb"</code> , with <code>"L-BFGS-B"</code> being another acceptable choice. See <a href="#">optimx</a> for details.
<code>itnmax</code>	An integer indicating the maximum number of iterations to allow for the optimization procedure.
<code>control</code>	A list of control parameters passed internally to <a href="#">optimx</a> . See <a href="#">optimx</a> for details.
<code>...</code>	Currently unimplemented.

## Details

In the case of a `geolmStd` object, the likelihood has been concentrated so that only the range parameter `r` and a scale parameter `lambda = nugget/psill` need to be estimated.

If object is a `geolmStd`, then `lower` is of length 2 if the covariance model of `cmo` is not `matern` or `amatern`. Otherwise, it should be of length 3. The first parameter is related to the range parameter `r`, the second to the scale parameter `lambda`, and the third to `par3`, if applicable. If `lower = NULL`, then the lower bounds are 0.001, 0, and 0.1, respectively. A similar pattern holds for `upper`, with the default being  $3 * \max(d)$ , where `d` is the matrix of distances between coordinates, 5, and 2.5.

The `kkt` argument in the `control` list is set to be `FALSE`.

## Author(s)

Joshua French

Joshua French

## See Also

[estimate](#),

## Examples

```
set.seed(10)
n = 100
set.seed(10)
n = 100
```

---

ploglik\_ycholv

*Compute the log-likelihood of a model*

---

## Description

`ll_ycholv` computes the log-likelihood of multivariate normal data using components typically found in a `geolm`. For `ploglik_ycholv`, `cholv` is the cholesky decomposition of the covariance matrix for the observed data after dividing the matrix by the (estimated) `psill`. See the examples below. Depending on parameter choices, the function can return the log-likelihood, the restricted log-likelihood, -2 times the log-likelihood or restricted log-likelihood, or the estimated partial sill for both a maximum likelihood and restricted maximum likelihood setting. This is intended to be an internal function, so minimal error checking is done.

## Usage

```
ploglik_ycholv(
  x,
  y,
  cholv,
  mu = NULL,
```

```

    reml = FALSE,
    minus2 = TRUE,
    return_ll = TRUE
  )

  ll_xycholv(
    x,
    y,
    cholv,
    mu = NULL,
    reml = FALSE,
    minus2 = TRUE,
    return_ll = TRUE
  )

```

### Arguments

<code>x</code>	The matrix of covariates.
<code>y</code>	The vector of observed responses.
<code>cholv</code>	The cholesky decomposition of the covariance matrix of <code>y</code> (or of that matrix divided by <code>psill</code> for <code>loglik_xycholv</code> ).
<code>mu</code>	A single numeric value indicating the assumed mean of the underlying process.
<code>reml</code>	A logical value. Should the Restricted Maximum Likelihood be returned. The default is <code>FALSE</code> .
<code>minus2</code>	A logical value. Should -2 times the log-likelihood be returned. The default is <code>TRUE</code> .
<code>return_ll</code>	A logical value. Should the log-likelihood be returned? Default is <code>TRUE</code> . If <code>FALSE</code> , the estimated partial sill is returned.

### Value

A likelihood value, -2 times the likelihood value, or the estimated partial sill, depending on the user's argument choices.

### References

Statistical Methods for Spatial Data Analysis. Oliver Schabenberger and Carol A. Gotway (Chapman & Hall/CRC Press) 2005. pp. 259-263

### Examples

```

y = rnorm(10)
x = matrix(rep(1, length(y)))
coords = matrix(runif(length(y) * 2), ncol = 2)
d = as.matrix(dist(coords))
pv = exp(-d/3) + 0.1 * diag(length(y))
est_psill = ploglik_xycholv(x, y, chol(pv), return_ll = FALSE)
v = pv * est_psill

```

```
# same result
ploglik_xychoLv(x, y, chol(pv), minus2 = FALSE)
ll_xychoLv(x, y, chol(v), minus2 = FALSE)
```

---

plot.evgram	<i>Plot evgram object</i>
-------------	---------------------------

---

### Description

Plots evgram object produced by [evgram](#) using the [plot](#) function.

### Usage

```
## S3 method for class 'evgram'
plot(x, ..., split = FALSE, add_legend = TRUE, args_legend = list())
```

### Arguments

x	An evgram object produced by the <a href="#">evgram</a> function.
...	Additional arguments to pass the <a href="#">plot</a> function to change aspects of the plot.
split	A logical value indicating whether, for a directional semivariogram, the directional semivariograms should be displayed in a single or split panels. Default is FALSE, for a single panel.
add_legend	A logical value indicating whether a legend should be included for a plot of directional semivariograms when <code>split = FALSE</code> .
args_legend	A list with arguments for <a href="#">legend</a> .

### Details

The arguments after ... are only considered if a directional semivariogram is provided, i.e., when `x$ndir != 1`.

`split` will split directional semivariograms into different plots automatically using [autosize](#).

`add_legend` and `args_legend` are only used for a directional semivariogram when `split = FALSE`

### Author(s)

Joshua French

### See Also

[xyplot](#), [evgram](#)

## Examples

```

data(co)
# omnidirectional example
v = evgram(A1 ~ 1, co, ~ easting + northing)
plot(v)
plot(v, main = "semivariogram of A1")

# directional semivariograms overlaid
v2 = evgram(A1 ~ 1, co, ~ easting + northing,
            angle = 22.5, ndir = 4)
plot(v2)
plot(v2, ylab = "semi-variance", pch = 2:5, type = "p")
plot(v2, lty = 2:5, type = "l",
     args_legend = list(x = "bottomright",
                       legend = c("22.5", "67.5", "112.5", "157.5")))
# directional semivariograms split
plot(v2, split = TRUE)
plot(v2, split = TRUE, col = 2, pch = 3, type = "b",
     main = c("(a)", "(b)", "(c)", "(d)"))

```

---

plot.geardf

*Plot geardf object*

---

## Description

Plot a geardf object produced by the [geardf](#) or `predict.geolm*` functions. See the [autopoints](#) and [autoimage](#) functions for advanced options.

## Usage

```

## S3 method for class 'geardf'
plot(x, zcol = names(x), interp = FALSE, common.legend = FALSE, ...)

```

## Arguments

<code>x</code>	A geardf object produced by the <a href="#">geardf</a> or <code>predict.geolm*</code> functions.
<code>zcol</code>	The names of the columns of <code>x</code> to plot. The coordinate columns are automatically stripped from <code>zcol</code> .
<code>interp</code>	A logical value indicating whether the values should be interpolated onto a grid. If FALSE, then the <a href="#">autopoints</a> function is used to construct a heated scatterplot. If TRUE, then the <a href="#">autoimage</a> function is used to create a heat map. The default is FALSE.
<code>common.legend</code>	A logical value indicating whether a common legend should be used for the scatterplots/images. The default is FALSE.
<code>...</code>	Additional arguments to be passed to the <a href="#">autopoints</a> or <a href="#">autoimage</a> functions.

**Author(s)**

Joshua French

**See Also**[autopoints](#), [autoimage](#)**Examples**

```

data(toydata)
# newdata must have columns with prediction coordinates
newdata = data.frame(x1 = runif(10), x2 = runif(10))

# specify a standard covariance model
mod = cmod_std(model = "exponential", psill = 1, r = 1)

# geolm for universal kriging
geolm_uk = geolm(y ~ x1 + x2, data = toydata, mod = mod,
                coordnames = c("x1", "x2"))

# prediction for universal kriging
pred_uk = predict(geolm_uk, newdata, return_type = "geardf")
# heated scatterplot
plot(pred_uk)
# interpolated image of results
plot(pred_uk, interp = TRUE)
# plot only predictions and rmspe with different colors
plot(pred_uk, c("pred", "rmspe"), col = cm.colors(5))
# plot only predictions with coarser interpolation grid
plot(pred_uk, "pred", interp = TRUE,
      interp.args = list(no.X = 10, no.Y = 10))

```

---

predict.geolm\_cmodMan *Predict method for geostatistical models*

---

**Description**

predict calculates the predicted values at specified locations. The method can additionally provide the mean square prediction error (mspe) and perform conditional simulation.

**Usage**

```

## S3 method for class 'geolm_cmodMan'
predict(
  object,
  newdata,
  nsim = 0,
  vop,
  vp,

```

```

return_type = "SpatialPointsDataFrame",
dmethod = "chol",
compute_mspe = TRUE,
sp = NULL,
...
)

```

## Arguments

object	An object produced by the <code>geolm</code> function.
newdata	An optional data frame in which to look for the coordinates at which to predict. If omitted, the observed data locations are used.
nsim	A non-negative integer indicating the number of realizations to sample at the specified coordinates using conditional simulation.
vop	The cross-covariance matrix between the observed responses and the responses to predict.
vp	The covariance matrix of the responses to predict.
return_type	A character string indicating the type of object that should be returned. The default is " <code>SpatialPointsDataFrame</code> " for easy plotting of results (see Examples). Other options include " <code>data.frame</code> ", " <code>geardf</code> ", and " <code>sf</code> ".
dmethod	The method used to decompose the covariance matrix for conditional simulation. Valid options are " <code>chol</code> ", " <code>eigen</code> ", and " <code>svd</code> ". The default is " <code>chol</code> ".
compute_mspe	A logical value indicating whether the mean square prediction error should be calculated. Default is TRUE.
sp	This argument will be deprecated in the future. Please use the <code>return_type</code> argument. A logical value indicating whether to object returned should be of class <code>SpatialPointsDataFrame</code> for easier plotting with the <code>sp</code> package. Default is NULL.
...	Currently unimplemented.

## Details

The `newdata` data frame must include the relevant covariates for the prediction locations, where the covariates are specified on the right side of the `~` in `object$formula`. `newdata` must also include the coordinates of the prediction locations, with these columns having the names provided in `object$coordnames`.

## Value

A `data.frame`, `SpatialPointsDataFrame`, `geardf`, or `sf` object with the kriging predictions `pred`, kriging variance/mean-square prediction error (`mspe`), the root mean-square prediction error `mspe` (`rmspe`), and the conditional simulations `sim.1`, `sim.2`, etc. `sim.1`, `sim.2`, etc.

## Author(s)

Joshua French



**Examples**

```

# generate response
y = rnorm(10)
# generate coordinates
x1 = runif(10); x2 = runif(10)

# data frame for observed data
data = data.frame(y, x1, x2)
coords = cbind(x1, x2)
d = as.matrix(dist(coords))
psill = 2 # partial sill
r = 4 # range parameter
evar = .1 # error variance
fvar = .1 # add finescale variance
# one can't generally distinguish between evan and fvar, but
# this is done for illustration purposes

# manually specify an exponential covariance model
v = psill * exp(-d/r) + (evan + fvar) * diag(10)
mod_man = cmod_man(v = v, evan = evan)

# coordinate names
cnames = c("x1", "x2")

# geolm for universal kriging
gearmod_uk = geolm(y ~ x1 + x2, data = data, mod = mod_man,
                  coordnames = cnames)

# newdata must have columns with prediction coordinates
# add 5 unsampled sites to sampled sites
newdata = data.frame(x1 = c(x1, runif(5)), x2 = c(x2, runif(5)))
newcoords = newdata[, cnames]
# create vop and vp using distances
dop = geodist(as.matrix(coords), as.matrix(newcoords))
dp = geodist(newcoords)

# manually create cross-covariance and covariance for
# prediction locations
vop = psill * exp(-dop/r) + fvar * (dop == 0)
vp = psill * exp(-dp/r) + fvar * diag(nrow(newcoords))

# prediction for universal kriging, with conditional simulation,
# using manual covariance matrices
pred_uk_man = predict(gearmod_uk, newdata, nsim = 2,
                    vop = vop, vp = vp, dmethod = "svd")

# do the same thing, but using cmod_std

# prediction for universal kriging, with conditional simulation
mod_std = cmod_std("exponential", psill = psill, r = r,
                  evan = evan, fvar = fvar)
gearmod_uk2 = geolm(y ~ x1 + x2, data = data, mod = mod_std,

```

```

      coordnames = c("x1", "x2"))
pred_uk_std = predict(gearmod_uk2, newdata, nsim = 2, dmethod = "svd")

# compare results
all.equal(pred_uk_man$pred, pred_uk_std$pred)
all.equal(pred_uk_man$mspe, pred_uk_std$mspe)

```

---

predict.geolm\_cmodStd *Predict method for geostatistical models*

---

## Description

predict calculates the predicted values at specified locations. The method can additionally provide the mean square prediction error (mspe) and perform conditional simulation.

## Usage

```

## S3 method for class 'geolm_cmodStd'
predict(
  object,
  newdata,
  nsim = 0,
  return_type = "SpatialPointsDataFrame",
  dmethod = "chol",
  compute_mspe = TRUE,
  sp = NULL,
  ...
)

```

## Arguments

object	An object produced by the geolm function.
newdata	An optional data frame in which to look for the coordinates at which to predict. If omitted, the observed data locations are used.
nsim	A non-negative integer indicating the number of realizations to sample at the specified coordinates using conditional simulation.
return_type	A character string indicating the type of object that should be returned. The default is " <a href="#">SpatialPointsDataFrame</a> " for easy plotting of results (see Examples). Other options include " <a href="#">data.frame</a> ", " <a href="#">geardf</a> ", and " <a href="#">sf</a> ".
dmethod	The method used to decompose the covariance matrix for conditional simulation. Valid options are "chol", "eigen", and "svd". The default is "chol".
compute_mspe	A logical value indicating whether the mean square prediction error should be calculated. Default is TRUE.
sp	This argument will be deprecated in the future. Please use the return_type argument. A logical value indicating whether to object returned should be of class <a href="#">SpatialPointsDataFrame</a> for easier plotting with the sp package. Default is NULL.
...	Currently unimplemented.

**Details**

The newdata data frame must include the relevant covariates for the prediction locations, where the covariates are specified on the right side of the  $\sim$  in `object$formula`. newdata must also include the coordinates of the prediction locations, with these columns having the names provided in `object$coordnames`.

**Value**

A `data.frame`, `SpatialPointsDataFrame`, `geardf`, or `sf` object with the kriging predictions `pred`, kriging variance/mean-square prediction error (`mspe`), the root mean-square prediction error `mspe` (`rmspe`), and the conditional simulations `sim.1`, `sim.2`, etc. `sim.1`, `sim.2`, etc.

**Author(s)**

Joshua French

**Examples**

```
# generate response
y = rnorm(10)
# generate coordinates
x1 = runif(10); x2 = runif(10)

# data frame for observed data
data = data.frame(y, x1, x2)
# newdata must have columns with prediction coordinates
newdata = data.frame(x1 = runif(5), x2 = runif(5))

# specify a standard covariance model
mod = cmod_std(model = "exponential", psill = 1, r = 1)

# geolm for universal kriging
gearmod_uk = geolm(y ~ x1 + x2, data = data, mod = mod,
                  coordnames = c("x1", "x2"))
# prediction for universal kriging, with conditional simulation
pred_uk = predict(gearmod_uk, newdata, nsim = 2)

# demonstrate plotting abilities if return_type == "geardf"
pred_geardf = predict(gearmod_uk, newdata,
                    return_type = "geardf")
plot(pred_geardf, "pred")
plot(pred_geardf, interp = TRUE)

# demonstrate plotting abilities if sp package installed
if (requireNamespace("sp", quietly = TRUE)) {
  pred_spdf = predict(gearmod_uk, newdata,
                    return_type = "SpatialPointsDataFrame")
  sp::splot(pred_spdf, "pred")
  sp::splot(pred_spdf)
}
# demonstrate plotting abilities if sf package installed
```

```
if (requireNamespace("sf", quietly = TRUE)) {
  pred_sfdf = predict(gearmod_uk, newdata,
    return_type = "sf")
  plot(pred_sfdf["pred"])
  plot(pred_sfdf)
}

# geolm for ordinary kriging
gearmod_ok = geolm(y ~ 1, data = data, mod = mod,
  coordnames = c("x1", "x2"))
# prediction for ordinary kriging
pred_ok = predict(gearmod_ok, newdata)

# geolm for simple kriging
gearmod_sk = geolm(y ~ 1, data = data, mod = mod,
  coordnames = c("x1", "x2"), mu = 1)
# prediction for simple kriging
pred_sk = predict(gearmod_sk, newdata)
```

---

print.evgram

*Print evgram object*

---

## Description

Print an object of class `evgram` produced by the `evgram` function.

## Usage

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

## Arguments

`x` An `evgram` object produced by the `evgram` function.  
`...` Not currently implemented.

## Author(s)

Joshua French

## Examples

```
data(co)
evgram(A1 ~ 1, co, ~ easting + northing)
```

---

residuals.geolm	<i>Extract residuals from a geolm object</i>
-----------------	--

---

## Description

Extract the residuals for an object produced by the [geolm](#).

## Usage

```
## S3 method for class 'geolm'  
residuals(object, ...)
```

## Arguments

object	An object produced by the <a href="#">geolm</a> function.
...	Not currently implemented.

## Value

The vector of residuals.

## Author(s)

Joshua French

## See Also

[residuals](#)

## Examples

```
data = data.frame(y = rnorm(10), x1 = runif(10),  
                 x2 = runif(10))  
d = as.matrix(dist(data[,c("x1", "x2")]))  
mod = cmod_man(v = exp(-d), evar = 1)  
gearmod = geolm(y ~ x1, data = data,  
               coordnames = ~ x1 + x2, mod = mod)  
# fitted values for original observations  
residuals(gearmod)
```

---

`solve_chol`*Solve using cholesky decomposition*

---

### Description

`solve_chol` solves a system of equations using the cholesky decomposition of a positive definite matrix `A`, i.e., using `a = chol(A)`.

### Usage

```
solve_chol(a, b, ...)
```

```
## S3 method for class 'chol'  
solve(a, b, ...)
```

### Arguments

<code>a</code>	The cholesky decomposition of a positive definite matrix.
<code>b</code>	A numeric or complex vector or matrix giving the right-hand side(s) of the linear system. If missing, <code>b</code> is taken to be an identity matrix and <code>solve</code> will return the inverse of <code>a</code> .
<code>...</code>	Currently unused.

### Details

Note: Unless you have good reason to suspect that the cholesky decomposition of your matrix will be stable, it is recommended that you use [solve](#) or perform the solve using the [qr](#) decomposition.

### Author(s)

Joshua French

### See Also

[solve](#), [chol](#), [qr](#), [solve.qr](#)

### Examples

```
set.seed(10)  
# create positive definite matrix a  
a = crossprod(matrix(rnorm(25^2), nrow = 25))  
# create vector x and matrix b  
# x can be used to check the stability of the solution  
x = matrix(rnorm(25))  
b = a %*% x  
  
# standard solve  
x1 = solve(a, b)
```

```
all.equal(x, x1)

# solve using cholesky decomposition
chola = chol(a)
x2 = solve_chol(chola, b)
all.equal(x, x2)

# solve using qr decomposition
qra = qr(a)
x3 = solve.qr(qra, b)
all.equal(x, x3)

# compare direct inversion
ai1 = solve(a)
ai2 = solve_chol(chola) #using cholesky decomposition
all.equal(ai1, ai2) # should be TRUE
```

---

toydata

*A toy data set for examples*

---

### Description

This is a toy data set of 25 observations generated on a  $[0, 1] \times [0, 1]$  domain.

### Usage

```
data(toydata)
```

### Format

A data frame with 25 rows and 3 columns:

**y** response

**x1** coordinate dimension 1 values

**x2** coordinate dimension 2 values

---

update.geolm

*Update linear model for geostatistical data*

---

### Description

update updates a geostatistical linear model based on the given model.

**Usage**

```
## S3 method for class 'geolm'  
update(object, mod, ...)  
  
## S3 method for class 'geolm_cmodMan'  
update(object, mod, ...)  
  
## S3 method for class 'geolm_cmodStd'  
update(object, mod, ...)
```

**Arguments**

object	An object produced by the geolm function.
mod	A spatial dependence model object obtained from one of the cmod_* functions or the estimate function.
...	Not implemented.

**Value**

Returns an object of the same class as object.

**Author(s)**

Joshua French

**See Also**

[update](#)

**Examples**

```
# generate response  
y = rnorm(10)  
# generate coordinates  
x1 = runif(10); x2 = runif(10)  
  
# data frame for observed data  
data = data.frame(y, x1, x2)  
coords = cbind(x1, x2)  
psill = 2 # partial sill  
r = 4 # range parameter  
evar = .1 # error variance  
fvar = .1 # add finescale variance  
# one can't generally distinguish between evan and fvar, but  
# this is done for illustration purposes  
  
cmod_std = cmod_std("exponential", psill = psill, r = r,  
                   evan = evan, fvar = fvar)  
  
cmod_std2 = cmod_std("exponential", psill = psill + 1,
```



```

      r = r + .5, evar = evar + .01,
      fvar = fvar)

# check geolm update for universal kriging
gear1 = geolm(y ~ x1 + x2, data = data, mod = cmod_std,
             coordnames = c("x1", "x2"))

gear2 = geolm(y ~ x1 + x2, data = data, mod = cmod_std2,
             coordnames = c("x1", "x2"))
gear2b = update(gear1, cmod_std2)
gear2b$call = NULL
gear2b$call = NULL
identical(gear2, gear2b)

```

---

vgram

*Empirical variogram*


---

### Description

vgram calculates an empirical variogram. Note that, by convention, the empirical variogram actually estimates the semivariogram, not the theoretical variogram (which is twice the semivariogram).

### Usage

```

vgram(
  formula,
  data,
  coordnames = NULL,
  nbins = 10,
  maxd = NULL,
  angle = 0,
  ndir = 1,
  type = "standard",
  npmin = 2,
  longlat = FALSE,
  verbose = TRUE,
  coords = NULL
)

```

### Arguments

formula	A formula describing the relationship between the response and any covariates of interest, e.g., <code>response ~ 1</code> . The variogram is computed for the residuals of the linear model <code>lm(formula, data)</code> .
data	A <code>data.frame</code> , <code>SpatialPointsDataFrame</code> , <code>SpatialPixelsDataFrame</code> , or <code>SpatialGridDataFrame</code> object.

coordnames	The columns of data containing the spatial coordinates, provided as a formula (e.g., $\sim x + y$ ), column numbers (e.g., <code>c(1, 2)</code> ), or column names (e.g., <code>c("x", "y")</code> ). The default is <code>NULL</code> , but this must be specified if data is of class <code>data.frame</code> .
nbins	The number of bins (tolerance regions) to use when estimating the sample semi-variogram.
maxd	The maximum distance used when calculating the semivariogram. Default is <code>NULL</code> , in which case half the maximum distance between coordinates is used.
angle	A single value (in degrees) indicating the starting direction for a directional variogram. The default is 0.
ndir	The number of directions for which to calculate a sample semivariogram. The default is 1, meaning calculate an omnidirection semivariogram.
type	The name of the estimator to use in the estimation process. The default is "standard", the typical method-of-moments estimator. Other options include "cressie" for the robust Cressie-Hawkins estimator, and "cloud" for a semivariogram cloud based on the standard estimator. If "cloud" is specified, the <code>nbins</code> argument is ignored.
npmin	The minimum number of pairs of points to use in the semivariogram estimator. For any bins with fewer points, the estimate for that bin is dropped.
longlat	A logical indicating whether Euclidean ( <code>FALSE</code> ) or Great Circle distance (WGS84 ellipsoid) ( <code>longlat = TRUE</code> ) should be used. Default is <code>FALSE</code> .
verbose	Print computation information. Default is <code>TRUE</code> .
coords	A deprecated argument.

### Details

Note that the directions may be different from other packages (e.g., `gstat` or `geoR` packages) because those packages calculate angles clockwise from the  $y$ -axis, which is a convention frequently seen in geostatistics (e.g., the `GSLIB` software library).

Additionally, note that calculating the empirical variogram for the residuals of `lm(response ~ 1)` will produce identical results to simply computing the sample semivariogram from the original response. However, if a trend is specified (the righthand side of  $\sim$  has non-trivial covariates), then the empirical variogram of the residuals will differ from that of the original response. A trend should be specified when the mean is non-stationary over the spatial domain.

### Value

Returns an `evgram` object with components:

### Author(s)

Joshua French

**Examples**

```

data(co)
v = vgram(A1 ~ 1, co, ~ easting + northing)
plot(v)
v2 = vgram(A1 ~ 1, co, c("easting", "northing"), angle = 22.5, ndir = 4)
plot(v2)

```

xyplot.evgram

*Plot evgram object***Description**

Plots evgram object produced by the [evgram](#) function using the [xyplot](#) function. Note: the `lattice` package must be loaded (i.e., `library(lattice)` or `lattice::xyplot` must be specifically called for this function to work. See Examples.

**Usage**

```
xyplot.evgram(x, ..., split = FALSE)
```

**Arguments**

<code>x</code>	An evgram object produced by the <a href="#">evgram</a> function.
<code>...</code>	Additional arguments to pass the <a href="#">xyplot</a> function to change aspects of the plot.
<code>split</code>	A logical value indicating whether, for a directional semivariogram, the directional semivariograms should be displayed in a single or split panels. Default is <code>FALSE</code> , for a single panel.

**Author(s)**

Joshua French

**See Also**

[xyplot](#), [evgram](#)

**Examples**

```

data(co)
v = evgram(A1 ~ 1, co, ~ easting + northing)
if (requireNamespace("lattice")) {
  lattice::xyplot(v)
}
v2 = evgram(A1 ~ 1, co, ~ easting + northing, angle = 22.5, ndir = 4)
if (requireNamespace("lattice")) {
  lattice::xyplot(v2)
  # show how attributes can be changed using different
  # arguments available in lattice::xyplot.
}

```

```
lattice::xyplot(v2, col = 2:5)
lattice::xyplot(v2, col = 2:5, pch = 1:4)
lattice::xyplot(v2, col = 2:5, pch = 1:4, lty = 2:5, type = "b")
lattice::xyplot(v2, col = 2:5, pch = 1:4, lty = 2:5, type = "b",
  key=list(text=list(levels(as.factor(v2$semi$angle))),
    space='right', points=list(pch=1:4, col=2:5),
    lines=list(col=2:5, lty = 2:5)))
lattice::xyplot(v2, split = TRUE)
}
```

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