

Package 'dsm'

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Title Density surface modelling of distance sampling data

LazyLoad yes

Author David L. Miller, Eric Rexstad, Louise Burt, Mark V. Bravington, Sharon Hedley.

Description This library implements density surface modelling of line transect data, based on the methods of Hedley et al. (2004). Some recent developments in the literature have also be incorporated. Outputs are point and interval estimates of population abundance and density. Please note that this version of dsm WILL NOT work with the Windows package DISTANCE earlier than version 6.2. For a version that works with DISTANCE 6.1, please go to: <https://github.com/DistanceDevelopment/dsm-distance-6.1> For the latest version of dsm, please use the github version at the URL listed below. Miller, D. L., M. L. Burt, E. Rexstad and L. Thomas. 2013. Spatial models for distance sampling data: recent developments and future directions. *Methods in Ecology and Evolution* Hedley, S.L., S.T. Buckland and D.L. Borchers. 2004. "Spatial distance sampling methods" pp 48-70 in *Advanced Distance Sampling*, Buckland, S.T. et al. (eds). Oxford University Press.

Version 2.2.5

URL <http://github.com/DistanceDevelopment/dsm>

Date 2014-10-15

Depends R (>= 3.0), mgcv (>= 1.7), mrds (>= 2.1.8)

Imports nlme, ggplot2, statmod

Suggests Distance

NeedsCompilation no

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 dsm-package

Density surface modelling

Description

Some blurb will eventually go here.

block.info.per.su *Find the block information*

Description

Takes the transect data and works out how many blocks of a given size (in segment terms) fit into each.

Usage

```
block.info.per.su(block.size, data, name.su)
```

Arguments

block.size	number of segments per block
data	data used to build the model
name.su	names of the sampling units (ie. transects)

Value

a data.frame with the following columns

name	the sample unit name (e.g. transect label)
num.seg	number of segments in that transect
num.block	number of blocks available
start.block	block # for first block
end.block	block # for last block
num.req	number of blocks needed for the unit

check.cols *Check column names exist*

Description

Internal function to check that supplied 'data.frames' have the correct columns and checks that sample labels are all unique.

Usage

```
check.cols(ddf.obj, segment.data, observation.data, strip.width, segment.area)
```

Arguments

ddf.obj a ddf object from ‘mrds’
segment.data segment data as defined in [dsm](#)
observation.data observation data as defined in [dsm](#)
strip.width strip width if strip transects are being used
segment.area area of segments

Value

nothing, but throws an error if something went wrong

Author(s)

David Lawrence Miller

dsm	<i>Fit a density surface model to segment-specific estimates of abundance or density.</i>
-----	---

Description

Fits a density surface model (DSM) to detection adjusted counts from a spatially-referenced distance sampling analysis. [dsm](#) takes observations of animals, allocates them to segments of line (or strip transects) and optionally adjusts the counts based on detectability using a supplied detection function model. A generalized additive model, generalized mixed model or generalised is then used to model these adjusted counts based on a formula involving environmental covariates.

Usage

```
dsm(formula, ddf.obj, segment.data, observation.data, engine = "gam",
     convert.units = 1, family = quasipoisson(link = "log"), group = FALSE,
     gamma = 1.4, control = list(keepData = TRUE), availability = 1,
     strip.width = NULL, segment.area = NULL, weights = NULL, ...)
```

Arguments

formula formula for the surface. This should be a valid [glm/gam/gamm](#) formula. See "Details", below, for how to define the response.
ddf.obj result from call to [ddf](#) or [ds](#). If ddf.obj is NULL then strip transects are assumed.
segment.data segment data, see [dsm-data](#).
observation.data observation data, see [dsm-data](#).
engine which fitting engine should be used for the DSM ([glm/gam/gamm/bam](#)).

<code>convert.units</code>	value to alter length or width for calculation of the offset, applied to ‘segment.area’ if used.
<code>family</code>	response distribution (popular choices include quasipoisson , Tweedie and negbin). Defaults to <code>quasipoisson</code> .
<code>group</code>	if TRUE the abundance of groups will be calculated rather than the abundance of individuals. Setting this option to TRUE is equivalent to setting the size of each group to be 1.
<code>control</code>	the usual <code>control</code> argument for a <code>gam</code> ; <code>keepData</code> must be TRUE for variance estimation to work.
<code>availability</code>	an availability bias used to scale the counts/estimated counts by. If we have N animals in a segment, then <code>N/availability</code> will be entered into the model. Uncertainty in the availability is not handled at present.
<code>gamma</code>	parameter to <code>gam()</code> set to a value of 1.4 (from advice in Wood (2006)) such that the <code>gam()</code> is inclined to not ‘overfit’ when GCV is used to select the smoothing parameter (ignored for REML, see <code>link{gam}</code> for further details).
<code>strip.width</code>	if <code>ddf.obj</code> , above, is NULL, then this is where the strip width is specified (i.e. for a strip transect survey). This is sometimes (and more correctly) referred to as the half-width, i.e. right truncation minus left truncation.
<code>segment.area</code>	if ‘NULL’ (default) segment areas will be calculated by multiplying the ‘Effort’ column in ‘segment.data’ by the (right minus left) truncation distance for the ‘ddf.obj’ or by ‘strip.width’. Alternatively a vector of segment areas can be provided (which must be the same length as the number of rows in ‘segment.data’) or a character string giving the name of a column in ‘segment.data’ which contains the areas. If <code>segment.area</code> is specified it takes precedent.
<code>weights</code>	weights for each observation used in model fitting. The default, <code>weights=NULL</code> , weights each observation by its area (see Details). Setting a scalar value (e.g. <code>weights=1</code>) all observations are equally weighted.
<code>...</code>	anything else to be passed straight to <code>glm/gam/gamm/bam</code> .

Details

The response (LHS of ‘formula’) can be one of the following:

<code>n</code> , <code>count</code> , <code>N</code> , <code>abundance</code>	count in each segment
<code>Nhat</code> , <code>abundance.est</code>	estimated abundance per segment, estimation is via a Horvitz-Thompson estimator. This should be used for density estimation.
<code>presence</code>	interpret the data as presence/absence (remember to change the family argument to <code>binomial</code>)
<code>D</code> , <code>density</code> , <code>Dhat</code> , <code>density.est</code>	density per segment

The offset used in the model is dependent on the response:

<code>count</code>	area of segment multiplied by average probability of detection in the segment
<code>estimated count</code>	area of the segment
<code>presence</code>	zero
<code>density</code>	zero

In the latter two cases (density and presence estimation) observations can be weighted by segment areas via the `weights=` argument. By default (`weights=NULL`), when density or presence are estimated the weights are set to the segment areas (using `segment.area` or by calculating $2 * (\text{strip width}) * \text{Effort}$) Alternatively `weights=1` will set the weights to all be equal. A third alternative is to pass in a vector of length equal to the number of segments, containing appropriate weights.

Value

a `glm/gam/gamm` object, with an additional element, `ddf` which holds the detection function object.

Large models

For large models, `engine="bam"` with `method="fREML"` may be useful. Models specified for `bam` should be as `gam`. READ `bam` before using this option; this option is considered EXPERIMENTAL at the moment. In particular note that the default basis choice (thin plate regression splines) will be slow and that in general fitting is less stable than when using `gam`. For negative binomial response, `theta` must be specified when using `bam`.

Author(s)

David L. Miller

References

Hedley, S. and S. T. Buckland. 2004. Spatial models for line transect sampling. *JABES* 9:181-199.

Miller, D. L., Burt, M. L., Rexstad, E. A., Thomas, L. (2013), Spatial models for distance sampling data: recent developments and future directions. *Methods in Ecology and Evolution*, 4: 1001-1010. doi: 10.1111/2041-210X.12105 (Open Access, available at <http://onlinelibrary.wiley.com/doi/10.1111/2041-210X.12105/abstract>)

Wood, S.N. 2006. *Generalized Additive Models: An Introduction with R*. CRC/Chapman & Hall.

Examples

```
library(Distance)
library(dsm)

# load the Gulf of Mexico dolphin data (see ?mexdolphins)
data(mexdolphins)

# fit a detection function and look at the summary
hr.model <- ds(mexdolphins$distdata, max(mexdolphins$distdata$distance),
              key = "hr", adjustment = NULL)
summary(hr.model)

# fit a simple smooth of x and y
mod1 <- dsm(N~s(x,y), hr.model, mexdolphins$segdata, mexdolphins$obsdata)
summary(mod1)

# create an offset (in metres)
# each prediction cell is 444km2
```

```

off.set <- 444*1000*1000

# predict over a grid
mod1.pred <- predict(mod1, mexdolphins$preddata, off.set)

# calculate the predicted abundance over the grid
sum(mod1.pred)

# plot the smooth
plot(mod1)

```

 dsm-data

Data format for DSM

Description

Two data.frames must be provided to `dsm`. They are referred to as `observation.data` and `segment.data` (for observation and segment data, respectively).

Details

`observation.data` - the observation data frame must have the following columns:

<code>object</code>	unique object identifier
<code>Sample.Label</code>	the identifier for the segment that the observation occurred in
<code>size</code>	the size of each observed group (i.e. 1 for individuals)
<code>distance</code>	perpendicular/radial distance to observation

`segment.data` - the segment data frame must have the following columns:

<code>Effort</code>	the effort (in terms of length of the segment)
<code>Sample.Label</code>	identifier for the segment (unique!)
<code>???</code>	environmental covariates, for example x and y

 dsm.cor

Check for autocorrelation in residuals

Description

Once a DSM has been fitted to data, this function can be used to check for autocorrelation in the residuals.

Usage

```
dsm.cor(dsm.obj, Transect.Label = "Transect.Label",
```

```
Segment.Label = "Segment.Label", max.lag = 10,
resid.type = "scaled.pearson", fun = cor, ylim = c(0, 1),
subset = "all", ...)
```

Arguments

dsm.obj	a fitted dsm object.
Transect.Label	label for the transect (default: Transect.Label). Using different labels can be useful when transects are split over geographical features or when transects are surveyed multiple times.
Segment.Label	label for the segments (default: Segment.Label).The result of calling order() must make sense.
resid.type	the type of residuals used, see residuals.gam and residuals.gam . Defaults to "scaled.pearson" in the GAM case and "normalized" in the GAMM case (which are equivalent).
fun	the function to use, by default cor , must take two column vectors as arguments.
max.lag	maximum lag to calculate at.
ylim	user defined limits in y direction.
subset	which subset of the data should the correlation function be calculated on?
...	other options to pass to plot.

Value

a plot or a vector of fun applied at the lags.

Details

Within each Transect.Label, segments will be sorted according to their Segment.Labels. This may require some time to get right for your particular data. If one has multiple surveys where transects are revisited, for example, one may want to make Transect.Label a unique transect-survey id. Neither label need to be included in the model, they must just be present in the \$data field in the model. This usually means that they have to be in the segment data passed to dsm.

The current iteration of this function will only plot correlations nicely, other things are up to you but you can get the function to return the data (by assigning the result to an object).

If there are NA values in the residuals then the correlogram will not be calculated. This usually occurs due to NA values in the covariates (so the smoother will not have fitted values there). Code like 'any(is.na(dsm.obj\$data))' might be helpful.

Author(s)

David L. Miller

Examples

```
library(Distance)
library(dsm)

data(mexdolphins)

hr.model <- ds(mexdolphins$distdata, max(mexdolphins$distdata$distance),
              key = "hr", adjustment = NULL)
mod1<-dsm(N~s(x,y), hr.model, mexdolphins$segdata, mexdolphins$obsdata)

dsm.cor(mod1,resid.type="d",max.lag=9,Segment.Label="Sample.Label")
```

 dsm.var.gam

Variance estimation via Bayesian results

Description

Use results from the Bayesian interpretation of the GAM to obtain uncertainty estimates. See Wood (2006).

Usage

```
dsm.var.gam(dsm.obj, pred.data, off.set = NULL, seglen.varname = "Effort",
            type.pred = "response")
```

Arguments

dsm.obj	an object returned from running dsm .
pred.data	either: a single prediction grid or list of prediction grids. Each grid should be a data.frame with the same columns as the original data.
off.set	a a vector or list of vectors with as many elements as there are in pred.data. Each vector is as long as the number of rows in the corresponding element of pred.data. These give the area associated with each prediction point.
seglen.varname	name for the column which holds the segment length (default value "Effort").
type.pred	should the predictions be on the "response" or "link" scale? (default "response").

Details

This is based on [dsm.var.prop](#) by Mark Bravington and Sharon Hedley.

Value

a list with elements

model	the fitted model object
pred.var	covariances of the regions given in pred.data. Diagonal elements are the variances in order
bootstrap	logical, always FALSE

pred.data	as above
off.set	as above
model	the fitted model with the extra term
dsm.object	the original model, as above

Author(s)

David L. Miller

dsm.var.movblk	<i>Variance estimation via parametric moving block bootstrap</i>
----------------	--

Description

Estimate the variance in abundance over an area using a moving block bootstrap. Two procedures are implemented, one incorporating detection function uncertainty, one not.

Usage

```
dsm.var.movblk(dsm.object, pred.data, n.boot, block.size, off.set,
  ds.uncertainty = FALSE, samp.unit.name = "Transect.Label",
  progress.file = NULL, bs.file = NULL, bar = TRUE)
```

Arguments

dsm.object	object returned from dsm .
pred.data	a data.frame that holds prediction points, must have the correct columns for other environmental covariates. It also MUST have a column called cell.area which gives the area for each prediction cell
n.boot	number of bootstrap resamples.
block.size	number of segments in each block.
off.set	offset to be used in the model, see offsets Note that this should NOT be log()'d.
ds.uncertainty	incorporate uncertainty in the detection function? See Details, below. Note that this feature is EXPERIMENTAL at the moment.
samp.unit.name	name sampling unit to resample (default 'Transect.Label').
progress.file	path to a file to be used (usually by Distance) to generate a progress bar (default NULL – no file written).
bs.file	path to a file to store each bootstrap round. This stores all of the bootstrap results rather than just the summaries, enabling outliers to be detected and removed. (Default NULL).
bar	should a progress bar be printed to screen? (Default TRUE).

Details

Setting `ds.uncertainty=TRUE` will incorporate detection function uncertainty directly into the bootstrap. This is done by generating observations from the fitted detection function and then re-fitting a new detection function (of the same form), then calculating a new effective strip width. Rejection sampling is used to generate the observations (except in the half-normal case) so the procedure can be rather slow. Note that this is currently not supported with covariates in the detection function.

Setting `ds.uncertainty=FALSE` will incorporate detection function uncertainty using the delta method. This assumes that the detection function and the spatial model are INDEPENDENT. This is probably not reasonable.

Examples

```
## Not run:
library(Distance)
library(dsm)

# load the Gulf of Mexico dolphin data (see ?mexdolphins)
data(mexdolphins)

# fit a detection function and look at the summary
hr.model <- ds(mexdolphins$distdata, max(mexdolphins$distdata$distance),
              key = "hr", adjustment = NULL)
summary(hr.model)

# fit a simple smooth of x and y
mod1<-dsm(N~s(x,y), hr.model, mexdolphins$segdata, mexdolphins$obsdata)
summary(mod1)

# create an offset (in metres)
# each prediction cell is 444km2
off.set <- 444*1000*1000

# calculate the variance by 500 moving block bootstraps
mod1.movblk <- dsm.var.movblk(mod1, mexdolphins$preddata, n.boot = 500,
                             block.size = 3, samp.unit.name = "Transect.Label", off.set = off.set,
                             bar = TRUE, bs.file = "mexico-bs.csv", ds.uncertainty = TRUE)

## End(Not run)
```

dsm.var.prop

Variance propagation for DSM models

Description

Rather than use a bootstrap to calculate the variance in a dsm model, use the clever variance propagation trick from Williams et al. (2011).

Usage

```
dsm.var.prop(dsm.obj, pred.data, off.set, seglen.varname = "Effort",
             type.pred = "response")
```

Arguments

dsm.obj	an object returned from running <code>dsm</code> .
pred.data	either: a single prediction grid or list of prediction grids. Each grid should be a <code>data.frame</code> with the same columns as the original data.
off.set	a a vector or list of vectors with as many elements as there are in <code>pred.data</code> . Each vector is as long as the number of rows in the corresponding element of <code>pred.data</code> . These give the area associated with each prediction point.
seglen.varname	name for the column which holds the segment length (default value "Effort").
type.pred	should the predictions be on the "response" or "link" scale? (default "response").

Details

The idea is to refit the spatial model but including the Hessian of the offset as an extra term. Variance estimates using this new model can then be used to calculate the variance of abundance estimates which incorporate detection function uncertainty. Further mathematical details are given in the paper in the references below.

Many prediction grids can be supplied by supplying a list of `data.frames` to the function.

Note that this routine is only useful if a detection function has been used in the DSM.

Based on (much more general) code from Mark Bravington and Sharon Hedley.

Value

a list with elements

model	the fitted model object
pred.var	covariances of the regions given in <code>pred.data</code> . Diagonal elements are the variances in order
bootstrap	logical, always FALSE
pred.data	as above
off.set	as above
model	the fitted model with the extra term
dsm.object	the original model, as above
model.check	simple check of subtracting the coefficients of the two models to see if there is a large difference
deriv	numerically calculated Hessian of the offset

Author(s)

Mark V. Bravington, Sharon L. Hedley. Bugs added by David L. Miller.

References

Williams, R., Hedley, S.L., Branch, T.A., Bravington, M.V., Zerbini, A.N. and Findlay, K.P. (2011). Chilean Blue Whales as a Case Study to Illustrate Methods to Estimate Abundance and Evaluate Conservation Status of Rare Species. *Conservation Biology* 25(3), 526-535.

Examples

```
## Not run:
library(Distance)
library(dsm)

# load the Gulf of Mexico dolphin data (see ?mexdolphins)
data(mexdolphins)

# fit a detection function and look at the summary
hr.model <- ds(mexdolphins$distdata, max(mexdolphins$distdata$distance),
              key = "hr", adjustment = NULL)
summary(hr.model)

# fit a simple smooth of x and y
mod1 <- dsm(N~s(x,y), hr.model, mexdolphins$segdata, mexdolphins$obsdata)

# Calculate the offset...
off.set <- 444*1000*1000

# Calculate the variance
mod1.var <- dsm.var.prop(mod1, mexdolphins$pred, off.set)

# summary over the whole area in mexdolphins$pred

# Plot a map of the CV
# need to format the prediction data with split
mod1.var.map <- dsm.var.prop(mod1,
                             split(mexdolphins$pred, 1:nrow(mexdolphins$pred)), off.set)
plot(mod1.var.map)

## End(Not run)
```

```
generate.ds.uncertainty
```

Generate data from a fitted detection function

Description

When ds.uncertainty is TRUE, this procedure generates data from the fitted detection function (assuming that it is correct).

Usage

```
generate.ds.uncertainty(ds.object)
```

Arguments

ds.object a fitted detection function object (as returned by a call to `ddf.ds()`).

Note

This function changes the random number generator seed. To avoid any potential side-effects, use something like: `seed <- get(".Random.seed",envir=.GlobalEnv) ### Run some code assign(".Random.seed",seed,envir=.GlobalEnv)`

Author(s)

David L. Miller

generate.mb.sample *Moving block bootstrap sampler*

Description

Not usually used on its own, called from within `dsm.var.movblk`.

Usage

```
generate.mb.sample(num.blocks.required, block.size, which.blocks, dsm.data,  
                  unit.info, n.units)
```

Arguments

num.blocks.required number of blocks that we need.
block.size number of segments per block.
which.blocks which blocks should be sampled.
dsm.data the \$data element of the result of a call to `dsm`.
unit.info result of calling `block.info.per.su`.
n.units number of sampling units.

Value

vector of log-residuals

latlong2km	<i>Convert latitude and longitude to Northings and Eastings</i>
------------	---

Description

Convert longitude and latitude co-ordinates to kilometres west-east and south-north from axes through (lon0,lat0) using the "spherical law of cosines".

Usage

```
latlong2km(lon, lat, lon0 = sum(range(lon))/2, lat0 = sum(range(lat))/2)
```

Arguments

lon	longitude
lat	latitude
lon0	longitude reference point (defaults to mean longitude)
lat0	latitude reference point (defaults to mean latitude)

Details

WARNING: This is an approximate procedure for converting between latitude/ longitude and Northing/Easting. Consider using projection conversions available in packages `sp` and `rgdal` for better results.

Value

list with elements `km.e` and `km.n`.

Author(s)

Simon N. Wood

make.soapgrid	<i>Create a knot grid for the internal part of a soap film smoother.</i>
---------------	--

Description

This routine simply creates a grid of knots (in the correct format) to be used as in the "internal" part of the soap film smoother

Usage

```
make.soapgrid(bnd, n.grid)
```

Arguments

- bnd list with elements x and y which give the locations of the boundary vertices. The first and last elements should be the same.
- n.grid either one number giving the number of points along the x and y axes that should be used to create the grid, or a vector giving the number in the x direction, then y direction.

Value

a list with elements x and y, containing the knot locations.

Author(s)

David L Miller

matrixnotposdef.handler

Handler to suppress the "matrix not positive definite" warning

Description

Internal function to suppress an annoying warnings from chol()

Usage

```
matrixnotposdef.handler(w)
```

Arguments

w a warning

Details

See: <https://stat.ethz.ch/pipermail/r-help/2012-February/302407.html> See: <http://romainfrancois.blog.free.fr/index.php?post/specific-warnings>

Value

not a warning if the warning was "matrix not positive definite" or "the matrix is either rank-deficient or indefinite"

Author(s)

David L. Miller

mexdolphins

*Pan-tropical spotted dolphins***Description**

Data from a combination of several NOAA shipboard surveys conducted on pan-tropical spotted dolphins in the Gulf of Mexico. 47 observations of groups of dolphins The group size was recorded, as well as the Beaufort sea state at the time of the observation. Coordinates for each observation and bathymetry data were also available as covariates for the analysis. A complete example analysis is provided at <https://github.com/dill/dsm/wiki/mexico-analysis>.

Format

List of 5 \$ obsdata : 'data.frame': 47 obs. of 5 variables: ..\$ object : int [1:47] 45 61 63 85 114 120 146 165 187 214\$ Sample.Label: Factor w/ 387 levels "19960417-1","19960417-2",...: 45 61 63 85 114 120 146 165 187 214\$ size : num [1:47] 21 150 125 75 50 45 40 220 100 35\$ distance : num [1:47] 3297 929 6051 5500 7259\$ Effort : num [1:47] 36300 17800 21000 21800 13400 20900 20600 21700 29400 22900 ... \$ preddata : 'data.frame': 1374 obs. of 7 variables: ..\$ latitude : num [1:1374] 30.1 30.1 30.1 30.1 30.1\$ longitude: num [1:1374] -87.6 -87.4 -87.2 -87.1 -86.9\$ depth : num [1:1374] 35 30 27 22 46 14 38 58 100 156\$ x : num [1:1374] 70832 86868 102904 118940 134976\$ y : num [1:1374] 341079 341079 341079 341079 341079\$ width : num [1:1374] 32072 32072 32072 32072 32072\$ height : num [1:1374] 37065 37065 37065 37065 37065 ... \$ segdata : 'data.frame': 387 obs. of 8 variables: ..\$ latitude : num [1:387] 29.9 29.8 29.8 29.7 29.6\$ longitude : num [1:387] -86.9 -86.8 -86.7 -86.7 -86.6\$ Effort : num [1:387] 13800 14000 14000 13900 13800 13800 14000 14000 13900 14800\$ Transect.Label: Factor w/ 45 levels "19960417","19960418",...: 1 1 1 1 1 1 1 1 1 2\$ Sample.Label : Factor w/ 387 levels "19960417-1","19960417-2",...: 1 2 3 4 5 6 7 8 9 10\$ depth : num [1:387] 135 148 152 164 180\$ x : num [1:387] 134159 143496 152050 161102 169553\$ y : num [1:387] 325561 314055 304324 293475 282984 ... \$ distdata : 'data.frame': 47 obs. of 10 variables: ..\$ object : int [1:47] 45 61 63 85 114 120 146 165 187 214\$ size : num [1:47] 21 150 125 75 50 45 40 220 100 35\$ distance : num [1:47] 3297 929 6051 5500 7259\$ Effort : num [1:47] 36300 17800 21000 21800 13400 20900 20600 21700 29400 22900\$ detected : num [1:47] 1 1 1 1 1 1 1 1 1 1\$ beaufort : num [1:47] 4 4 2 1 3 5 1 5 2 5\$ latitude : num [1:47] 27.7 26 26 27.5 27.4\$ longitude: num [1:47] -86 -87.6 -87.9 -90.4 -95\$ x : num [1:47] 228139 69199 37046 -210016 -658878\$ y : num [1:47] 79258 -113083 -112197 54208 43337 ... \$ survey.area: 'data.frame': 45 obs. of 2 variables: ..\$ longitude: num [1:45] -87.4 -94.9 -96.3 -96.5 -96.6\$ latitude : num [1:45] 25.8 25.8 25.8 26.2 27.2 ...

References

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offsets

Offsets

Description

This will be documentation on calculating offsets.

plot.dsm

Plot a density surface model.

Description

See [plot.gam](#).

Usage

```
## S3 method for class 'dsm'  
plot(x, ...)
```

Arguments

x a dsm object
... other arguments passed to [plot.gam](#).

Value

a plot!

Author(s)

David L. Miller

See Also

dsm [plot.gam](#)

plot.dsm.var *Create plots of abundance uncertainty*

Description

Note that the prediction data set must have x and y columns even if these were not used in the model.

Usage

```
## S3 method for class 'dsm.var'
plot(x, poly = NULL, limits = NULL, breaks = NULL,
     legend.breaks = NULL, xlab = "x", ylab = "y", observations = TRUE,
     plot = TRUE, boxplot.coef = 1.5, x.name = "x", y.name = "y",
     gg.grad = NULL, ...)
```

Arguments

x	a dsm.var object
poly	a list or data.frame with columns x and y, which gives the coordinates of a polygon to draw. It may also optionally have a column group, if there are many polygons.
limits	limits for the fill colours
breaks	breaks for the colour fill
legend.breaks	breaks as they should be displayed
xlab	label for the x axis
ylab	label for the y axis
observations	should observations be plotted?
plot	actually plot the map, or just return a ggplot2 object?
boxplot.coef	control trimming (as in summary.dsm.var), only has an effect if the bootstrap file was saved.
x.name	name of the variable to plot as the x axis.
y.name	name of the variable to plot as the y axis.
gg.grad	optional ggplot gradient object.
...	any other arguments

Value

a plot

Details

In order to get plotting to work with `dsm.var.prop` and `dsm.var.gam`, one must first format the data correctly since these functions are designed to compute very general summaries. One summary is calculated for each element of the list `pred` supplied to `dsm.var.prop` and `dsm.var.gam`.

For a plot of uncertainty over a prediction grid, `pred` (a `data.frame`), say, we can create the correct format by simply using `pred.new <- split(pred, 1:nrow(pred))`.

Author(s)

David L. Miller

See Also

`dsm.var.prop`, `dsm.var.gam`, `dsm.var.movblk`

predict.dsm

Predict from a fitted density surface model

Description

Make predictions outside (or inside) the covered area.

Usage

```
## S3 method for class 'dsm'
predict(object, newdata = NULL, off.set = NULL,
        type = "response", ...)
```

Arguments

<code>object</code>	a fitted <code>dsm</code> object as produced by <code>dsm()</code> .
<code>newdata</code>	spatially referenced covariates e.g. altitude, depth, distance to shore, etc. Note covariates in this dataframe must have names <i>identical</i> to variable names used in fitting the DSM.
<code>off.set</code>	area of each of the cells in the prediction grid. Ignored if there is already a column in <code>newdata</code> called <code>off.set</code> .
<code>type</code>	what scale should the results be on. The default is "response", see <code>predict.gam</code> for an explanation of other options (usually not necessary).
<code>...</code>	any other arguments passed to <code>predict.gam</code> .

Value

predicted values on the response scale (density/abundance).

Author(s)

David L. Miller

print.dsm	<i>Print a description of a density surface model object</i>
-----------	--

Description

This method just gives a short description of the fitted model. Use the [summary.dsm](#) method for more information.

Usage

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

Arguments

x	a dsm object
...	unspecified and unused arguments for S3 consistency

Author(s)

David L. Miller

See Also

[summary.ds](#)

print.dsm.var	<i>Print a description of a density surface model variance object</i>
---------------	---

Description

This method only provides a short summary, use the [summary.dsm.var](#) method for information.

Usage

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

Arguments

x	a dsm variance object
...	unspecified and unused arguments for S3 consistency

Author(s)

David L. Miller

See Also

[summary.dsm.var](#)

`print.summary.dsm.var` *Print summary of density surface model variance object*

Description

See [summary.dsm.var](#) for information.

Usage

```
## S3 method for class 'summary.dsm.var'
print(x, ...)
```

Arguments

`x` a summary of dsm variance object
`...` unspecified and unused arguments for S3 consistency

Author(s)

David L. Miller

See Also

[summary.dsm.var](#)

`rqgam.check` *Randomised quantile residuals check plots for GAMs/DSMs*

Description

Function operates as [gam.check](#) but using randomised quantile residuals, a la Dunn and Smyth (1996). Checks of k are not computed, these need to be done using [gam.check](#).

Usage

```
rqgam.check(gam.obj, ...)
```

Arguments

`gam.obj` a gam, glm or dsm object.
`...` arguments passed on to all plotting function

Value

just plots!

Author(s)

Based on code provided by Natalie Kelly, bugs added by Dave Miller

Examples

```
library(Distance)
library(dsm)

# load the Gulf of Mexico dolphin data (see ?mexdolphins)
data(mexdolphins)

# fit a detection function and look at the summary
hr.model <- ds(mexdolphins$distdata, max(mexdolphins$distdata$distance),
              key = "hr", adjustment = NULL)

# fit a simple smooth of x and y
mod1 <- dsm(N~s(x,y), hr.model, mexdolphins$segdata, mexdolphins$obsdata)
rqgam.check(mod1)
```

summary.dsm

Summarize a fitted density surface model

Description

Gives a brief summary of a fitted dsm object.

Usage

```
## S3 method for class 'dsm'
summary(object, ...)
```

Arguments

object a dsm object
 ... other arguments passed to `summary.gam`.

Value

a summary object

Author(s)

David L. Miller

See Also

dsm

`summary.dsm.var`*Summarize the variance of a density surface model*

Description

Gives a brief summary of a fitted dsm variance object.

Usage

```
## S3 method for class 'dsm.var'  
summary(object, alpha = 0.05, boxplot.coef = 1.5,  
        bootstrap.subregions = NULL, ...)
```

Arguments

<code>object</code>	a <code>dsm.var</code> object
<code>alpha</code>	alpha level for confidence intervals
<code>boxplot.coef</code>	the value of <code>coef</code> used to calculate the outliers see boxplot .
<code>bootstrap.subregions</code>	list of vectors of logicals or indices for subregions for which variances need to be calculated (only for bootstraps (see dsm.var.prop for how to use subregions with variance propagation).
<code>...</code>	unused arguments for S3 compatibility

Value

a summary object

Author(s)

David L. Miller

See Also

`dsm.var.movblk` `dsm.var.prop`

trim.var	<i>Trimmed variance</i>
----------	-------------------------

Description

Trim the variance estimates from the bootstrap. This is defined as the percentage defined as amount necessary to bring median and trimmed mean within 8

Usage

```
trim.var(untrimmed.bootstraps, boxplot.coef = 1.5)
```

Arguments

untrimmed.bootstraps

(usually the `$study.area.total` element of a returned dsm bootstrap object.

boxplot.coef the value of coef used to calculate the outliers see [boxplot](#).

Details

Code originally by Louise Burt.

Value

trimmed variance

Author(s)

Louise Burt

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