

Package ‘dsm’

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Title Density Surface Modelling of Distance Sampling Data

LazyLoad yes

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Description Density surface modelling of line transect data. A Generalized Additive Model-based approach is used to calculate spatially-explicit estimates of animal abundance from distance sampling (also presence/absence and strip transect) data. Several utility functions are provided for model checking, plotting and variance estimation.

Version 2.2.17

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

BugReports <https://github.com/DistanceDevelopment/dsm/issues>

Depends R (>= 3.0), mgcv (>= 1.8-23), mrds (>= 2.1.16), numDeriv

Imports nlme, ggplot2, plyr, statmod

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 dsm-package

Density surface modelling

Description

dsm implements spatial models for distance sampling data.

Details

Further information on distance sampling methods and example code is available at <http://distancesampling.org/R/>.

For help with distance sampling and this package, there is a Google Group <https://groups.google.com/forum/#!forum/distance-sampling>.

A example analysis is available at <http://distancesampling.org/R/vignettes/mexico-analysis.html>.

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 generalized linear model 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, control = list(keepData = TRUE), availability = 1,
     strip.width = NULL, segment.area = NULL, weights = NULL,
     transect = "line", method = "REML", ...)
```

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). |
| convert.units | conversion factor to multiply the area of the segments by. See 'Units' below. |
| family | response distribution (popular choices include quasipoisson , Tweedie/tw and negbin/nb). Defaults to quasipoisson . |
| group | 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. |
| control | the usual control argument for a gam; keepData must be TRUE for variance estimation to work (though this option cannot be set for GLMs or GAMMs). |
| availability | an availability bias used to scale the counts/estimated counts by. If we have N animals in a segment, then N/availability will be entered into the model. Uncertainty in the availability is not handled at present. |
| strip.width | if ddf.obj, 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. |
| segment.area | 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 segment.area is specified it takes precedent. |
| weights | weights for each observation used in model fitting. The default, weights=NULL, weights each observation by its area (see Details). Setting a scalar value (e.g. weights=1) all observations are equally weighted. |
| transect | type of transect ("line", the default or "point"). This is overridden by the detection function transect type, this is usually only necessary when no detection function is specified. |
| method | The smoothing parameter estimation method. Default is "REML", using Restricted Maximum Likelihood. See gam for other options. Ignored for engine="glm". |
| ... | anything else to be passed straight to glm/gam/gamm/bam . |

Details

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

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

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

| | |
|-----------------|---|
| count | area of segment multiplied by average probability of detection in the segment |
| estimated count | area of the segment |

| | |
|----------|------|
| presence | zero |
| density | 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.

Units

It is often the case that distances are collected in metres and segment lengths are recorded in kilometres. `dsm` allows you to provide a conversion factor (`convert.units`) to multiply the areas by. For example: if distances are in metres and segment lengths are in kilometres setting `convert.units=1000` will lead to the analysis being in metres. Setting `convert.units=1/1000` will lead to the analysis being in kilometres. The conversion factor will be applied to ‘`segment.area`’ if that is specified.

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

```
## 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(distdata, max(distdata$distance),
              key = "hr", adjustment = NULL)
summary(hr.model)

# fit a simple smooth of x and y to counts
mod1 <- dsm(counts(x,y), hr.model, segdata, obsdata)
summary(mod1)

# predict over a grid
mod1.pred <- predict(mod1, preddata, preddata$area)

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

# plot the smooth
plot(mod1)

## End(Not run)
```

 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`.

Details

The `segment.data` table has the sample identifiers which define the segments, the corresponding effort (line length) expended and the environmental covariates that will be used to model abundance/density. `observation.data` provides a link table between the observations used in the detection function and the samples (segments), so that we can aggregate the observations to the segments (i.e. `observation.data` is a "look-up table" between the observations and the segments).

`observation.data` - the observation data.frame must have (at least) 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 (e.g 1 if all animals occurred individually) |
| <code>distance</code> | distance to observation |

One can often also use `observation.data` to fit a detection function (so additional columns for detection function covariates are allowed in this table).

`segment.data`: the segment data.frame must have (at least) 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 location (projected latitude and longitude), and other relevant covariates |

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

| | |
|-----------------------------|--|
| <code>dsm.obj</code> | a fitted dsm object. |
| <code>Transect.Label</code> | label for the transect (default: <code>Transect.Label</code>). Using different labels can be useful when transects are split over geographical features or when transects are surveyed multiple times. |
| <code>Segment.Label</code> | label for the segments (default: <code>Segment.Label</code>).The result of calling <code>order()</code> must make sense. |
| <code>max.lag</code> | maximum lag to calculate at. |
| <code>resid.type</code> | 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). |
| <code>fun</code> | the function to use, by default <code>cor</code> , must take two column vectors as arguments. |
| <code>ylim</code> | user defined limits in y direction. |
| <code>subset</code> | which subset of the data should the correlation function be calculated on? |
| <code>...</code> | 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)

# load the data, see ?mexdolphins
data(mexdolphins)

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

# look at lag 1 differences up to a maximum of lag 9, using deviance
# residuals
dsm.cor(mod1, resid.type="deviance", max.lag=9,
        Segment.Label="Sample.Label")
```

dsm.var.gam

Prediction variance estimation assuming independence

Description

If one is willing to assume the the detection function and spatial model are independent, this function will produce estimates of variance of predictions of abundance, using the result that squared coefficients of variation will add.

Usage

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

Arguments

| | |
|----------------|--|
| dsm.obj | a model 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 cell. If a single number is supplied it will be replicated for the length of <code>pred.data</code> . |
| 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` taken from code by Mark Bravington and Sharon Hedley.

Value

a list with elements

| | |
|------------|---|
| model | the fitted model object |
| pred.var | variance of the regions given in <code>pred.data</code> . |
| bootstrap | logical, always FALSE |
| model | the fitted model with the extra term |
| dsm.object | the original model, as above |

Author(s)

David L. Miller

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(distdata, max(distdata$distance),
              key = "hr", adjustment = NULL)
summary(hr.model)

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

# Calculate the variance
# this will give a summary over the whole area in mexdolphins$preddata
mod1.var <- dsm.var.gam(mod1, preddata, off.set=preddata$area)
```

```
## End(Not run)
```

dsm.var.movblk *Variance estimation via parametric moving block bootstrap*

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 | 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. |
| n.boot | number of bootstrap resamples. |
| block.size | number of segments in each block. |
| 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 cell. If a single number is supplied it will be replicated for the length of <code>pred.data</code> . |
| 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 <code>Distance</code>) 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(distdata, max(distdata$distance),
              key = "hr", adjustment = NULL)
summary(hr.model)

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

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

## End(Not run)
```

dsm.var.prop

Prediction variance propagation for DSMs

Description

To ensure that uncertainty from the detection function is correctly propagated to the final variance estimate of abundance, this function uses a method first detailed in Williams et al (2011).

Usage

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

Arguments

| | |
|----------------|--|
| dsm.obj | a model 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 cell. If a single number is supplied it will be replicated for the length of <code>pred.data</code> . |
| 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 an extra random effect. This random effect has zero mean and hence to effect on point estimates. Its variance is the Hessian of the detection function. Variance estimates then 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 simply calls `dsm_varprop`. If you don't require multiple prediction grids, the other routine will probably be faster.

This routine is only useful if a detection function with covariates has been used in the DSM.

Value

a list with elements

| | |
|-------------|--|
| model | the fitted model object |
| pred.var | variance of each region given in <code>pred.data</code> . |
| 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 |

Diagnostics

The summary output from the function includes a simply diagnostic that shows the average probability of detection from the "original" fitted model (the model supplied to this function; column `Fitted.model`) and the probability of detection from the refitted model (used for variance propagation; column `Refitted.model`) along with the standard error of the probability of detection from the fitted model (`Fitted.model.se`), at the unique values of any factor covariates used in the detection function (for continuous covariates the 5

Limitations

Note that this routine is only useful if a detection function has been used in the DSM. It cannot be used when the `Nhat`, `abundance.est` responses are used. Importantly this requires that if the detection function has covariates, then these do not vary within a segment (so, for example covariates like sex cannot be used).

Negative binomial models fitted using the `nb` family will give strange results (overly big variance estimates due to scale parameter issues) so nb models are automatically refitted with `negbin` (with a warning). It is probably worth refitting these models with `negbin` manually (perhaps giving a smallish range of possible values for the negative binomial parameter) to check that convergence was reached.

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.

| | |
|-------------|--|
| dsm_varprop | <i>Variance propagation for density surface models</i> |
|-------------|--|

Description

Calculate the uncertainty in predictions from a fitted DSM, including uncertainty from the detection function.

Usage

```
dsm_varprop(model, newdata, trace = FALSE, var_type = "Vp")
```

Arguments

| | |
|----------|--|
| model | a fitted <code>dsm</code> |
| newdata | the prediction grid |
| trace | for debugging, see how the scale parameter estimation is going |
| var_type | which variance-covariance matrix should be used (" <code>Vp</code> " for variance-covariance conditional on smoothing parameter(s), " <code>Vc</code> " for unconditional). See <code>gamObject</code> for an details/explanation. If in doubt, stick with the default, " <code>Vp</code> ". |

Details

When we make predictions from a spatial model, we also want to know the uncertainty about that abundance estimate. Since density surface models are 2 (or more) stage models, we need to incorporate the uncertainty from the earlier stages (i.e. the detection function) into our "final" uncertainty estimate.

This function will refit the spatial model but include the Hessian of the offset as an extra term. Variance estimates using this new model can then be used to calculate the variance of predicted abundance estimates which incorporate detection function uncertainty. Importantly this requires that if the detection function has covariates, then these do not vary within a segment (so, for example covariates like sex cannot be used).

For more information on how to construct the prediction grid `data.frame`, `newdata`, see [predict.dsm](#).

This routine is only useful if a detection function with covariates has been used in the DSM.

Note that we can use `var_type="Vc"` here (see [gamObject](#)), which is the variance-covariance matrix for the spatial model, corrected for smoothing parameter uncertainty. See Wood, Pya & Svanbäck (2016) for more information.

Negative binomial models fitted using the `nb` family will give strange results (overly big variance estimates due to scale parameter issues) so `nb` models are automatically refitted with `negbin` (with a warning). It is probably worth refitting these models with `negbin` manually (perhaps giving a smallish range of possible values for the negative binomial parameter) to check that convergence was reached.

Value

a list with elements

| | |
|------------------------|---|
| <code>old_model</code> | fitted model supplied to the function as <code>model</code> |
| <code>refit</code> | refitted model object, with extra term |
| <code>pred</code> | point estimates of predictions at <code>newdata</code> |
| <code>var</code> | total variance calculated over all of <code>newdata</code> |
| <code>ses</code> | standard error for each prediction cell in <code>newdata</code> |

Diagnostics

The summary output from the function includes a simple diagnostic that shows the average probability of detection from the "original" fitted model (the model supplied to this function; column `Fitted.model`) and the probability of detection from the refitted model (used for variance propagation; column `Refitted.model`) along with the standard error of the probability of detection from the fitted model (`Fitted.model.se`), at the unique values of any factor covariates used in the detection function (for continuous covariates the 5

Author(s)

David L. Miller, based on code from Mark V. Bravington and Sharon L. Hedley.

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.
- Wood, S.N., Pya, N. and S"afken, B. (2016) Smoothing parameter and model selection for general smooth models. *Journal of the American Statistical Association*, 1-45.

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) before running code and assign(".Random.seed",seed,envir=.GlobalEnv) after.

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`*Create a knot grid for the internal part of a soap film smoother.*

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**

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

| | |
|----------------|-----------|
| <code>w</code> | 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 in the Gulf of Mexico

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 (and description of the data) is provided at <http://distancesampling.org/R/vignettes/mexico-analysis.html>.

References

Halpin, P.N., A.J. Read, E. Fujioka, B.D. Best, B. Donnelly, L.J. Hazen, C. Kot, K. Urian, E. LaBrecque, A. Dimatteo, J. Cleary, C. Good, L.B. Crowder, and K.D. Hyrenbach. 2009. OBIS-SEAMAP: The world data center for marine mammal, sea bird, and sea turtle distributions. *Oceanography* 22(2):104-115

NOAA Southeast Fisheries Science Center. 1996. Report of a Cetacean Survey of Oceanic and Selected Continental Shelf Waters of the Northern Gulf of Mexico aboard NOAA Ship Oregon II (Cruise 220)

obs_exp

Observed versus expected diagnostics for fitted DSMs

Description

Given a covariate, calculate the observed and expected counts for each unique value of the covariate. This can be a useful goodness of fit check for DSMs.

Usage

```
obs_exp(model, covar, cut = NULL)
```

Arguments

| | |
|-------|---|
| model | a fitted dsm model object |
| covar | covariate to aggregate by (character) |
| cut | vector of cut points to aggregate at. If not supplied, the unique values of covar are used. |

Details

One strategy for model checking is to calculate observed and expected counts at different aggregations of the variable. If these match well then the model fit is good.

Value

data.frame with values of observed and expected counts.

Author(s)

David L Miller, on the suggestion of Mark Bravington.

Examples

```
library(Distance)
library(dsm)

# example with the Gulf of Mexico dolphin data
data(mexdolphins)
hr.model <- ds(distdata, max(distdata$distance),
              key = "hr", adjustment = NULL)
mod1 <- dsm(count~s(x,y), hr.model, segdata, obsdata)
```

offsets

Offsets

Description

This will be documentation on calculating offsets.

| | |
|----------|--------------------------------------|
| plot.dsm | <i>Plot a density surface model.</i> |
|----------|--------------------------------------|

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 | <i>Create plots of abundance uncertainty</i> |
|--------------|--|

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 <code>dsm.var</code> object |
| poly | a <code>list</code> or <code>data.frame</code> with columns <code>x</code> and <code>y</code> , which gives the coordinates of a polygon to draw. It may also optionally have a column <code>group</code> , 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 <code>ggplot2</code> object? |
| boxplot.coef | control trimming (as in <code>summary.dsm.var</code>), 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 <code>ggplot</code> 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`

| | |
|-------------------|--|
| plot_pred_by_term | <i>Spatially plot predictions per model term</i> |
|-------------------|--|

Description

Plot the effect of each smooth in the model spatially. For each term in the model, plot its effect in space. Plots are made on the same scale, so that the relative influence of each smooth can be seen.

Usage

```
plot_pred_by_term(dsm.obj, data, location_cov = c("x", "y"))
```

Arguments

| | |
|--------------|--|
| dsm.obj | fitted dsm object |
| data | data to use to plot (often the same as the precision grid), data should also include width and height columns for plotting |
| location_cov | which covariates to plot by (usually 2, spatial covariates, by default =c("x", "y")) |

Value

a ggplot2 plot

Author(s)

David L Miller (idea taken from inlabru)

Examples

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

# load the Gulf of Mexico dolphin data and fit a model
data(mexdolphins)
hr.model <- ds(distdata, max(distdata$distance),
              key = "hr", adjustment = NULL)
mod1 <- dsm(count~s(x,y) + s(depth), hr.model, segdata, obsdata)

preddata$width <- preddata$height <- sqrt(preddata$area)

# make the plot
plot_pred_by_term(mod1, preddata, c("x", "y"))

# better plot would be
# library(viridis)
# plot_pred_by_term(mod1, preddata, c("x", "y")) + scale_fill_viridis()

## End(Not run)
```

`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 dsm object as produced by <code>dsm()</code> . |
| <code>newdata</code> | spatially referenced covariates e.g. altitude, depth, distance to shore, etc. Covariates in the <code>data.frame</code> 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. Should be in the same units as the segments/distances given to <code>dsm</code> . 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 predict.gam for an explanation of other options (usually not necessary). |
| <code>...</code> | any other arguments passed to predict.gam . |

Details

If `newdata` is not supplied, predictions are made for the data that built the model. Note that the order of the results will not necessarily be the same as the `segdata` (segment data) `data.frame` that was supplied (it will be sorted by the `Segment.Label` field).

Value

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

Author(s)

David L. Miller

See Also

[predict.gam](#) [dsm.var.gam](#) [dsm.var.prop](#) [dsm.var.movblk](#)

| | |
|-----------|--|
| 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](#)

| | |
|--------------------------------|---|
| <code>print.dsm_varprop</code> | <i>Print a description of a density surface model variance object</i> |
|--------------------------------|---|

Description

This method only provides a short summary, see [summary.dsm_varprop](#).

Usage

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

Arguments

| | |
|------------------|---|
| <code>x</code> | a dsm variance object |
| <code>...</code> | unspecified and unused arguments for S3 consistency |

Author(s)

David L. Miller

See Also[summary.dsm_varprop](#)

| | |
|------------------------------------|---|
| <code>print.summary.dsm.var</code> | <i>Print summary of density surface model variance object</i> |
|------------------------------------|---|

Description

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

Usage

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

Arguments

| | |
|------------------|---|
| <code>x</code> | a summary of dsm variance object |
| <code>...</code> | unspecified and unused arguments for S3 consistency |

Author(s)

David L. Miller

See Also[summary.dsm.var](#)

```
print.summary.dsm_varprop
```

Print summary of density surface model variance object

DescriptionSee [summary.dsm_varprop](#) for information.**Usage**

```
## S3 method for class 'summary.dsm_varprop'  
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 plot for GAMs/DSMs

Description

Reproduces the "Resids vs. linear pred" plot from [gam.check](#) but using randomised quantile residuals, a la Dunn and Smyth (1996). Checks for heteroskedasticity as usual, looking for "funnel"-type structures in the points, which is much easier with randomised quantile residuals than with deviance residuals, when your model uses a count distribution as the response.

Usage

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

Arguments

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

Details

Note that this function only works with negative binomial and Tweedie response distributions.

Earlier versions of this function produced the full gam.check output, but this was confusing as only one of the plots was really useful. Checks of k are not computed, these need to be done using [gam.check](#).

Value

just plots!

Author(s)

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

Examples

```
library(Distance)
library(dsm)
library(tweedie)

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

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

# fit a simple smooth of x and y with a Tweedie response with estimated
# p parameter
mod1 <- dsm(count~s(x, y), hr.model, segdata, obsdata, family=tw())
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 | <i>Summarize the variance of a density surface model</i> |
|-----------------|--|

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

object a dsm.var object
 alpha alpha level for confidence intervals (default 0.05 to give a 95% confidence interval, i.e. we generate $100 * c(\alpha/2, 1-\alpha/2)$ confidence intervals)
 boxplot.coef the value of coef used to calculate the outliers see [boxplot](#).
 bootstrap.subregions
 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).
 ... unused arguments for S3 compatibility

Value

a summary object

Author(s)

David L. Miller

See Also

dsm.var.movblk dsm.var.prop

summary.dsm_varprop *Summarize the variance of a density surface model*

Description

Gives a brief summary of a fitted dsm_varprop variance object.

Usage

```
## S3 method for class 'dsm_varprop'  
summary(object, alpha = 0.05, ...)
```

Arguments

| | |
|--------|---|
| object | a dsm.var object |
| alpha | alpha level for confidence intervals (default 0.05 to give a 95% confidence interval) |
| ... | unused arguments for S3 compatibility |

Value

a summary object

Author(s)

David L. Miller

See Also

dsm_varprop summary.dsm.var

| | |
|----------|-------------------------|
| 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](#).

Value

trimmed variance

Author(s)

Louise Burt

| | |
|----------------|--|
| vis.concurvity | <i>Visualise concurvity between terms in a GAM</i> |
|----------------|--|

Description

Plot measures of how much one term in the model could be explained by another. When values are high, one should consider re-running variable selection with one of the offending variables removed to check for stability in term selection.

Usage

```
vis.concurvity(model, type = "estimate")
```

Arguments

model fitted model
type concurvity measure to plot, see [concurvity](#)

Details

These methods are considered somewhat experimental at this time. Consult [concurvity](#) for more information on how concurvity measures are calculated.

Author(s)

David L Miller

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(distdata, max(distdata$distance),
              key = "hr", adjustment = NULL)

# fit a simple smooth of x and y to counts
mod1 <- dsm(count~s(x,y)+s(depth), hr.model, segdata, obsdata)

# visualise concurvity using the "estimate" metric
vis.concurvity(mod1)

## End(Not run)
```


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