## Package 'mrds’

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Title Mark-Recapture Distance Sampling
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Description Animal abundance estimation via conventional, multiple covariate and mark-recapture distance sampling (CDS/MCDS/MRDS). Detection function fitting is performed via maximum likelihood. Also included are diagnostics and plotting for fitted detection functions. Abundance estimation is via a Horvitz-Thompson-like estimator.
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mrds-package Mark-Recapture Distance Sampling (mrds)

## Description

This package implements mark-recapture distance sampling methods as described in D.L. Borchers, W. Zucchini and Fewster, R.M. (1988), "Mark-recapture models for line transect surveys", Biometrics 54: 1207-1220. and Laake, J.L. (1999) "Distance sampling with independent observers: Reducing bias from heterogeneity by weakening the conditional independence assumption." in Amstrup, G.W., Garner, S.C., Laake, J.L., Manly, B.F.J., McDonald, L.L. and Robertson, D.G. (eds) "Marine mammal survey and assessment methods", Balkema, Rotterdam: 137-148 and Borchers, D.L., Laake, J.L., Southwell, C. and Paxton, C.L.G. "Accommodating unmodelled heterogeneity in double-observer distance sampling surveys". 2006. Biometrics 62:372-378.)

## Details

Examples of distance sampling analyses are available at http://examples.distancesampling. org/.

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

## Author(s)

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add.df.covar.line Add covariate levels detection function plots

## Description

Add a line or lines to a plot of the detection function which correspond to a a given covariate combination. These can be particularly useful when there is a small number of factor levels or if quantiles of a continuous covariate are specified.

## Usage

add.df.covar.line(ddf, data, ndist $=250$, pdf = FALSE, breaks = "Sturges", ...)
add_df_covar_line(ddf, data, ndist = 250, pdf = FALSE, breaks = "Sturges", ...)

## Arguments

## ddf

data a data.frame with the covariate combination you want to plot.
ndist number of distances at which to evaluate the detection function.
pdf should the line be drawn on the probability density scale; ignored for line transects.
breaks required to ensure that PDF lines are the right size, should match what is supplied to original plot command. Defaults to "Sturges" breaks, as in hist. Only used if pdf=TRUE.
... extra arguments to give to line (lty, lwd, col).

## Details

All covariates must be specified in data. Plots can become quite busy when this approach is used. It may be useful to fix some covariates at their median level and plot set values of a covariate of interest. For example setting weather (e.g., Beaufort) to its median and plotting levels of observer, then creating a second plot for a fixed observer with levels of weather.
Arguments to lines are supplied in ... and aesthetics like line type (lty), line width (lwd) and colour (col) are recycled. By default lty is used to distinguish between the lines. It may be useful to add a legend to the plot (lines are plotted in the order of data).

## Value

invisibly, the values of detectability over the truncation range.

## Author(s)

David L Miller

## Examples

```
## Not run:
# fit an example model
data(book.tee.data)
egdata <- book.tee.data$book.tee.dataframe
result <- ddf(dsmodel = ~mcds(key = "hn", formula = ~sex),
    data = egdata[egdata$observer==1, ], method = "ds",
    meta.data = list(width = 4))
# make a base plot, showpoints=FALSE makes the plot less busy
plot(result, showpoints=FALSE)
# add lines for sex one at a time
add.df.covar.line(result, data.frame(sex=0), lty=2)
add.df.covar.line(result, data.frame(sex=1), lty=3)
# add a legend
legend(3, 1, c("Average", "sex==0", "sex==1"), lty=1:3)
# alternatively we can add both at once
# fixing line type and varying colour
plot(result, showpoints=FALSE)
add.df.covar.line(result, data.frame(sex=c (0,1)), lty=1,
    col=c("red", "green"))
# add a legend
legend(3, 1, c("Average", "sex==0", "sex==1"), lty=1,
    col=c("black", "red", "green"))
## End(Not run)
```

adj.check.order

Check order of adjustment terms

## Description

'adj.check.order' checks that the Cosine, Hermite or simple polynomials are of the correct order.

## Usage

adj.check.order(adj.series, adj.order, key)

## Arguments

adj.series Adjustment series used ('cos','herm','poly')
adj.order Integer to check
key
key function to be used with this adjustment series

## Details

Only even functions are allowed as adjustment terms, per p. 47 of Buckland et al (2001). If incorrect terms are supplied then an error is throw via stop.

## Value

Nothing! Just calls stop if something goes wrong.

## Author(s)

David Miller

## References

S.T.Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake. 1993. Robust Models. In: Distance Sampling, eds. S.T.Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake. Chapman \& Hall.

## See Also

adjfct.cos, adjfct. poly, adjfct.herm, detfct, mcds, cds

## Description

Extract the AIC from a fitted detection function.

## Usage

\#\# S3 method for class 'ddf'
AIC(object, ..., $k=2$ )

## Arguments

object a fitted detection function object
... optionally more fitted model objects.
$k \quad$ penalty per parameter to be used; the default $k=2$ is the "classical" AIC

## Author(s)

David L Miller
apex.gamma Get the apex for a gamma detection function

## Description

Get the apex for a gamma detection function

## Usage

apex.gamma(ddfobj)

## Arguments

ddfobj ddf object

## Value

the distance at which the gamma peaks

## Author(s)

Jeff Laake

```
assign.default.values Assign default values to list elements that have not been already as-
                        signed
```


## Description

Assigns default values for argument in list $\times$ from argument=value pairs in ...if $\times \$$ argument doesn't already exist

## Usage

assign.default.values(x, ...)

## Arguments

| $x$ | generic list |
| :--- | :--- |
| $\ldots$ | unspecified list of argument=value pairs that are used to assign values |

## Value

x - list with filled values

## Author(s)

Jeff Laake

## Description

For models with covariates the detection probability for each observation can vary. This function computes an average value for a set of distances to plot an average line to graphically represent the fitted model in plots that compare histograms and the scatter of individual estimated detection probabilities. Averages are calculated over the observed covariate combinations.

## Usage

average.line(finebr, obs, model)

## Arguments

| finebr | set of fine breaks in distance over which detection function values are averaged <br> and plotted |
| :--- | :--- |
| obs | value of observer for averaging (1-2 individual observers; 3 duplicates; 4 pooled <br> observation team $)$ |
| model | ddf model object |

## Value

list with 2 elements

$$
\begin{array}{ll}
\text { xgrid } & \text { vector of gridded distance values } \\
\text { values } & \text { vector of average detection function values at the xgrid values }
\end{array}
$$

## Note

Internal function called from plot functions for ddf objects

## Author(s)

Jeff Laake
average.line.cond Average conditional detection function line for plotting

## Description

For models with covariates the detection probability for each observation can vary. This function computes an average value for a set of distances to plot an average line to graphically represent the fitted model in plots that compare histograms and the scatter of individual estimated detection probabilities.

## Usage

average.line.cond(finebr, obs, model)

## Arguments

finebr set of fine breaks in distance over which detection function values are averaged and plotted
obs value of observer for averaging (1-2 individual observers)
model ddf model object

## Value

list with 2 elements:

$$
\begin{array}{ll}
\text { xgrid } & \text { vector of gridded distance values } \\
\text { values } & \text { vector of average detection function values at the xgrid values }
\end{array}
$$

## Note

Internal function called from plot functions for ddf objects

## Author(s)

Jeff Laake

book. tee. data $\quad$| Golf tee data used in chapter 6 of Advanced Distance Sampling exam- |
| :--- |
| ples |

## Description

Double platform data collected in a line transect survey of golf tees by 2 observers at St. Andrews. Field sex was actually colour of the golf tee: 0 - green; 1 - yellow. Exposure was either low (0) or high(1) depending on height of tee above the ground. size was the number of tees in an observed cluster.

## Format

The format is: List of $4 \$$ book.tee.dataframe:'data.frame': 324 obs. of 7 variables: .. $\$$ object : num [1:324] 1122334455 ... .. \$ observer: Factor w/ 2 levels "1","2": 1212121212 ... .. $\$$ detected: num [1:324] 1010101010 ... ..\$ distance: num [1:324] 2.682 .683 .333 .330 .340 .34 2.532 .531 .461 .46 ... .. \$ size : num [1:324] 2222112222 ... .. $\$$ sex : num [1:324] 11110 $01111 \ldots$.. $\$$ exposure: num [1:324] $1100001100 \ldots \$$ book.tee.region :'data.frame': 2 obs. of 2 variables: ..\$ Region.Label: Factor w/ 2 levels "1","2": 12 ..\$ Area : num [1:2] 1040640 \$ book.tee.samples :'data.frame': 11 obs. of 3 variables: ..\$ Sample.Label: num [1:11] 1234567

8910 ... .. \$ Region.Label: Factor w/ 2 levels "1","2": 1111112222 ... ..\$ Effort : num [1:11] 10303027211223231512 ... \$ book.tee.obs :'data.frame': 162 obs. of 3 variables: .. $\$$ object : int [1:162] 12321222324596061 ... .. $\$$ Region.Label: int [1:162] 1111111111 ... ..\$ Sample.Label: int [1:162] $1111111111 \ldots$

```
calc.se.Np Find se of average p and N
```


## Description

Find se of average p and N

## Usage

```
calc.se.Np(model, avgp, n, average.p)
```


## Arguments

| model | a ddf model object |
| :--- | :--- |
| avgp | average p function |
| n | sample size |
| average.p | the average probability of detection for the model |

## Author(s)

David L. Miller

```
cdf.ds
```

Cumulative distribution function (cdf) for fitted distance sampling detection function

## Description

Computes cdf values of observed distances from fitted distribution. For a set of observed x it returns the integral of $f(x)$ for the range $=($ inner, $x$ ), where inner is the innermost distance which is observable (either 0 or left if left truncated). In terms of $g(x)$ this is the integral of $g(x)$ over range divided by the integral of $g(x)$ over the entire range of the data (inner, W ).

## Usage

cdf.ds(model, newdata $=$ NULL)

## Arguments

model fitted distance sampling model
newdata new data values if computed for values other than the original observations

## Value

vector of cdf values for each observation

## Note

This is an internal function that is not intended to be invoked directly. It is called by qqplot.ddf to compute values for Kolmogorov-Smirnov and Cramer-von Mises tests and the Q-Q plot.

## Author(s)

Jeff Laake

## See Also

qqplot.ddf

```
cds CDS function definition
```


## Description

Creates model formula list for conventional distance sampling using values supplied in call to ddf

```
Usage
    cds(
        key = NULL,
        adj.series = NULL,
        adj.order = NULL,
        adj.scale = "width",
        adj.exp = FALSE,
        formula = ~1,
        shape.formula = ~1
    )
```


## Arguments

key string identifying key function (currently either "hn" (half-normal),"hr" (hazardrate), "unif" (uniform) or "gamma" (gamma distribution)
adj. series string identifying adjustment functions cos (Cosine), herm (Hermite polynomials), poly (simple polynomials) or NULL
adj. order vector of order of adjustment terms to include
adj.scale whether to scale the adjustment terms by "width" or "scale"
adj.exp if TRUE uses $\exp (\operatorname{adj})$ for adjustment to keep $f(x)>0$
formula formula for scale function (included for completeness only only formula=~1 for cds)
shape.formula formula for shape function

## Value

A formula list used to define the detection function model

| fct | string "cds" |
| :--- | :--- |
| key | key function string |
| adj.series | adjustment function string |
| adj.order | adjustment function orders |
| adj.scale | adjustment function scale type |
| formula | formula for scale function |
| shape.formula | formula for shape function |

## Author(s)

Jeff Laake; Dave Miller
check.bounds Check parameters bounds during optimisations

## Description

Simple internal function to check that the optimisation didn't hit bounds. Based on code that used to live in detfct.fit.opt.

## Usage

check.bounds(lt, lowerbounds, upperbounds, ddfobj, showit, setlower, setupper)

## Arguments

| lt | optimisation object |
| :--- | :--- |
| lowerbounds | current lower bounds |
| upperbounds | current upper bounds |
| ddfobj | ddf object |
| showit | debug level |
| setlower | were lower bounds set by the user |
| setupper | were upper bounds set by the user |

## Value

TRUE if parameters are close to the bound, else FALSE

## Author(s)

Dave Miller; Jeff Laake

## check.mono $\quad$ Check that a detection function is monotone

## Description

Check that a fitted detection function is monotone non-increasing.

## Usage

```
check.mono(
    df,
    strict = TRUE,
    n.pts = 100,
    tolerance = 1e-06,
    plot = FALSE,
    max.plots = 6
    )
```


## Arguments

df a fitted detection function object
strict if TRUE (default) the detection function must be "strictly" monotone, that is that $(g(x[i])<=g(x[i-1]))$ over the whole range (left to right truncation points).
n.pts number of equally-spaced points between left and right truncation at which to evaluate the detection function (default 100)
tolerance numerical tolerance for monotonicity checks (default 1e-6)
plot plot a diagnostic highlighting the non-monotonic areas (default FALSE)
max.plots when plot=TRUE, what is the maximum number of plots of non-monotone covariate combinations that should be plotted? Plotted combinations are a random sample of the non-monotonic subset of evaluations. No effect for non-covariate models.

## Details

Evaluates a series of points over the range of the detection function (left to right truncation) then determines:

1. If the detection function is always less than or equal to its value at the left truncation $(g(x)<=g(l e f t)$, or usually $g(x)<=g(0))$. 2. (Optionally) The detection function is always monotone decreasing $(g(x[i])<=g(x[i-1]))$. This check is only performed when strict=TRUE (the default). 3. The detection function is never less than $0(g(x)>=0)$. 4. The detection function is never greater than 1 ( $g(x)<=1$ ).

For models with covariates in the scale parameter of the detection function is evaluated at all observed covariate combinations.
Currently covariates in the shape parameter are not supported.
coef.ds

## Value

TRUE if the detection function is monotone, FALSE if it's not. warnings are issued to warn the user that the function is non-monotonic.

## Author(s)

David L. Miller

```
coef.ds Extract coefficients
```


## Description

Extract coefficients and provide a summary of parameters and estimates from the output of ddf model objects.

## Usage

```
\#\# S3 method for class 'ds'
    coef(object,...)
        \#\# S3 method for class 'io'
    coef(object,...)
            \#\# S3 method for class 'io.fi'
    coef(object,...)
            \#\# S3 method for class 'trial'
    coef(object,...)
            \#\# S3 method for class 'trial.fi'
    coef(object,...)
            \#\# S3 method for class 'rem'
    coef(object,...)
            \#\# S3 method for class 'rem.fi'
    coef(object,...)
```


## Arguments

object ddf model object of class ds, io, io.fi, trial, trial.fi, rem, or rem.fi.
... unspecified arguments that are unused at present

## Value

For coef.ds List of data frames for coefficients (scale and exponent (if hazard))
scale dataframe of scale coefficient estimates and standard errors
exponent dataframe with exponent estimate and standard error if hazard detection function
For all others Data frame containing each coefficient and standard error

Note
These functions are called by the generic function coef for any ddf model object. It can be called directly by the user, but it is typically safest to use coef which calls the appropriate function based on the type of model.

## Author(s)

Jeff Laake
compute.Nht Horvitz-Thompson estimates $1 / p \_i$ or $s_{-} i / p \_i$

## Description

Compute individual components of Horvitz-Thompson abundance estimate in covered region for a particular subset of the data depending on value of group = TRUE (do group abundance); FALSE (do individual abundance)

## Usage

compute.Nht(pdot, group = TRUE, size = NULL)

## Arguments

$$
\begin{array}{ll}
\text { pdot } & \text { vector of estimated detection probabilities } \\
\text { group } & \text { if TRUE (do group abundance); FALSE(do individual abundance) } \\
\text { size } & \text { vector of group size values for clustered populations }
\end{array}
$$

## Value

vector of $\mathrm{H}-\mathrm{T}$ components for abundance estimate

Note
Internal function called by covered.region. dht

## Author(s)

Jeff Laake

```
covered.region.dht
    Covered region estimate of abundance from Horvitz-Thompson-like
    estimator
```


## Description

Computes H-T abundance within covered region by sample.

## Usage

covered.region.dht(obs, samples, group)

## Arguments

| obs | observations table |
| :--- | :--- |
| samples | samples table |
| group | if TRUE compute abundance of group otherwise abundance of individuals |

## Value

Nhat.by.sample - dataframe of abundance by sample

## Note

Internal function called by dht and related functions

## Author(s)

Jeff Laake

## create.bins

Create bins from a set of binned distances and a set of cutpoints.

## Description

This is an internal routine and shouldn't be necessary in normal analyses.

## Usage

create.bins(data, cutpoints)

## Arguments

| data | 'data.frame' with at least the column 'distance'. |
| :--- | :--- |
| cutpoints | vector of cutpoints for the bins |

## Value

argument 'data' with two extra columns 'distbegin' and 'distend'.

## Author(s)

David L. Miller

```
create.model.frame Create a model frame for ddf fitting
```


## Description

Creates a model.frame for distance detection function fitting. It includes some pre-specified and computed variables with those included in the model specified by user (formula)

## Usage

create.model.frame(xmat, scale.formula, meta.data, shape.formula = NULL)

## Arguments

| xmat | dataframe for ddf |
| :--- | :--- |
| scale.formula | user specified formula for scale of distance detection function |
| meta.data | user-specified meta.data (see ddf |
| shape.formula | user specified formula for shape parameter of distance detection function |

## Details

The following fields are always included: detected, observer, binned, and optionally distance (unless null), timesdetected (if present in data). If the distance data were binned, include distbegin and distend point fields. If the integration width varies also include int.begin and int.end and include an offset field for an iterative glm, if used. Beyond these fields only fields used in the model formula are included.

## Value

model frame for analysis

## Note

Internal function and not called by user

## Author(s)

Jeff Laake
create.varstructure Creates structures needed to compute abundance and variance

## Description

Creates samples and obs dataframes used to compute abundance and its variance based on a structure of geographic regions and samples within each region. The intent is to generalize this routine to work with other sampling structures.

## Usage

create.varstructure(model, region, sample, obs, dht.se)

## Arguments

| model | fitted ddf object |
| :--- | :--- |
| region | region table |
| sample | sample table |
| obs | table of object \#'s and links to sample and region table |
| dht.se | is uncertainty going to be calculated later? |

## Details

The function performs the following tasks: 1)tests to make sure that region labels are unique, 2) merges sample and region tables into a samples table and issue a warning if not all samples were used, 3) if some regions have no samples or if some values of Area were not valid areas given then issue error and stop, then an error is given and the code stops, 4) creates a unique region/sample label in samples and in obs, 5) merges observations with sample and issues a warning if not all observations were used, 6) sorts regions by its label and merges the values with the predictions from the fitted model based on the object number and limits it to the data that is appropriate for the fitted detection function.

## Value

List with 2 elements:
samples merged dataframe containing region and sample info - one record per sample
obs merged observation data and links to region and samples

## Note

Internal function called by dht

## Author(s)

Jeff Laake

## Description

Generic function for fitting detection functions for distance sampling with single and double observer configurations. Independent observer, trial and dependent observer (removal) configurations are included. This is a generic function which does little other than to validate the calling arguments and methods and then calls the appropriate method specific function to do the analysis.

```
Usage
    ddf(
        dsmodel = call(),
        mrmodel \(=\) call(),
        data,
        method = "ds",
        meta.data \(=\) list(),
        control = list()
    )
```


## Arguments

dsmodel distance sampling model specification
mrmodel mark-recapture model specification
data dataframe containing data to be analyzed
method analysis method
meta.data list containing settings controlling data structure
control list containing settings controlling model fitting

## Details

The fitting code has certain expectations about data. It should be a dataframe with at least the following fields named and defined as follows:

$$
\begin{array}{ll}
\text { object } & \text { object number } \\
\text { observer } & \text { observer number (1 or } 2) \text { for double observer; only } 1 \text { if single observer } \\
\text { detected } & 1 \text { if detected by the observer and } 0 \text { if missed; always } 1 \text { for single observer } \\
\text { distance } & \text { perpendicular distance }
\end{array}
$$

If the data are for clustered objects, the dataframe should also contain a field named size that gives the observed number in the cluster. If the data are for a double observer survey, then there are two records for each observation and each should have the same object number. The code assumes the observations are listed in the same order for each observer such that if the data are subsetted by
observer there will be the same number of records in each and each subset will be in the same object order. In addition to these predefined and pre-named fields, the dataframe can have any number and type of fields that are used as covariates in the dsmodel and mrmodel. At present, discrepancies between observations in distance, size and any user-specified covariates cannot be assimilated into the uncertainty of the estimate. The code presumes the values for those fields are the same for both records (observer=1 and observer=2) and it uses the value from observer 1. Thus it makes sense to make the values the same for both records in each pair even when both detect the object or when observer 1 doesn't detect the object the data would have to be taken from observer 2 and would not be consistent.

Five different fitting methods are currently available and these in turn define whether dsmodel and mrmodel need to be defined.

| Method | Single/Double | dsmodel | mrmodel |
| :--- | :--- | :--- | :--- |
| ds | Single | yes | no |
| io | Double | yes | yes |
| io.fi | Double | no | yes |
| trial | Double | yes | yes |
| trial.fi | Double | no | yes |
| rem | Double | yes | yes |
| rem.fi | Double | no | yes |

Methods with the suffix ". fi" use the assumption of full independence and do not use the distance sampling portion of the likelihood which is why a dsmodel is not needed. An mrmodel is only needed for double observer surveys and thus is not needed for method ds.

The dsmodel specifies the detection function $\mathrm{g}(\mathrm{y})$ for the distance sampling data and the models restrict $g(0)=1$. For single observer data $g(y)$ is the detection function for the single observer and if it is a double observer survey it is the relative detection function (assuming $g(0)=1$ ) of both observers as a team (the unique observations from both observers). In double observer surveys, the detection function is $p(y)=p(0) g(y)$ such that $p(0)<1$. The detection function $g(y)$ is specified by dsmodel and $p(0)$ estimated from the conditional detection functions (see mrmodel below). The value of dsmodel is specified using a hybrid formula/function notation. The model definition is prefixed with $\mathrm{a} \sim$ and the remainder is a function definition with specified arguments. At present there are two different functions, cds and mcds, for conventional distance sampling and multi-covariate distance sampling. Both functions have the same required arguments (key,formula). The first specifies the key function this can be half-normal ("hn"), hazard-rate ("hr"), gamma ("gamma") or uniform ("unif"). The argument formula specifies the formula for the log of the scale parameter of the key function (e.g., the equivalent of the standard deviation in the half-normal). The variable distance should not be included in the formula because the scale is for distance. See Marques, F.F.C. and S.T. Buckland (2004) for more details on the representation of the scale formula. For the hazard rate and gamma functions, an additional shape.formula can be specified for the model of the shape parameter. The default will be $\sim 1$. Adjustment terms can be specified by setting adj. series which can have the values: "none", "cos" (cosine), "poly" (polynomials), and "herm" (Hermite polynomials). One must also specify a vector of orders for the adjustment terms (adj.order) and a scaling (adj. scale) which may be "width" or "scale" (for scaling by the scale parameter). Note that the uniform key can only be used with adjustments (usually cosine adjustments for a Fourier-type analysis).

The mrmodel specifies the form of the conditional detection functions (i.e.,probability it is seen by observer j given it was seen by observer $3-\mathrm{j}$ ) for each observer ( $\mathrm{j}=1,2$ ) in a double observer survey. The value is specified using the same mix of formula/function notation but in this case the functions are glm and gam. The arguments for the functions are formula and link. At present, only glm is allowed and it is restricted to link=logit. Thus, currently the only form for the conditional detection functions is logistic as expressed in eq 6.32 of Laake and Borchers (2004). In contrast to dsmodel, the argument formula will typically include distance and all other covariates that affect detection probability. For example, mrmodel=~glm(formula=~distance+size+sex) constructs a conditional detection function based on the logistic form with additive factors, distance, size, and sex. As another example, mrmodel=~glm(formula=~distance*size+sex) constructs the same model with an added interaction between distance and size.
The argument meta. data is a list that enables various options about the data to be set. These options include:
point if TRUE the data are from point counts and FALSE (default) implies line transect data
width distance specifying half-width of the transect
left distance specifying inner truncation value
binned TRUE or FALSE to specify whether distances should be binned for analysis
breaks if binned=TRUE, this is a required sequence of break points that are used for plotting/gof. They should match distbegin, distend values if bins are fixed
int. range an integration range for detection probability; either a vector of 2 or matrix with 2 columns
mono constrain the detection function to be weakly monotonically decreasing (only applicable when there are no covariates in the detection function)
mono.strict when TRUE constrain the detection function to be strictly monotonically decreasing (again, only applicable when there are no covariates in the detection function)

Using meta. data=list(int. range $=c(1,10)$ ) is the same as meta. data=list(left=1, width=10). If meta.data=list (binned=TRUE) is used, the dataframe needs to contain the fields distbegin and distend for each observation which specify the left and right hand end points of the distance interval containing the observation. This is a general data structure that allows the intervals to change rather than being fixed as in the standard distance analysis tools. Typically, if the intervals are changing so is the integration range. For example, assume that distance bins are generated using fixed angular measurements from an aircraft in which the altitude is varying. Because all analyses are truncated (i.e., the last interval does not go to infinity), the transect width (and the left truncation point if there is a blindspot below the aircraft) can potentially change for each observation. The argument int. range can also be entered as a matrix with 2 columns (left and width) and a row for each observation.

The argument control is a list that enables various analysis options to be set. It is not necessary to set any of these for most analyses. They were provided so the user can optionally see intermediate fitting output and to control fitting if the algorithm doesn't converge which happens infrequently. The list values include:
showit Integer ( $0-3$, default 0 ) controls the (increasing)amount of information printed during fitting. 0 - none, $>=1$ - information about refitting and bound changes is printed, $>=2$ - information about adjustment term fitting is printed, $==3$-per-iteration parameter estimates and log-likelihood printed.
estimate if FALSE fits model but doesn't estimate predicted probabilities
refit if TRUE the algorithm will attempt multiple optimizations at different starting values if it doesn't converge
nrefits number of refitting attempts
initial a named list of starting values for the parameters (e.g. \$scale, \$shape, \$adjustment)
lowerbounds a vector of lowerbounds for the parameters
upperbounds a vector of upperbounds for the parameters
limit if TRUE restrict analysis to observations with detected=1
debug if TRUE, if fitting fails, return an object with fitting information
nofit if TRUE don't fit a model, but use the starting values and generate an object based on those values
optimx.method one (or a vector of) string(s) giving the optimisation method to use. If more than one is supplied, the results from one are used as the starting values for the next. See optimx
optimx.maxit maximum number of iterations to use in the optimisation.
mono. random.start By default when monotonicity constraints are enforced, a grid of starting values are tested. Instead random starting values can be used (uniformly distributed between the upper and lower bounds). Set TRUE for random start, FALSE (default) uses the grid method
silent silences warnings within ds fitting method (helpful for running many times without generating many warning/error messages).

Examples of distance sampling analyses are available at http://examples.distancesampling. org/.
Hints and tips on fitting (particularly optimisation issues) are on the mrds-opt manual page.

## Value

model object of class=(method, "ddf")

## Author(s)

Jeff Laake

## References

Laake, J.L. and D.L. Borchers. 2004. Methods for incomplete detection at distance zero. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.
Marques, F.F.C. and S.T. Buckland. 2004. Covariate models for the detection function. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.

## See Also

ddf.ds, ddf.io, ddf.io.fi, ddf.trial, ddf.trial.fi, ddf.rem, ddf.rem.fi, mrds-opt

## Examples

```
# load data
data(book.tee.data)
region <- book.tee.data$book.tee.region
egdata <- book.tee.data$book.tee.dataframe
samples <- book.tee.data$book.tee.samples
obs <- book.tee.data$book.tee.obs
# fit a half-normal detection function
result <- ddf(dsmodel=~mcds(key="hn", formula=~1), data=egdata, method="ds",
    meta.data=list(width=4))
# fit an independent observer model with full independence
result.io.fi <- ddf(mrmodel=~glm(~distance), data=egdata, method="io.fi",
    meta.data=list(width = 4))
# fit an independent observer model with point independence
result.io <- ddf(dsmodel=~cds(key = "hn"), mrmodel=~glm(~distance),
        data=egdata, method="io", meta.data=list(width=4))
## Not run:
# simulated single observer point count data (see ?ptdata.single)
data(ptdata.single)
ptdata.single$distbegin <- (as.numeric(cut(ptdata.single$distance,
                        10*(0:10)))-1)*10
ptdata.single$distend <- (as.numeric(cut(ptdata.single$distance,
    10*(0:10))))*10
model <- ddf(data=ptdata.single, dsmodel=~ cds(key="hn"),
        meta.data=list(point=TRUE,binned=TRUE,breaks=10*(0:10)))
summary(model)
plot(model,main="Single observer binned point data - half normal")
model <- ddf(data=ptdata.single, dsmodel=~cds(key="hr"),
    meta.data=list(point=TRUE, binned=TRUE, breaks=10*(0:10)))
summary(model)
plot(model, main="Single observer binned point data - hazard rate")
dev.new()
# simulated double observer point count data (see ?ptdata.dual)
# setup data
data(ptdata.dual)
ptdata.dual$distbegin <- (as.numeric(cut(ptdata.dual$distance,
    10*(0:10)))-1)*10
ptdata.dual$distend <- (as.numeric(cut(ptdata.dual$distance,
    10*(0:10))))*10
model <- ddf(method="io", data=ptdata.dual, dsmodel=~cds(key="hn"),
```

```
mrmodel=~glm(formula=~distance*observer),
meta.data=list(point=TRUE, binned=TRUE, breaks=10*(0:10)))
```

summary (model)
plot(model, main="Dual observer binned point data", new=FALSE, pages=1)
model <- ddf(method="io", data=ptdata.dual,
dsmodel=~cds(key="unif", adj.series="cos", adj.order=1),
mrmodel=~glm(formula=~distance*observer),
meta.data=list(point=TRUE, binned=TRUE, breaks=10*(0:10)))
summary (model)
$\operatorname{par}(m f r o w=c(2,3))$
plot(model,main="Dual observer binned point data", new=FALSE)
\#\# End(Not run)
ddf.ds CDS/MCDS Distance Detection Function Fitting

## Description

Fits a conventional distance sampling (CDS) (likelihood eq 6.6 in Laake and Borchers 2004) or multi-covariate distance sampling (MCDS)(likelihood eq 6.14 in Laake and Borchers 2004) model for the detection function of observed distance data. It only uses key functions and does not incorporate adjustment functions as in CDS/MCDS analysis engines in DISTANCE (Marques and Buckland 2004). Distance can be grouped (binned), ungrouped (unbinned) or mixture of the two. This function is not called directly by the user and is called from ddf,ddf.io, or ddf.trial.

## Usage

\#\# S3 method for class 'ds'
ddf(model, data, meta.data = list(), control = list(), call, method = "ds")

## Arguments

model model list with key function and scale formula if any
data data.frame; see ddf for details
meta.data list containing settings controlling data structure
control list containing settings controlling model fitting
call original function call if this function not called directly from ddf (e.g., called via ddf.io)
method analysis method; only needed if this function called from ddf.io or ddf.trial

## Details

For a complete description of each of the calling arguments, see ddf. The argument model in this function is the same as dsmodel in ddf. The argument dataname is the name of the dataframe specified by the argument data in ddf. The arguments control,meta. data, and method are defined the same as in ddf.

## Value

result: a ds model object

## Note

If mixture of binned and unbinned distance, width must be set to be $>=$ largest interval endpoint; this could be changed with a more complicated analysis; likewise, if all binned and bins overlap, the above must also hold; if bins don't overlap, width must be one of the interval endpoints; same holds for left truncation Although the mixture analysis works in principle it has not been tested via simulation.

## Author(s)

Jeff Laake

## References

Laake, J.L. and D.L. Borchers. 2004. Methods for incomplete detection at distance zero. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.
Marques, F.F.C. and S.T. Buckland. 2004. Covariate models for the detection function. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.

## See Also

flnl, summary.ds, coef.ds, plot.ds,gof.ds

## Examples

```
# ddf.ds is called when ddf is called with method="ds"
data(book.tee.data)
region <- book.tee.data$book.tee.region
egdata <- book.tee.data$book.tee.dataframe
samples <- book.tee.data$book.tee.samples
obs <- book.tee.data$book.tee.obs
result <- ddf(dsmodel = ~mcds(key = "hn", formula = ~1),
    data = egdata[egdata$observer==1, ], method = "ds",
    meta.data = list(width = 4))
summary(result,se=TRUE)
plot(result,main="cds - observer 1")
```

```
print(dht(result,region, samples,obs,options=list(varflag=0,group=TRUE),
            se=TRUE))
print(ddf.gof(result))
```

ddf.gof Goodness of fit tests for distance sampling models

## Description

Generic function that computes chi-square goodness of fit test for detection function models with binned data and Cramer-von Mises and Kolmogorov-Smirnov (if ks=TRUE)tests for exact distance data. By default a Q-Q plot is generated for exact data (and can be suppressed using the qq=FALSE argument).

## Usage

ddf.gof(
model,
breaks = NULL,
$\mathrm{nc}=\mathrm{NULL}$,
qq = TRUE,
nboot $=100$,
$\mathrm{ks}=\mathrm{FALSE}$,
)

## Arguments

| model | model object |
| :--- | :--- |
| breaks | Cutpoints to use for binning data |
| nc | Number of distance classes |
| qq | Flag to indicate whether quantile-quantile plot is desired <br> nboot |
| number of replicates to use to calculate p-values for the Kolmogorov-Smirnov |  |
| goodness of fit test statistics |  |$\quad$| perform the Kolmogorov-Smirnov test (this involves many bootstraps so can |
| :--- |
| take a while) |$\quad$| Graphics parameters to pass into qqplot function |
| :--- |

## Details

Formal goodness of fit testing for detection function models using Kolmogorov-Smirnov and Cramervon Mises tests. Both tests are based on looking at the quantile-quantile plot produced by qqplot.ddf and deviations from the line $x=y$.

The Kolmogorov-Smirnov test asks the question "what's the largest vertical distance between a point and the $y=x$ line?" It uses this distance as a statistic to test the null hypothesis that the samples
(EDF and CDF in our case) are from the same distribution (and hence our model fits well). If the deviation between the $y=x$ line and the points is too large we reject the null hypothesis and say the model doesn't have a good fit.
Rather than looking at the single biggest difference between the $\mathrm{y}=\mathrm{x}$ line and the points in the $\mathrm{Q}-\mathrm{Q}$ plot, we might prefer to think about all the differences between line and points, since there may be many smaller differences that we want to take into account rather than looking for one large deviation. Its null hypothesis is the same, but the statistic it uses is the sum of the deviations from each of the point to the line. Note that a bootstrap procedure is required for the KolmogorovSmirnov test to ensure that the p-values from the procedure are correct as the we are comparing the cumulative distribution function (CDF) and empirical distribution function (EDF) and we have estimated the parameters of the detection function. The nboot parameter controls the number of bootstraps to use. Set to 0 to avoid computing bootstraps (much faster but with no KolmogorovSmirnov results, of course).
One can change the precision of printed values by using the print.ddf.gof method's digits argument.

```
Value
List of class ddf.gof containing
chi-square Goodness of fit test statistic
df Degrees of freedom associated with test statistic
\(p\)-value \(\quad\) Significance level of test statistic
```


## Author(s)

Jeff Laake

## See Also

qqplot.ddf

```
ddf.io
```

Mark-Recapture Distance Sampling (MRDS) IO - PI

## Description

Mark-Recapture Distance Sampling (MRDS) Analysis of Independent Observer Configuration and Point Independence

## Usage

\#\# S3 method for class 'io'
ddf(dsmodel, mrmodel, data, meta.data = list(), control = list(), call = "")

## Arguments

| dsmodel | distance sampling model specification; model list with key function and scale <br> formula if any |
| :--- | :--- |
| mrmodel | mark-recapture model specification; model list with formula and link |
| data | analysis dataframe |
| meta.data | list containing settings controlling data structure |
| control | list containing settings controlling model fitting |
| call | original function call used to call ddf |

## Details

MRDS analysis based on point independence involves two separate and independent analyses of the mark-recapture data and the distance sampling data. For the independent observer configuration, the mark-recapture data are analysed with a call to ddf.io.fi (see likelihood eq 6.8 and 6.16 in Laake and Borchers 2004) to fit conditional distance sampling detection functions to estimate $p(0)$, detection probability at distance zero for the independent observer team based on independence at zero (eq 6.22 in Laake and Borchers 2004). Independently, the distance data, the union of the observations from the independent observers, are used to fit a conventional distance sampling (CDS) (likelihood eq 6.6) or multi-covariate distance sampling (MCDS) (likelihood eq 6.14) model for the detection function, $g(y)$, such that $g(0)=1$. The detection function for the observer team is then created as $p(y)=p(0) * g(y)$ (eq 6.28 of Laake and Borchers 2004) from which predictions are made. ddf. io is not called directly by the user and is called from ddf with method="io".
For a complete description of each of the calling arguments, see ddf. The argument dataname is the name of the dataframe specified by the argument data in ddf. The arguments dsmodel, mrmodel, control and meta.data are defined the same as in ddf.

## Value

result: an io model object which is composed of io.fi and ds model objects

## Author(s)

Jeff Laake

## References

Laake, J.L. and D.L. Borchers. 2004. Methods for incomplete detection at distance zero. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.

## See Also

ddf.io.fi, ddf.ds,summary.io,coef.io,plot.io, gof.io

```
ddf.io.fi
```


## Description

Mark-Recapture Analysis of Independent Observer Configuration with Full Independence

## Usage

\#\# S3 method for class 'io.fi'
ddf(model, data, meta.data = list(), control = list(), call = "", method)

## Arguments

| model | mark-recapture model specification |
| :--- | :--- |
| data | analysis dataframe |
| meta.data | list containing settings controlling data structure |
| control | list containing settings controlling model fitting |
| call | original function call used to call ddf |
| method | analysis method; only needed if this function called from ddf.io |

## Details

The mark-recapture data derived from an independent observer distance sampling survey can be used to derive conditional detection functions ( $p_{-} j(y)$ ) for both observers $(j=1,2)$. They are conditional detection functions because detection probability for observer $j$ is based on seeing or not seeing observations made by observer $3-j$. Thus, $p_{-} 1(y)$ is estimated by $p_{-} 1 / 2(y)$.
If detections by the observers are independent (full independence) then $p \_1(y)=p \_112(y), p \_2(y)=p \_2 \mid 1(y)$ and for the union, full independence means that $p(y)=p \_1(y)+p_{-} 2(y)-p \_1(y) * p \_2(y)$ for each distance $y$. In fitting the detection functions the likelihood given by eq 6.8 and 6.16 in Laake and Borchers (2004) is used. That analysis does not require the usual distance sampling assumption that perpendicular distances are uniformly distributed based on line placement that is random relative to animal distribution. However, that assumption is used in computing predicted detection probability which is averaged based on a uniform distribution (see eq 6.11 of Laake and Borchers 2004).
For a complete description of each of the calling arguments, see ddf. The argument model in this function is the same as mrmodel in ddf. The argument dataname is the name of the dataframe specified by the argument data in ddf. The arguments control,meta. data, and method are defined the same as in ddf.

## Value

result: an io.fi model object

## Author(s)

Jeff Laake

## References

Laake, J.L. and D.L. Borchers. 2004. Methods for incomplete detection at distance zero. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.

## See Also

ddf.io,summary.io.fi,coef.io.fi, plot.io.fi,gof.io.fi,io.glm

```
ddf.rem Mark-Recapture Distance Sampling (MRDS) Removal - PI
```


## Description

Mark-Recapture Distance Sampling (MRDS) Analysis of Removal Observer Configuration and Point Independence

## Usage

\#\# S3 method for class 'rem'
ddf(dsmodel, mrmodel, data, meta.data = list(), control = list(), call = "")

## Arguments

dsmodel distance sampling model specification; model list with key function and scale formula if any
mrmodel mark-recapture model specification; model list with formula and link
data analysis dataframe
meta.data list containing settings controlling data structure
control list containing settings controlling model fitting
call original function call used to call ddf

## Details

MRDS analysis based on point independence involves two separate and independent analyses of the mark-recapture data and the distance sampling data. For the removal observer configuration, the mark-recapture data are analysed with a call to ddf. rem.fi (see Laake and Borchers 2004) to fit conditional distance sampling detection functions to estimate $p(0)$, detection probability at distance zero for the primary observer based on independence at zero (eq 6.22 in Laake and Borchers 2004). Independently, the distance data, the observations from the primary observer, are used to fit a conventional distance sampling (CDS) (likelihood eq 6.6) or multi-covariate distance sampling (MCDS) (likelihood eq 6.14) model for the detection function, $g(y)$, such that $g(0)=1$. The detection function for the primary observer is then created as $p(y)=p(0) * g(y)$ (eq 6.28 of Laake and Borchers 2004) from which predictions are made. ddf. rem is not called directly by the user and is called from ddf with method="rem".

For a complete description of each of the calling arguments, see ddf. The argument data is the dataframe specified by the argument data in ddf. The arguments dsmodel, mrmodel, control and meta.data are defined the same as in ddf.

## Value

result: an rem model object which is composed of rem.fi and ds model objects

## Author(s)

Jeff Laake

## References

Laake, J.L. and D.L. Borchers. 2004. Methods for incomplete detection at distance zero. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.

## See Also

ddf.rem.fi, ddf.ds

```
ddf.rem.fi
Mark-Recapture Distance Sampling (MRDS) Removal - FI
```


## Description

Mark-Recapture Distance Sampling (MRDS) Analysis of Removal Observer Configuration with Full Independence

## Usage

\#\# S3 method for class 'rem.fi'
$\operatorname{ddf}($ model, data, meta.data $=$ list(), control $=$ list(), call = "", method)

## Arguments

model mark-recapture model specification
data analysis dataframe
meta.data list containing settings controlling data structure
control list containing settings controlling model fitting
call original function call used to call ddf
method analysis method; only needed if this function called from ddf.io

## Details

The mark-recapture data derived from an removal observer distance sampling survey can only derive conditional detection functions ( $p_{-} \mathrm{j}(\mathrm{y})$ ) for both observers $(\mathrm{j}=1)$ because technically it assumes that detection probability does not vary by occasion (observer in this case). It is a conditional detection function because detection probability for observer 1 is conditional on the observations seen by either of the observers. Thus, $p_{-} 1(y)$ is estimated by $p_{-} 1 / 2(y)$.
If detections by the observers are independent (full independence) then $p_{-} 1(y)=p \_112(y)$ and for the union, full independence means that $p(y)=p \_1(y)+p \_2(y)-p \_1(y) * p \_2(y)$ for each distance y. In fitting the detection functions the likelihood from Laake and Borchers (2004) are used. That analysis does not require the usual distance sampling assumption that perpendicular distances are uniformly distributed based on line placement that is random relative to animal distribution. However, that assumption is used in computing predicted detection probability which is averaged based on a uniform distribution (see eq 6.11 of Laake and Borchers 2004).
For a complete description of each of the calling arguments, see ddf. The argument model in this function is the same as mrmodel in ddf. The argument dataname is the name of the dataframe specified by the argument data in ddf. The arguments control,meta.data, and method are defined the same as in ddf.

## Value

result: an rem.fi model object

## Author(s)

Jeff Laake

## References

Laake, J.L. and D.L. Borchers. 2004. Methods for incomplete detection at distance zero. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.

## See Also

> ddf.io,rem.glm

> ddf.trial Mark-Recapture Distance Sampling (MRDS) Trial Configuration - PI

## Description

Mark-Recapture Distance Sampling (MRDS) Analysis of Trial Observer Configuration and Point Independence

## Usage

\#\# S3 method for class 'trial'
ddf(dsmodel, mrmodel, data, meta.data = list(), control = list(), call = "")

## Arguments

dsmodel distance sampling model specification; model list with key function and scale formula if any
mrmodel mark-recapture model specification; model list with formula and link
data analysis data.frame
meta.data list containing settings controlling data structure
control list containing settings controlling model fitting
call original function call used to call ddf

## Details

MRDS analysis based on point independence involves two separate and independent analyses of the mark-recapture data and the distance sampling data. For the trial configuration, the markrecapture data are analysed with a call to ddf.trial.fi (see likelihood eq 6.12 and 6.17 in Laake and Borchers 2004) to fit a conditional distance sampling detection function for observer 1 based on trials (observations) from observer 2 to estimate $p \_1(0)$, detection probability at distance zero for observer 1. Independently, the distance data from observer 1 are used to fit a conventional distance sampling (CDS) (likelihood eq 6.6) or multi-covariate distance sampling (MCDS) (likelihood eq 6.14) model for the detection function, $g(y)$, such that $g(0)=1$. The detection function for observer 1 is then created as $p_{-} 1(y)=p_{-} 1(0) * g(y)$ (eq 6.28 of Laake and Borchers 2004) from which predictions are made. ddf.trial is not called directly by the user and is called from ddf with method="trial".
For a complete description of each of the calling arguments, see ddf. The argument dataname is the name of the dataframe specified by the argument data in ddf. The arguments dsmodel, mrmodel, control and meta.data are defined the same as in ddf.

## Value

result: a trial model object which is composed of trial.fi and ds model objects

## Author(s)

Jeff Laake

## References

Laake, J.L. and D.L. Borchers. 2004. Methods for incomplete detection at distance zero. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.

## See Also

ddf.trial.fi, ddf.ds, summary.trial, coef.trial, plot.trial, gof.trial

```
ddf.trial.fi Mark-Recapture Analysis of Trial Configuration - FI
```


## Description

Mark-Recapture Analysis of Trial Observer Configuration with Full Independence

## Usage

```
## S3 method for class 'trial.fi'
ddf(model, data, meta.data = list(), control = list(), call = "", method)
```


## Arguments

| model | mark-recapture model specification |
| :--- | :--- |
| data | analysis dataframe |
| meta.data | list containing settings controlling data structure |
| control | list containing settings controlling model fitting |
| call | original function call used to call ddf |
| method | analysis method; only needed if this function called from ddf.trial |

## Details

The mark-recapture data derived from a trial observer distance sampling survey can be used to derive a conditional detection function ( $p_{-} 1(\mathrm{y})$ ) for observer 1 based on trials (observations) from observer 2. It is a conditional detection function because detection probability for observer 1 is based on seeing or not seeing observations made by observer 2. Thus, $p_{-} 1(y)$ is estimated by $p \_1 / 2(y)$. If detections by the observers are independent (full independence) then $p \_1(y)=p \_1 / 2(y)$ for each distance y. In fitting the detection functions the likelihood given by eq 6.12 or 6.17 in Laake and Borchers (2004) is used. That analysis does not require the usual distance sampling assumption that perpendicular distances are uniformly distributed based on line placement that is random relative to animal distribution. However, that assumption is used in computing predicted detection probability which is averaged based on a uniform distribution (see eq 6.13 of Laake and Borchers 2004).
For a complete description of each of the calling arguments, see ddf. The argument model in this function is the same as mrmodel in ddf. The argument dataname is the name of the dataframe specified by the argument data in ddf. The arguments control,meta. data, and method are defined the same as in ddf.

## Value

result: a trial.fi model object

## Author(s)

Jeff Laake

## References

Laake, J.L. and D.L. Borchers. 2004. Methods for incomplete detection at distance zero. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.

| See Also |
| :--- |
| ddf.trial, summary.trial.fi, coef.trial.fi, plot.trial.fi, gof.trial.fi |
| DeltaMethod |

## Description

Computes delta method variance-covariance matrix of results of any generic function fct that computes a vector of estimates as a function of a set of estimated parameters par.

## Usage

DeltaMethod(par, fct, vcov, delta, ...)

## Arguments

| par | vector of parameter values at which estimates should be constructed |
| :--- | :--- |
| fct | function that constructs estimates from parameters par |
| vcov | variance-covariance matrix of the parameters |
| delta | proportional change in parameters used to numerically estimate first derivative <br> with central-difference formula (ignored) |
| $\ldots$ | any additional arguments needed by fct |

## Details

The delta method (aka propagation of errors is based on Taylor series approximation - see Seber's book on Estimation of Animal Abundance). It uses the first derivative of fct with respect to par. It also uses the variance-covariance matrix of the estimated parameters which is derived in estimating the parameters and is an input argument.
The first argument of fct should be par which is a vector of parameter estimates. It should return a single value (or vector) of estimate(s). The remaining arguments of fct if any can be passed to fct by including them at the end of the call to DeltaMethod as name=value pairs.

## Value

a list with values
variance estimated variance-covariance matrix of estimates derived by fct
partial matrix (or vector) of partial derivatives of fct with respect to the parameters par

## Note

This is a generic function that can be used in any setting beyond the mrds package. However this is an internal function for mrds and the user does not need to call it explicitly.

## Author(s)

Jeff Laake and David L Miller

```
det.tables Observation detection tables
```


## Description

Creates a series of tables for dual observer data that shows the number missed and detected for each observer within defined distance classes.

## Usage

```
    det.tables(model, nc = NULL, breaks = NULL)
```


## Arguments

| model | fitted model from ddf |
| :--- | :--- |
| nc | number of equal-width bins for histogram |
| breaks | user define breakpoints |

## Value

list object of class "det.tables"
Observer1 table for observer 1
Observer2 table for observer 2
Duplicates histogram counts for duplicates
Pooled histogram counts for all observations by either observer
Obs1_2 table for observer 1 within subset seen by observer 2
Obs2_1 table for observer 2 within subset seen by observer 1

## Author(s)

Jeff Laake

## Examples

```
data(book.tee.data)
region <- book.tee.data$book.tee.region
egdata <- book.tee.data$book.tee.dataframe
samples <- book.tee.data$book.tee.samples
obs <- book.tee.data$book.tee.obs
xx <- ddf(mrmodel=~glm(formula=~distance*observer),
    dsmodel=~mcds(key="hn", formula=~sex),
    data=egdata, method="io", meta.data=list(width=4))
tabs <- det.tables(xx, breaks=c(0, 0.5, 1, 2, 3, 4))
par(mfrow=c(2, 2))
plot(tabs, new=FALSE, which=c(1, 2, 5, 6))
```

detfct.fit Fit detection function using key-adjustment functions

## Description

Fit detection function to observed distances using the key-adjustment function approach. If adjustment functions are included it will alternate between fitting parameters of key and adjustment functions and then all parameters much like the approach in the CDS and MCDS Distance FORTRAN code. To do so it calls detfct.fit.opt which uses the R optim function which does not allow non-linear constraints so inclusion of adjustments does allow the detection function to be non-monotone.

## Usage

detfct.fit(ddfobj, optim.options, bounds, misc.options)

## Arguments

| ddfobj | detection function object |
| :--- | :--- |
| optim.options | control options for optim |
| bounds | bounds for the parameters |
| misc.options | miscellaneous options |

## Value

fitted detection function model object with the following list structure
par final parameter vector
value final negative log likelihood value
counts number of function evaluations
convergence see codes in optim

```
detfct.fit.opt
message string about convergence
hessian hessian evaluated at final parameter values
aux
a list with 20 elements
- maxit: maximum number of iterations allowed for optimization
- lower: lower bound values for parameters
- upper: upper bound values for parameters
- setlower: TRUE if they are user set bounds
- setupper: TRUE if they are user set bounds
- point: TRUE if point counts and FALSE if line transect
- int.range: integration range values
- showit: integer value that determines information printed during iteration
- silent: option to silence errors from detfct.fit.opt
- integral.numeric if TRUE compute logistic integrals numerically
- breaks: breaks in distance for defined fixed bins for analysis
- maxiter: maximum iterations used
- refit: if TRUE, detection function will be fitted more than once if parameters are at a boundary or when convergence is not achieved
- nrefits: number of refittings
- mono: if TRUE monotonicity will be enforced
- mono.strict: if TRUE, then strict monotonicity is enforced; otherwise weak
- width: radius of point count or half-width of strip
- standardize: if TRUE, detection function is scaled so \(g(0)=1\)
- ddfobj: distance detection function object; see create.ddfobj
- bounded: TRUE if parameters ended up a boundary (I think)
- model: list of formulas for detection function model (probably can remove this)
```


## Author(s)

Dave Miller; Jeff Laake
detfct.fit.opt Fit detection function using key-adjustment functions

## Description

Fit detection function to observed distances using the key-adjustment function approach. If adjustment functions are included it will alternate between fitting parameters of key and adjustment functions and then all parameters much like the approach in the CDS and MCDS Distance FORTRAN code. This function is called by the driver function detfct. fit, then calls optimx function.

## Usage

detfct.fit.opt(ddfobj, optim.options, bounds, misc.options, fitting = "all")

## Arguments

| ddfobj | detection function object |
| :--- | :--- |
| optim.options | control options for optim |
| bounds | bounds for the parameters |
| misc.options | miscellaneous options <br> character string with values "all","key","adjust" to determine which parameters <br> fitting |
|  | are allowed to vary in the fitting |

## Value

fitted detection function model object with the following list structure

| par | final parameter vector |
| :--- | :--- |
| value | final negative log likelihood value |
| counts | number of function evaluations |
| convergence | see codes in optim |
| message | string about convergence |
| hessian | hessian evaluated at final parameter values |
| aux | a list with 20 elements |
|  | • maxit: maximum number of iterations allowed for optimization |
|  | - lower: lower bound values for parameters |

- upper: upper bound values for parameters
- setlower: TRUE if they are user set bounds
- setupper: TRUE if they are user set bounds
- point: TRUE if point counts and FALSE if line transect
- int.range: integration range values
- showit: integer value that determines information printed during iteration
- integral.numeric if TRUE compute logistic integrals numerically
- breaks: breaks in distance for defined fixed bins for analysis
- maxiter: maximum iterations used
- refit: if TRUE, detection function will be fitted more than once if parameters are at a boundary or when convergence is not achieved
- nrefits: number of refittings
- mono: if TRUE, monotonicity will be enforced
- mono.strict: if TRUE, then strict monotonicity is enforced; otherwise weak
- width: radius of point count or half-width of strip
- standardize: if TRUE, detection function is scaled so $g(0)=1$
- ddfobj: distance detection function object; see create.ddfobj
- bounded: TRUE if estimated parameters are at the bounds
- model: list of formulas for detection function model (probably can remove this)


## Author(s)

Dave Miller; Jeff Laake; Lorenzo Milazzo

## dht

Density and abundance estimates and variances

## Description

Compute density and abundance estimates and variances based on Horvitz-Thompson-like estimator.

```
Usage
    dht(
        model,
        region.table,
        sample.table,
        obs.table = NULL,
        subset = NULL,
        se = TRUE,
        options = list()
    )
```


## Arguments

| model | ddf model object |
| :--- | :--- |
| region.table | data.frame of region records. Two columns: Region.Label and Area. If only <br> density is required, one can set Area=0 for all regions. |
| sample.table | data.frame of sample records. Three columns: Region.Label, Sample.Label, <br> Effort. |
| obs.table | data.frame of observation records with fields: object, Region.Label, and <br> Sample.Label which give links to sample.table, region.table and the data <br> records used in model. Not necessary if the data.frame used to create the <br> model contains Region.Label, Sample.Label columns. |
| subset | subset statement to create obs.table <br> if TRUE computes standard errors, coefficient of variation and confidence inter- <br> se |
| options (based on log-normal approximation). See "Uncertainty" below. |  |

## Details

Density and abundance within the sampled region is computed based on a Horvitz-Thompsonlike estimator for groups and individuals (if a clustered population) and this is extrapolated to the entire survey region based on any defined regional stratification. The variance is based on replicate samples within any regional stratification. For clustered populations, $E(s)$ and its standard error are also output.
Abundance is estimated with a Horvitz-Thompson-like estimator (Huggins 1989, 1991; Borchers et al 1998; Borchers and Burnham 2004). The abundance in the sampled region is simply $1 / p_{1}+$
$1 / p_{2}+\ldots+1 / p_{n}$ where $p_{i}$ is the estimated detection probability for the $i$ th detection of $n$ total observations. It is not strictly a Horvitz-Thompson estimator because the $p_{i}$ are estimated and not known. For animals observed in tight clusters, that estimator gives the abundance of groups (group=TRUE in options) and the abundance of individuals is estimated as $s_{1} / p_{1}+s_{2} / p_{2}+\ldots+$ $s_{n} / p_{n}$, where $s_{i}$ is the size (e.g., number of animals in the group) of each observation (group=FALSE in options).
Extrapolation and estimation of abundance to the entire survey region is based on either a random sampling design or a stratified random sampling design. Replicate samples (lines) are specified within regional strata region.table, if any. If there is no stratification, region.table should contain only a single record with the Area for the entire survey region. The sample. table is linked to the region.table with the Region.Label. The obs.table is linked to the sample.table with the Sample.Label and Region.Label. Abundance can be restricted to a subset (e.g., for a particular species) of the population by limiting the list the observations in obs.table to those in the desired subset. Alternatively, if Sample.Label and Region.Label are in the data.frame used to fit the model, then a subset argument can be given in place of the obs. table. To use the subset argument but include all of the observations, use subset $=1==1$ to avoid creating an obs. table.

In extrapolating to the entire survey region it is important that the unit measurements be consistent or converted for consistency. A conversion factor can be specified with the convert.units variable in the options list. The values of Area in region. table, must be made consistent with the units for Effort in sample.table and the units of distance in the data.frame that was analyzed. It is easiest to do if the units of Area is the square of the units of Effort and then it is only necessary to convert the units of distance to the units of Effort. For example, if Effort was entered in kilometres and Area in square kilometres and distance in metres then using options=list (convert.units=0.001) would convert metres to kilometres, density would be expressed in square kilometres which would then be consistent with units for Area. However, they can all be in different units as long as the appropriate composite value for convert.units is chosen. Abundance for a survey region can be expressed as: $A * N / a$ where $A$ is Area for the survey region, $N$ is the abundance in the covered (sampled) region, and a is the area of the sampled region and is in units of Effort * distance. The sampled region a is multiplied by convert.units, so it should be chosen such that the result is in the same units of Area. For example, if Effort was entered in kilometres, Area in hectares $(100 \mathrm{~m} \times 100 \mathrm{~m})$ and distance in metres, then using options=list (convert.units=10) will convert a to units of hectares ( 100 to convert metres to 100 metres for distance and .1 to convert km to 100 m units).

The argument options is a list of variable=value pairs that set options for the analysis. All but two of these have been described above. pdelta should not need to be changed but was included for completeness. It controls the precision of the first derivative calculation for the delta method variance. If the option areas. supplied is TRUE then the covered area is assumed to be supplied in the CoveredArea column of the sample data.frame.

## Value

list object of class dht with elements:

| clusters | result list for object clusters |
| :--- | :--- |
| individuals | result list for individuals |
| Expected.S | data.frame of estimates of expected cluster size with fields Region, Expected.S <br> and se.Expected.S If each cluster size $=1$, then the result only includes indi- <br> viduals and not clusters and Expected.. |

The list structure of clusters and individuals are the same:
\(\left.$$
\begin{array}{ll}\text { bysample } & \begin{array}{l}\text { data.frame giving results for each sample; Nchat is the estimated abundance } \\
\text { within the sample and Nhat is scaled by surveyed area/covered area within that } \\
\text { region }\end{array} \\
\text { summary } & \begin{array}{l}\text { data.frame of summary statistics for each region and total }\end{array} \\
\text { N } & \begin{array}{l}\text { data.frame of estimates of abundance for each region and total } \\
\text { D }\end{array}
$$ <br>

data.frame of estimates of density for each region and total\end{array}\right]\)| average detection probability estimate |
| :--- |

Nhat. by . sample another summary of Nhat by sample used by dht. se

## Uncertainty

If the argument se=TRUE, standard errors for density and abundance is computed. Coefficient of variation and log-normal confidence intervals are constructed using a Satterthwaite approximation for degrees of freedom (Buckland et al. 2001 p. 90). The function dht. se computes the variance and interval estimates.

The variance has two components:

- variation due to uncertainty from estimation of the detection function parameters;
- variation in abundance due to random sample selection;

The first component (model parameter uncertainty) is computed using a delta method estimate of variance (Huggins 1989, 1991, Borchers et al. 1998) in which the first derivatives of the abundance estimator with respect to the parameters in the detection function are computed numerically (see DeltaMethod).

The second component (encounter rate variance) can be computed in one of several ways depending on the form taken for the encounter rate and the estimator used. To begin with there three possible values for varflag to calculate encounter rate:

- 0 uses a binomial variance for the number of observations (equation 13 of Borchers et al. 1998). This estimator is only useful if the sampled region is the survey region and the objects are not clustered; this situation will not occur very often;
- 1 uses the encounter rate $n / L$ (objects observed per unit transect) from Buckland et al. (2001) pg 78-79 (equation 3.78) for line transects (see also Fewster et al, 2009 estimator R2). This variance estimator is not appropriate if size or a derivative of size is used in the detection function;
- 2 is the default and uses the encounter rate estimator $\hat{N} / L$ (estimated abundance per unit transect) suggested by Innes et al (2002) and Marques \& Buckland (2004).

In general if any covariates are used in the models, the default varflag=2 is preferable as the estimated abundance will take into account variability due to covariate effects. If the population is clustered the mean group size and standard error is also reported.
For options 1 and 2, it is then possible to choose one of the estimator forms given in Fewster et al (2009) for line transects: "R2", "R3", "R4", "S1", "S2", "01", "02" or "03" by specifying the ervar= option (default "R2"). For points estimator "P3" is the only option. See varn and Fewster et al (2009) for further details on these estimators.

## dht options

Several options are available to control calculations and output:
ci.width Confidence interval width, expressed as a decimal between 0 and 1 (default 0.95 , giving a $95 \%$ CI)
pdelta delta value for computing numerical first derivatives (Default: 0.001)
varflag 0,1,2 (see "Uncertainty") (Default: 2)
convert. units multiplier for width to convert to units of length (Default: 1)
ervar encounter rate variance type (see "Uncertainty" and type argument of varn). (Default: "R2" for lines and "P3" for points)

## Author(s)

Jeff Laake, David L Miller

## References

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Innes, S., M.P. Heide-Jorgensen, J.L. Laake, K.L. Laidre, H.J. Cleator, P. Richard, and R.E.A. Stewart. 2002. Surveys of belugas and narwhals in the Canadian High Arctic in 1996. NAMMCO Scientific Publications 4: 169-190.

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## See Also

print.dht dht.se

$$
\begin{array}{ll}
\text { dht. deriv } & \text { Computes abundance estimates at specified parameter values using } \\
& \text { Horvitz-Thompson-like estimator }
\end{array}
$$

## Description

Computes abundance at specified values of parameters for numerical computation of first derivative with respect to parameters in detection function. An internal function called by DeltaMethod which is invoked by dht.se

## Usage

dht.deriv(par, model, obs, samples, options = list())

## Arguments

par detection function parameter values
model ddf model object
obs observations table
samples samples table
options list of options as specified in dht

## Value

vector of abundance estimates at values of parameters specified in par

## Note

Internal function; not intended to be called by user

## Author(s)

Jeff Laake

## See Also

dht, dht.se, DeltaMethod

## Description

Computes standard error, cv, and log-normal confidence intervals for abundance and density within each region (if any) and for the total of all the regions. It also produces the correlation matrix for regional and total estimates.

## Usage

dht.se(
model,
region.table, samples, obs, options, numRegions, estimate.table, Nhat.by.sample
)

## Arguments

model ddf model object
region.table table of region values
samples table of samples(replicates)
obs table of observations
options list of options that can be set (see dht)
numRegions number of regions
estimate.table table of estimate values
Nhat.by.sample estimated abundances by sample

## Details

The variance has two components:

- variation due to uncertainty from estimation of the detection function parameters;
- variation in abundance due to random sample selection;

The first component (model parameter uncertainty) is computed using a delta method estimate of variance (Huggins 1989, 1991, Borchers et al. 1998) in which the first derivatives of the abundance estimator with respect to the parameters in the detection function are computed numerically (see DeltaMethod).

The second component (encounter rate variance) can be computed in one of several ways depending on the form taken for the encounter rate and the estimator used. To begin with there three possible values for varflag to calculate encounter rate:

- 0 uses a binomial variance for the number of observations (equation 13 of Borchers et al. 1998). This estimator is only useful if the sampled region is the survey region and the objects are not clustered; this situation will not occur very often;
- 1 uses the encounter rate $n / L$ (objects observed per unit transect) from Buckland et al. (2001) pg 78-79 (equation 3.78) for line transects (see also Fewster et al, 2009 estimator R2). This variance estimator is not appropriate if size or a derivative of size is used in the detection function;
- 2 is the default and uses the encounter rate estimator $\hat{N} / L$ (estimated abundance per unit transect) suggested by Innes et al (2002) and Marques \& Buckland (2004).

In general if any covariates are used in the models, the default varflag=2 is preferable as the estimated abundance will take into account variability due to covariate effects. If the population is clustered the mean group size and standard error is also reported.

For options 1 and 2, it is then possible to choose one of the estimator forms given in Fewster et al (2009). For line transects: "R2", "R3", "R4", "S1", "S2", "01", "02" or "03" can be used by specifying the ervar= option (default "R2"). For point transects only the "P3" estimator may be used. See varn and Fewster et al (2009) for further details on these estimators.
Exceptions to the above occur if there is only one sample in a stratum. In that case it uses Poisson assumption $(\operatorname{Var}(x)=x)$ and it assumes a known variance so $z=1.96$ is used for critical value. In all other cases the degrees of freedom for the $t$-distribution assumed for the $\log$ (abundance) or $\log$ (density) is based on the Satterthwaite approximation (Buckland et al. 2001 pg 90 ) for the degrees of freedom (df). The df are weighted by the squared cv in combining the two sources of variation because of the assumed log-normal distribution because the components are multiplicative. For combining df for the sampling variance across regions they are weighted by the variance because it is a sum across regions.
A non-zero correlation between regional estimates can occur from using a common detection function across regions. This is reflected in the correlation matrix of the regional and total estimates which is given in the value list. It is only needed if subtotals of regional estimates are needed.

## Value

List with 2 elements:
estimate.table completed table with $\mathrm{se}, \mathrm{cv}$ and confidence limits
vc correlation matrix of estimates

## Note

This function is called by dht and it is not expected that the user will call this function directly but it is documented here for completeness and for anyone expanding the code or using this function in their own code.

## Author(s)

Jeff Laake

## References

see dht

## See Also

dht, print.dht

```
ds.function Distance Sampling Functions
```


## Description

Computes values of conditional and unconditional detection functions and probability density functions for for line/point data for single observer or dual observer in any of the 3 configurations (io,trial,rem).

## Usage

ds.function( model,
newdata $=$ NULL,
obs = "All",
conditional = FALSE,
pdf = TRUE,
finebr
)

## Arguments

| model | model object |
| :--- | :--- |
| newdata | dataframe at which to compute values; if NULL uses fitting data <br> obs |
| 1 or 2 for observer 1 or 2,3 for duplicates, "." for combined and "All" to return <br> all of the values |  |
| conditional | if FALSE, computes $\mathrm{p}(\mathrm{x})$ based on distance detection function and if TRUE <br> based on mr detection function <br> if FALSE, returns $\mathrm{p}(\mathrm{x})$ and if TRUE, returns $\mathrm{p}(\mathrm{x}) * \mathrm{pi}(\mathrm{x}) /$ integral $\mathrm{p}(\mathrm{x}) * \mathrm{pi}(\mathrm{x})$ |
| pdf | fine break values over which line is averaged |
| finebr |  |

## Details

Placeholder - Not functional -

## Value

List containing
xgrid grid of distance values
values average detection fct values at the xgrid values

## Author(s)

Jeff Laake

## flnl

Log-likelihood computation for distance sampling data

## Description

For a specific set of parameter values, it computes and returns the negative log-likelihood for the distance sampling likelihood for distances that are unbinned, binned and a mixture of both. The function flnl is the function minimized using optim from within ddf. ds.

## Usage

flnl(fpar, ddfobj, misc.options, fitting = "all")

## Arguments

fpar parameter values for detection function at which negative log-likelihood should be evaluated
ddfobj distance sampling object
misc.options a list with the following elements: width transect width; int.range the integration range for observations; showit 0 to 3 controls level debug output; integral. numeric if TRUE integral is computed numerically rather than analytically; point is this a point transect?
fitting character "key" if only fitting key function parameters, "adjust" if fitting adjustment parameters or "all" to fit both

## Details

Most of the computation is in flpt. lnl in which the negative log-likelihood is computed for each observation. flnl is a wrapper that optionally outputs intermediate results and sums the individual log-likelihood values.
flnl is the main routine that manipulates the parameters using getpar to handle fitting of key, adjustment or all of the parameters. It then calls flpt. lnl to do the actual computation of the likelihood. The probability density function for point counts is fr and for line transects is fx . $f x=g(x) / m u$ (where $g(x)$ is the detection function); whereas, $f(r)=r^{*} g(r) / m u$ where mu in both cases is the normalizing constant. Both functions are in source code file for link\{detfct\} and are called from distpdf and the integral calculations are made with integratepdf.

## Value

negative log-likelihood value at the parameter values specified in fpar

## Note

These are internal functions used by ddf. ds to fit distance sampling detection functions. It is not intended for the user to invoke these functions but they are documented here for completeness.

## Author(s)

Jeff Laake, David L Miller

## See Also

flt.var, detfct

flt.var | Hessian computation for fitted distance detection function model pa- |
| :--- |
| rameters |

## Description

Computes hessian to be used for variance-covariance matrix. The hessian is the outer product of the vector of first partials (see pg 62 of Buckland et al 2002).

## Usage

flt.var(ddfobj, misc.options)

## Arguments

ddfobj distance sampling object
misc.options width-transect width (W); int.range-integration range for observations; showit0 to 3 controls level of iteration printing; integral.numeric-if TRUE integral is computed numerically rather than analytically

## Value

variance-covariance matrix of parameters in the detection function

## Note

This is an internal function used by ddf. ds to fit distance sampling detection functions. It is not intended for the user to invoke this function but it is documented here for completeness.

## Author(s)

Jeff Laake and David L Miller

## References

Buckland et al. 2002

## See Also

flnl,flpt.lnl,ddf.ds
g0 Compute value of $p(0)$ using a logit formulation

## Description

Compute value of $\mathrm{p}(0)$ using a logit formulation

## Usage

g0(beta, z)

## Arguments

| beta | logistic parameters |
| :--- | :--- |
| $z$ | design matrix of covariate values |

## Value

vector of $p(0)$ values

## Author(s)

Jeff Laake
getpar Extraction and assignment of parameters to vector

## Description

Extracts parameters of a particular type (scale, shape, adjustments or $\mathrm{g} 0(\mathrm{p}(0))$ ) from the vector of parameters in ddfobj. All of the parameters are kept in a single vector for optimization even though they have very different uses. assign. par parses them from the vector based on a known structure and assigns them into ddfobj. getpar extracts the requested types to be extracted from ddfobj.

## Usage

```
getpar(ddfobj, fitting = "all", index = FALSE)
```


## Arguments

| ddfobj | distance sampling object (see create.ddfobj) |
| :--- | :--- |
| fitting | character string which is either "all","key","adjust" which determines which pa- <br> rameters are retrieved |
| index | logical that determines whether parameters are returned (FALSE) or starting <br> indices in parameter vector for scale, shape, adjustment parameters |

## Value

index $==$ FALSE, vector of parameters that were requested or index $==$ TRUE, vector of 3 indices for shape, scale, adjustment

## Note

Internal functions not intended to be called by user.

## Author(s)

Jeff Laake

See Also
assign.par

```
gof.ds
```


## Description

Compute chi-square goodness-of-fit test for ds models

## Usage

gof.ds(model, breaks = NULL, nc = NULL)

## Arguments

| model | ddf model object |
| :--- | :--- |
| breaks | distance cut points |
| nc | number of distance classes |

## Value

list with chi-square value, df and p-value

## Author(s)

Jeff Laake

## See Also

ddf.gof

```
gstdint Integral of pdf of distances
```


## Description

Computes the integral of distpdf with scale=1 (stdint=TRUE) or specified scale (stdint=FALSE).

## Usage

```
    gstdint(
        x,
        ddfobj,
        index = NULL,
        select = NULL,
        width,
        standardize = TRUE,
        point = FALSE,
        stdint = TRUE,
        doeachint = FALSE,
        left = left
    )
```


## Arguments

x
ddfobj distance detection function specification
index specific data row index
select logical vector for selection of data values
width
standardize if TRUE, divide through by the function evaluated at 0
point logical to determine if point (TRUE) or line transect(FALSE)
stdint if TRUE, scale=1 otherwise specified scale used
doeachint if TRUE perform integration using integrate
left left truncation width

## Value

vector of integral values of detection function

## Note

This is an internal function that is not intended to be invoked directly.

## Author(s)

Jeff Laake and David L Miller

```
histline Plot histogram line
```


## Description

Takes bar heights (height) and cutpoints (breaks), and constructs a line-only histogram from them using the function $\operatorname{plot}($ ) (if lineonly==FALSE) or lines() (if lineonly==TRUE).

## Usage

```
histline(
        height,
        breaks,
        lineonly = FALSE,
        outline = FALSE,
        ylim = range(height),
        xlab = "x",
        ylab = "y",
        det.plot = FALSE,
        add = FALSE,
    )
```


## Arguments

| height | heights of histogram bars |
| :--- | :--- |
| breaks | cutpoints for x |
| lineonly | if TRUE, drawn with plot; otherwise with lines to allow addition of current plot |
| outline | if TRUE, only outline of histogram is plotted |
| ylim | limits for $y$ axis |
| xlab | label for $x$ axis |
| ylab | label for y axis |
| det.plot | if TRUE, plot is of detection so yaxis limited to unit interval |
| add | should this plot add to a previous window |
| $\ldots$ | Additional unspecified arguments for plot |

## Value

None

## Author(s)

Jeff Laake and David L Miller

```
integratedetfct.logistic
```

Integrate a logistic detection function

## Description

Integrates a logistic detection function; a separate function is used because in certain cases the integral can be solved analytically and also because the scale trick used with the half-normal and hazard rate doesn't work with the logistic.

## Usage

integratedetfct.logistic(x, scalemodel, width, theta1, integral.numeric, w)

## Arguments

x
scalemodel
width
theta1 parameters for logistic
integral.numeric
if TRUE computes numerical integral value
w

## Value

vector of integral values

## Author(s)

Jeff Laake

## Description

Computes integral (analytically) over x from 0 to width of a logistic detection function; For reference see integral \#526 in CRC Std Math Table 24th ed

## Usage

integratelogistic.analytic(x, models, beta, width)

## Arguments

| $x$ | matrix of data |
| :--- | :--- |
| models | list of model formulae |
| beta | parameters of logistic detection function |
| width | transect half-width |

## Author(s)

Jeff Laake
integratepdf Numerically integrate pdf of observed distances over specified ranges

## Description

Computes integral of pdf of observed distances over $x$ for each observation. The method of computation depends on argument switches set and the type of detection function.

## Usage

```
integratepdf(
    ddfobj,
    select,
    width,
    int.range,
    standardize = TRUE,
    point = FALSE,
    left = 0,
    doeachint = FALSE
)
```


## Arguments

| ddfobj | distance detection function specification |
| :--- | :--- |
| select | logical vector for selection of data values |
| width | truncation width |
| int.range | integration range matrix; vector is converted to matrix |
| standardize | logical used to decide whether to divide through by the function evaluated at 0 |
| point | logical to determine if point count (TRUE) or line transect (FALSE) |
| left | left truncation width |
| doeachint | calculate each integral numerically |

## Value

vector of integral values - one for each observation

## Author(s)

Jeff Laake \& Dave Miller

## Description

Provides an iterative algorithm for finding the MLEs of detection (capture) probabilities for a twooccasion (double observer) mark-recapture experiment using standard algorithms GLM/GAM and an offset to compensate for conditioning on the set of observations. While the likelihood can be formulated and solved numerically, the use of GLM/GAM provides all of the available tools for fitting, predictions, plotting etc without any further development.

## Usage

```
io.glm(
    datavec,
    fitformula,
    eps = 1e-05,
    iterlimit = 500,
    GAM = FALSE,
    gamplot = TRUE
)
```


## Arguments

| datavec | dataframe |
| :--- | :--- |
| fitformula | logit link formula |
| eps | convergence criterion |
| iterlimit | maximum number of iterations allowed |
| GAM | uses GAM instead of GLM for fitting |
| gamplot | set to TRUE to get a gam plot object if GAM=TRUE |

## Details

Note that currently the code in this function for GAMs has been commented out until the remainder of the mrds package will work with GAMs. This is an internal function that is used as by ddf.io.fi to fit mark-recapture models with 2 occasions. The argument mrmodel is used for fitformula.

## Value

list of class("ioglm","glm","lm") or class("ioglm","gam")
glmobj GLM or GAM object
offsetvalue offsetvalues from iterative fit
plotobj gam plot object (if GAM \& gamplot==TRUE, else NULL)

## Author(s)

Jeff Laake, David Borchers, Charles Paxton

## References

Buckland, S.T., J.M. breiwick, K.L. Cattanach, and J.L. Laake. 1993. Estimated population size of the California gray whale. Marine Mammal Science, 9:235-249.
Burnham, K.P., S.T. Buckland, J.L. Laake, D.L. Borchers, T.A. Marques, J.R.B. Bishop, and L. Thomas. 2004. Further topics in distance sampling. pp: 360-363. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.

## is.linear.logistic Collection of functions for logistic detection functions

## Description

These functions are used to test whether a logistic detection function is a linear function of distance (is.linear.logistic) or is constant (varies by distance but no other covariates) is.logistic. constant). Based on these tests, the most appropriate manner for integrating the detection function with respect to distance is chosen. The integrals are needed to estimate the average detection probability for a given set of covariates.

## Usage

is.linear.logistic(xmat, g0model, zdim, width)

## Arguments

| xmat | data matrix |
| :--- | :--- |
| g0model | logit model |
| zdim | number of columns in design matrix |
| width | transect width |

## Details

If the logit is linear in distance then the integral can be computed analytically. If the logit is constant or only varies by distance then only one integral needs to be computed rather than an integral for each observation.

## Value

Logical TRUE if condition holds and FALSE otherwise

## Author(s)

Jeff Laake

```
is.logistic.constant Is a logit model constant for all observations?
```


## Description

Determines whether the specified logit model is constant for all observations. If it is constant then only one integral needs to be computed.

## Usage

```
is.logistic.constant(xmat, g0model, width)
```


## Arguments

| xmat | data |
| :--- | :--- |
| g0model | logit model |
| width | transect width |

## Value

logical value

Author(s)
Jeff Laake
keyfct.th1 Threshold key function

## Description

Threshold key function

## Usage

keyfct.th1 (distance, key.scale, key.shape)

## Arguments

distance perpendicular distance vector
key.scale vector of scale values
key.shape vector of shape values

## Value

vector of probabilities
keyfct.th2 Threshold key function

## Description

Threshold key function

## Usage

keyfct.th2(distance, key.scale, key.shape)

## Arguments

distance perpendicular distance vector
key.scale vector of scale values
key.shape vector of shape values

## Value

vector of probabilities

```
    keyfct.tpn Two-part normal key function
```


## Description

The two-part normal detection function of Becker and Christ (2015). Either side of an estimated apex in the distance histogram has a half-normal distribution, with differing scale parameters. Covariates may be included but affect both sides of the function.

## Usage <br> keyfct.tpn(distance, ddfobj)

## Arguments

distance perpendicular distance vector
ddfobj meta object containing parameters, design matrices etc

## Details

Two-part normal models have 2 important parameters:

- The apex, which estimates the peak in the detection function (where $g(x)=1$ ). The log apex is reported in summary results, so taking the exponential of this value should give the peak in the plotted function (see examples).
- The parameter that controls the difference between the sides .dummy_apex_side, which is automatically added to the formula for a two-part normal model. One can add interactions with this variable as normal, but don't need to add the main effect as it will be automatically added.


## Value

a vector of probabilities that the observation were detected given they were at the specified distance and assuming that $\mathrm{g}(\mathrm{mu})=1$

## Author(s)

Earl F Becker, David L Miller

## References

Becker, E. F., \& Christ, A. M. (2015). A Unimodal Model for Double Observer Distance Sampling Surveys. PLOS ONE, 10(8), e0136403. doi:10.1371/journal.pone. 0136403
lfbcvi
Black-capped vireo mark-recapture distance sampling analysis

## Description

These data represent avian point count surveys conducted at 453 point sample survey locations on the 24,000 (approx) live-fire region of Fort Hood in central Texas. Surveys were conducted by independent double observers ( 2 per survey occasion) and as such we had a maximum of 3 paired survey histories, giving a maximum of 6 sample occasions (see MacKenzie et al. 2006, MacKenzie and Royle 2005, and Laake et al. 2011 for various sample survey design details). At each point, we surveyed for 5 minutes (technically broken into 3 time intervals of 2 , 2 , and 1 minutes; not used here) and we noted detections by each observer and collected distance to each observation within a set of distance bins ( $0-25,25-50,50-75,75-100 \mathrm{~m}$ ) of the target species (Black-capped vireo's in this case) for each surveyor. Our primary focus was to use mark-recapture distance sampling methods to estimate density of Black-capped vireo's, and to estimate detection rates for the mark-recapture, distance, and composite model.

## Format

The format is a data frame with the following covariate metrics.
PointID Unique identifier for each sample location; locations are the same for both species
VisitNumber Visit number to the point
Species Species designation, either Golden-cheeked warbler (GW) or Black-capped Vireo (BV)
Distance Distance measure, which is either NA (representing no detection), or the median of the binned detection distances
PairNumber ID value indicating which observers were paired for that sampling occasion
Observer Observer ID, either primary (1), or secondary (2)
Detected Detection of a bird, either $1=$ detected, or $0=$ not detected
Date Date of survey since 15 march 2011
Pred Predicted occupancy value for that survey hexagon based on Farrell et al. (2013)
Category Region.Label categorization, see mrds help file for details on data structure
Effort Amount of survey effort at the point
Day Number of days since 15 March 2011
ObjectID Unique ID for each paired observations

## Details

In addition to detailing the analysis used by Collier et al. (2013, In Review), this example documents the use of mrds for avian point count surveys and shows how density models can be incorporated with occupancy models to develop spatially explicit density surface maps. For those that are interested, for the distance sampling portion of our analysis, we used both conventional distance sampling (cds) and multiple covariate distance sampling (mcds) with uniform and half-normal key
functions. For the mark-recapture portion of our analysis, we tended to use covariates for distance (median bin width), observer, and date of survey (days since 15 March 2011).
We combined our mrds density estimates via a Horvitz-Thompson styled estimator with the resource selection function gradient developed in Farrell et al. (2013) and estimated density on an $\sim 3.14 \mathrm{ha}$ hexagonal grid across our study area, which provided a density gradient for the Fort Hood military installation. Because there was considerable data manipulation needed for each analysis to structure the data appropriately for use in mrds, rather than wrap each analysis in a single function, we have provided both the Golden-cheeked warbler and Black-capped vireo analyses in their full detail. The primary differences you will see will be changes to model structures and model outputs between the two species.

## Author(s)

Bret Collier and Jeff Laake

## References

Farrell, S.F., B.A. Collier, K.L. Skow, A.M. Long, A.J. Campomizzi, M.L. Morrison, B. Hays, and R.N. Wilkins. 2013. Using LiDAR-derived structural vegetation characteristics to develop highresolution, small-scale, species distribution models for conservation planning. Ecosphere 43(3): 42. http://dx.doi.org/10.1890/ES12-000352.1

Laake, J.L., B.A. Collier, M.L. Morrison, and R.N. Wilkins. 2011. Point-based mark recapture distance sampling. Journal of Agricultural, Biological and Environmental Statistics 16: 389-408.
Collier, B.A., S.L. Farrell, K.L. Skow, A. M. Long, A.J. Campomizzi, K.B. Hays, J.L. Laake, M.L. Morrison, and R.N. Wilkins. 2013. Spatially explicit density of endangered avian species in a disturbed landscape. Auk, In Review.

## Examples

```
## Not run:
data(lfbcvi)
xy=cut(lfbcvi$Pred, c(-0.0001, .1, .2, .3, .4, . 5, . 6, .7, .8, .9, 1),
    labels=c("1", "2", "3", "4", "5", "6", "7", "8", "9", "10"))
x=data.frame(lfbcvi, New=xy)
# Note that I scaled the individual covariate of day-helps with
# convergence issues
bird.data <- data.frame(object=x$ObjectID, observer=x$Observer,
                    detected=x$Detected, distance=x$Distance,
                Region.Label=x$New, Sample.Label=x$PointID,
                Day=(x$Day/max(x$Day)))
# make observer a factor variable
bird.data$observer=factor(bird.data$observer)
# Jeff Laake suggested this snippet to quickly create distance medians
# which adds bin information to the bird.data dataframe
bird.data$distbegin=0
```

```
bird.data$distend=100
bird.data$distend[bird.data$distance==12.5]=25
bird.data$distbegin[bird.data$distance==37.5]=25
bird.data$distend[bird.data$distance==37.5]=50
bird.data$distbegin[bird.data$distance==62.5]=50
bird.data$distend[bird.data$distance==62.5]=75
bird.data$distbegin[bird.data$distance==87.5]=75
bird.data$distend[bird.data$distance==87.5]=100
# Removed all survey points with distance=NA for a survey event;
# hence no observations for use in ddf() but needed later
bird.data=bird.data[complete.cases(bird.data),]
# Manipulations on full dataset for various data.frame creation for
# use in density estimation using dht()
#Samples dataframe
xx=x
x=data.frame(PointID=x$PointID, Species=x$Species,
    Category=x$New, Effort=x$Effort)
x=x[!duplicated(x$PointID),]
point.num=table(x$Category)
samples=data.frame(PointID=x$PointID, Region.Label=x$Category,
                    Effort=x$Effort)
final.samples=data.frame(Sample.Label=samples$PointID,
                        Region.Label=samples$Region.Label,
                        Effort=samples$Effort)
#obs dataframe
obs=data.frame(ObjectID=xx$ObjectID, PointID=xx$PointID)
#used to get Region and Sample assigned to ObjectID
obs=merge(obs, samples, by=c("PointID", "PointID"))
obs=obs[!duplicated(obs$ObjectID),]
obs=data.frame(object=obs$ObjectID, Region.Label=obs$Region.Label,
            Sample.Label=obs$PointID)
region.data=data.frame(Region.Label=c(1, 2, 3,4,5,6,7,8,9, 10),
Area=c(point.num[1]*3.14, point.num[2]*3.14,
    point.num[3]*3.14, point.num[4]*3.14,
    point.num[5]*3.14, point.num[6]*3.14,
    point.num[7]*3.14, point.num[8]*3.14,
    point.num[9]*3.14, point.num[10]*3.14))
# Candidate Models
BV1=ddf(
    dsmodel=~mcds(key="hn",formula=~1),
    mrmodel=~glm(~distance),
    data=bird.data,
    method="io",
    meta.data=list(binned=TRUE,point=TRUE,width=100,breaks=c(0,50,100)))
BV1FI=ddf(
    dsmodel=~mcds(key="hn",formula=~1),
```

```
    mrmodel=~glm(~distance),
    data=bird.data,
    method="io.fi",
    meta.data=list(binned=TRUE, point=TRUE,width=100,breaks=c(0,50,100)))
BV2=ddf(
    dsmodel=~mcds(key="hr",formula=~1),
    mrmodel=~glm(~distance),
    data=bird.data,
    method="io",
    meta.data=list(binned=TRUE, point=TRUE,width=100,breaks=c(0,50,100)))
BV3=ddf(
    dsmodel=~mcds(key="hn",formula=~1),
    mrmodel=~glm(~distance+observer),
    data=bird.data,
    method="io",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
BV3FI=ddf(
    dsmodel=~mcds(key="hn",formula=~1),
    mrmodel=~glm(~distance+observer),
    data=bird.data,
    method="io.fi",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
BV4=ddf(
    dsmodel=~mcds(key="hr",formula=~1),
    mrmodel=~glm(~distance+observer),
    data=bird.data,
    method="io",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
BV5=ddf(
    dsmodel=~mcds(key="hn",formula=~1),
    mrmodel=~glm(~distance*observer),
    data=bird.data,
    method="io",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
BV5FI=ddf(
    dsmodel=~mcds(key="hn",formula=~1),
    mrmodel=~glm(~distance*observer),
    data=bird.data,
    method="io.fi",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
BV6=ddf(
    dsmodel=~mcds(key="hr",formula=~1),
    mrmodel=~glm(~distance*observer),
    data=bird.data,
    method="io",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
BV7=ddf(
    dsmodel=~cds(key="hn",formula=~1),
    mrmodel=~glm(~distance*Day),
    data=bird.data,
    method="io",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
BV7FI=ddf(
```

```
    dsmodel=~cds(key="hn",formula=~1),
    mrmodel=~glm(~distance*Day),
    data=bird.data,
    method="io.fi",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
BV8=ddf(
    dsmodel=~cds(key="hr",formula=~1),
    mrmodel=~glm(~distance*Day),
    data=bird.data,
    method="io",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
BV9=ddf(
    dsmodel=~mcds(key="hn",formula=~1),
    mrmodel=~glm(~distance*observer*Day),
    data=bird.data,
    method="io",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
BV9FI=ddf(
    dsmodel=~mcds(key="hn",formula=~1),
    mrmodel=~glm(~distance*observer*Day),
    data=bird.data,
    method="io.fi",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
BV10=ddf(
    dsmodel=~mcds(key="hr",formula=~1),
    mrmodel=~glm(~distance*observer*Day),
    data=bird.data,
    method="io",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
#BV.DS=ddf(
# dsmodel=~mcds(key="hn",formula=~1),
# data=bird.data,
# method="ds",
# meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
#AIC table building code.
AIC = c(BV1$criterion, BV1FI$criterion, BV2$criterion, BV3$criterion,
    BV3FI$criterion, BV4$criterion, BV5$criterion, BV5FI$criterion,
    BV6$criterion, BV7$criterion, BV7FI$criterion, BV8$criterion,
    BV9$criterion, BV9FI$criterion, BV10$criterion)
#creates a set of row names for me to check my grep() call below
rn = c("BV1", "BV1FI", "BV2", "BV3", "BV3FI", "BV4", "BV5", "BV5FI",
    "BV6", "BV7", "BV7FI", "BV8", "BV9", "BV9FI", "BV10")
#Number parameters
k = c(length(BV1$par), length(BV1FI$par), length(BV2$par),
    length(BV3$par), length(BV3FI$par), length(BV4$par),
    length(BV5$par),length(BV5FI$par), length(BV6$par),
    length(BV7$par), length(BV7FI$par), length(BV8$par),
\#build AIC table
AIC.table=data.frame \((A I C=A I C, r n=r n, k=k, d A I C=a b s(m i n(A I C)-A I C)\),
```

$$
\operatorname{likg}=\exp (-.5 *(\operatorname{abs}(\min (A I C)-A I C))))
$$

\#row.names(AIC.table)=grep("BV", ls(), value=TRUE)
AIC.table=AIC.table[with(AIC.table, order(-likg, -dAIC, AIC, k)),]
AIC.table=data.frame(AIC.table, wi=AIC.table\$likg/sum(AIC.table\$likg))
AIC.table
\# Model average N_hat_covered estimates
\# not very clean, but I wanted to show full process, need to use
\# collect.models and model.table here later on
estimate <- c(BV1\$Nhat, BV1FI\$Nhat, BV2\$Nhat, BV3\$Nhat, BV3FI\$Nhat, BV4\$Nhat, BV5\$Nhat, BV5FI\$Nhat, BV6\$Nhat, BV7\$Nhat, BV7FI\$Nhat, BV8\$Nhat, BV9\$Nhat, BV9FI\$Nhat, BV10\$Nhat)

AIC.values=AIC
\# had to use str() to extract here as Nhat.se is calculated in \# mrds:::summary.io, not in ddf(), so it takes a bit std.err <- c(summary (BV1)\$Nhat.se, summary (BV1FI) \$Nhat.se, summary (BV2)\$Nhat. se, summary (BV3)\$Nhat.se, summary (BV3FI) \$Nhat.se, summary (BV4)\$Nhat.se, summary (BV5) \$Nhat.se, summary (BV5FI)\$Nhat.se, summary (BV6)\$Nhat.se, summary (BV7)\$Nhat.se, summary (BV7FI)\$Nhat. se, summary (BV8)\$Nhat. se, summary (BV9) \$Nhat. se, summary (BV9FI) \$Nhat.se, summary (BV10)\$Nhat.se)
\#\# End(Not run)
\#\# Not run:
\#Not Run
\#requires RMark
library (RMark)
\#uses model.average structure to model average real abundance estimates for
\#covered area of the surveys
mmi.list=list(estimate=estimate, AIC=AIC.values, se=std.err)
model. average(mmi.list, revised=TRUE)
\#Not Run
\#Summary for the top 2 models
\#summary (BV5, se=TRUE)
\#summary (BV5FI, se=TRUE)
\#Not Run
\#Best Model
\#best.model=AIC.table[1,]
\#Not Run
\#GOF for models
\#ddf.gof(BV5, breaks=c(0, 25, 50, 75, 100))
\#Not Run
\#Density estimation across occupancy categories
\#out.BV=dht(BV5, region.data, final.samples, obs, se=TRUE,
options=list(convert.units=.01))

```
#Plot--Not Run
```

\#Composite Detection Function
\#plot(BV5, which=3, showpoints=FALSE, angle=0, density=0, col="black", lwd=3,
\# main="Black-capped Vireo", xlab="Distance (m)", las=1, cex.axis=1.25,
\# cex.lab=1.25)
\#\# End(Not run)
lfgcwa Golden-cheeked warbler mark-recapture distance sampling analysis

## Description

These data represent avian point count surveys conducted at 453 point sample survey locations on the 24,000 (approx) live-fire region of Fort Hood in central Texas. Surveys were conducted by independent double observers ( 2 per survey occasion) and as such we had a maximum of 3 paired survey histories, giving a maximum of 6 sample occasions (see MacKenzie et al. 2006, MacKenzie and Royle 2005, and Laake et al. 2011 for various sample survey design details). At each point, we surveyed for 5 minutes (technically broken into 3 time intervals of 2,2 , and 1 minutes; not used here) and we noted detections by each observer and collected distance to each observation within a set of distance bins ( $0-50,50-100 \mathrm{~m}$; Laake et al. 2011) of the target species (Golden-cheeked warblers in this case) for each surveyor. Our primary focus was to use mark-recapture distance sampling methods to estimate density of Golden-cheeked warblers, and to estimate detection rates for the mark-recapture, distance, and composite model.

## Format

The format is a data frame with the following covariate metrics.
PointID Unique identifier for each sample location; locations are the same for both species
VisitNumber Visit number to the point
Species Species designation, either Golden-cheeked warbler (GW) or Black-capped Vireo (BV)
Distance Distance measure, which is either NA (representing no detection), or the median of the binned detection distances
PairNumber ID value indicating which observers were paired for that sampling occasion
Observer Observer ID, either primary (1), or secondary (2)
Detected Detection of a bird, either $1=$ detected, or $0=$ not detected
Date Date of survey since 15 March 2011, numeric value
Pred Predicted occupancy value for that survey hexagon based on Farrell et al. (2013)
Category Region.Label categorization, see R package mrds help file for details on data structure
Effort Amount of survey effort at the point
Day Number of days since 15 March 2011, numeric value
ObjectID Unique ID for each paired observations

## Details

In addition to detailing the analysis used by Collier et al. (2013, In Review), this example documents the use of mrds for avian point count surveys and shows how density models can be incorporated with occupancy models to develop spatially explicit density surface maps. For those that are interested, for the distance sampling portion of our analysis, we used both conventional distance sampling (cds) and multiple covariate distance sampling (mcds) with uniform and half-normal key functions. For the mark-recapture portion of our analysis, we tended to use covariates for distance (median bin width), observer, and date of survey (days since 15 March 2011).
We combined our mrds density estimates via a Horvitz-Thompson styled estimator with the resource selection function gradient developed in Farrell et al. (2013) and estimated density on an $\sim 3.14$ ha hexagonal grid across our study area, which provided a density gradient for Fort Hood. Because there was considerable data manipulation needed for each analysis to structure the data appropriately for use in mrds, rather than wrap each analysis in a single function, we have provided both the Golden-cheeked warbler and Black-capped vireo analyses in their full detail. The primary differences you will see will be changes to model structures and model outputs between the two species.

## Author(s)

Bret Collier and Jeff Laake

## References

Farrell, S.F., B.A. Collier, K.L. Skow, A.M. Long, A.J. Campomizzi, M.L. Morrison, B. Hays, and R.N. Wilkins. 2013. Using LiDAR-derived structural vegetation characteristics to develop highresolution, small-scale, species distribution models for conservation planning. Ecosphere 43(3): 42. http://dx.doi.org/10.1890/ES12-000352.1

Laake, J.L., B.A. Collier, M.L. Morrison, and R.N. Wilkins. 2011. Point-based mark recapture distance sampling. Journal of Agricultural, Biological and Environmental Statistics 16: 389-408.
Collier, B.A., S.L. Farrell, K.L. Skow, A.M. Long, A.J. Campomizzi, K.B. Hays, J.L. Laake, M.L. Morrison, and R.N. Wilkins. 2013. Spatially explicit density of endangered avian species in a disturbed landscape. Auk, In Review.

## Examples

```
## Not run:
data(lfgcwa)
xy <- cut(lfgcwa$Pred, c(-0.0001, .1, . 2, . 3, .4, . 5, . 6, . 7, . 8, . 9, 1),
    labels=c("1", "2", "3", "4", "5", "6", "7", "8", "9", "10"))
x <- data.frame(lfgcwa, New=xy)
# Note that I scaled the individual covariate of day-helps with
# convergence issues
bird.data <- data.frame(object=x$ObjectID, observer=x$Observer,
                        detected=x$Detected, distance=x$Distance,
                Region.Label=x$New, Sample.Label=x$PointID,
                Day=(x$Day/max(x$Day)))
```

```
# make observer a factor variable
bird.data$observer=factor(bird.data$observer)
# Jeff Laake suggested this snippet to quickly create distance medians
# which adds bin information to the \code{bird.data} dataframe
bird.data$distbegin=0
bird.data$distend=100
bird.data$distend[bird.data$distance==12.5]=50
bird.data$distbegin[bird.data$distance==37.5]=0
bird.data$distend[bird.data$distance==37.5]=50
bird.data$distbegin[bird.data$distance==62.5]=50
bird.data$distend[bird.data$distance==62.5]=100
bird.data$distbegin[bird.data$distance==87.5]=50
bird.data$distend[bird.data$distance==87.5]=100
# Removed all survey points with distance=NA for a survey event;
# hence no observations for use in \code{ddf()} but needed later
bird.data=bird.data[complete.cases(bird.data),]
# Manipulations on full dataset for various data.frame creation
# for use in density estimation using \code{dht()}
# Samples dataframe
xx <- x
x <- data.frame(PointID=x$PointID, Species=x$Species,
    Category=x$New, Effort=x$Effort)
x <- x[!duplicated(x$PointID),]
point.num <- table(x$Category)
samples <- data.frame(PointID=x$PointID, Region.Label=x$Category,
        Effort=x$Effort)
final.samples=data.frame(Sample.Label=samples$PointID,
                                    Region.Label=samples$Region.Label,
                                    Effort=samples$Effort)
# obs dataframe
obs <- data.frame(ObjectID=xx$ObjectID, PointID=xx$PointID)
# used to get Region and Sample assigned to ObjectID
obs <- merge(obs, samples, by=c("PointID", "PointID"))
obs <- obs[!duplicated(obs$ObjectID),]
obs <- data.frame(object=obs$ObjectID, Region.Label=obs$Region.Label,
    Sample.Label=obs$PointID)
#Region.Label dataframe
region.data <- data.frame(Region.Label=c(1, 2, 3, 4, 5, 6,7,8,9),
Area=c(point.num[1]*3.14,
point.num[2]*3.14,
    point.num[3]*3.14,
    point.num[4]*3.14,
    point.num[5]*3.14,
    point.num[6]*3.14,
    point.num[7]*3.14,
    point.num[8]*3.14,
```

```
point.num[9]*3.14))
```

\# Candidate Models
GW1 $=\operatorname{ddf}($
dsmodel=~cds(key="unif", adj.series="cos", adj.order=1,adj.scale="width"),
mrmodel=~glm(~distance),
data=bird.data,
method="io",
meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))

## GW2=ddf(

dsmodel=~cds(key="unif", adj.series="cos", adj.order=1,adj.scale="width"),
mrmodel=~glm(~distance+observer),
data=bird.data,
method="io",
meta.data=list(binned=TRUE, point=TRUE, width=100, breaks=c(0,50,100)))
GW3=ddf(
dsmodel=~cds(key="unif", adj.series="cos", adj.order=1,adj.scale="width"), mrmodel=~glm(~distance*observer),
data=bird.data,
method="io",
meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
GW4=ddf(
dsmodel=~mcds(key="hn", formula=~1), mrmodel=~glm(~distance),
data=bird.data,
method="io", meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
GW4FI=ddf(
dsmodel=~mcds(key="hn",formula=~1),
mrmodel=~glm(~distance),
data=bird.data,
method="io.fi",
meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
GW5=ddf(
dsmodel=~mcds(key="hn", formula=~1),
mrmodel=~glm(~distance+observer),
data=bird.data,
method="io", meta.data=list(binned=TRUE, point=TRUE, width=100, breaks=c (0,50, 100)))
GW5FI=ddf(
dsmodel=~mcds(key="hn", formula=~1),
mrmodel=~glm(~distance+observer),
data=bird.data,
method="io.fi",
meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
GW6=ddf(
dsmodel=~mcds(key="hn", formula=~1),
mrmodel=~glm(~distance*observer),
data=bird.data,
method="io",
meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c $(0,50,100))$ )
GW6FI=ddf(

```
    dsmodel=~mcds(key="hn",formula=~1),
    mrmodel=~glm(~distance*observer),
    data=bird.data,
    method="io.fi",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
GW7=ddf(
    dsmodel=~cds(key="hn",formula=~1),
    mrmodel=~glm(~distance*Day),
    data=bird.data,
    method="io",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
GW7FI=ddf(
    dsmodel=~cds(key="hn",formula=~1),
    mrmodel=~glm(~distance*Day),
    data=bird.data,
    method="io.fi",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
GW8=ddf(
    dsmodel=~mcds(key="hn",formula=~1),
    mrmodel=~glm(~distance*observer*Day),
    data=bird.data,
    method="io",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
GW8FI=ddf(
    dsmodel=~mcds(key="hn",formula=~1),
    mrmodel=~glm(~distance*observer*Day),
    data=bird.data,
    method="io.fi",
    meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
#GWDS=ddf(
# dsmodel=~mcds(key="hn",formula=~1),
# data=bird.data,
# method="ds",
# meta.data=list(binned=TRUE, point=TRUE, width=100,breaks=c(0,50,100)))
```


## \#\#\#\# GCWA Summary Metrics

```
\#AIC table building code, not exactly elegant, but I did not
want to add more package dependencies
AIC = c(GW1$criterion, GW2$criterion, GW3$criterion, GW4$criterion,
    GW4FI$criterion, GW5$criterion, GW5FI$criterion,
    GW6$criterion, GW6FI$criterion, GW7$criterion, GW7FI$criterion,
    GW8$criterion, GW8FI$criterion)
#creates a set of row names for me to check my grep() call below
rn <- c("GW1", "GW2", "GW3", "GW4", "GW4FI", "GW5", "GW5FI", "GW6",
    "GW6FI", "GW7","GW7FI", "GW8", "GW8FI")
# number of parameters for each model
k <- c(length(GW1$par), length(GW2$par), length(GW3$par), length(GW4$par),
    length(GW4FI$par), length(GW5$par), length(GW5FI$par),
```

```
    length(GW6$par), length(GW6FI$par), length(GW7$par),
    length(GW7FI$par), length(GW8$par), length(GW8FI$par))
# build AIC table and
AIC.table <- data.frame(AIC = AIC, rn=rn, k=k, dAIC = abs(min(AIC)-AIC),
            likg = exp(-.5*(abs(min(AIC)-AIC))))
# row.names(AIC.table)=grep("GW", ls(), value=TRUE)
AIC.table <- AIC.table[with(AIC.table, order(-likg, -dAIC, AIC, k)),]
AIC.table <- data.frame(AIC.table, wi=AIC.table$likg/sum(AIC.table$likg))
AIC.table
# Model average N_hat_covered estimates
# not very clean, but I wanted to show full process, need to use
# collect.models and model.table here
estimate <- c(GW1$Nhat, GW2$Nhat, GW3$Nhat, GW4$Nhat, GW4FI$Nhat,
                GW5$Nhat, GW5FI$Nhat, GW6$Nhat, GW6FI$Nhat, GW7$Nhat,
    GW7FI$Nhat, GW8$Nhat, GW8FI$Nhat)
AIC.values <- AIC
# Nhat.se is calculated in mrds:::summary.io, not in ddf(), so
# it takes a bit to pull out
std.err <- c(summary(GW1)$Nhat.se, summary(GW2)$Nhat.se,
    summary(GW3)$Nhat. se, summary(GW4)$Nhat.se,
    summary(GW4FI)$Nhat.se, summary(GW5)$Nhat.se,
    summary(GW5FI)$Nhat.se, summary(GW6)$Nhat.se,
    summary(GW6FI)$Nhat.se, summary(GW7)$Nhat.se,
    summary(GW7FI)$Nhat.se, summary(GW8)$Nhat.se,
    summary(GW8FI)$Nhat.se)
## End(Not run)
## Not run:
#Not Run
#requires RMark
library(RMark)
#uses model.average structure to model average real abundance estimates for
#covered area of the surveys
mmi.list=list(estimate=estimate, AIC=AIC.values, se=std.err)
model.average(mmi.list, revised=TRUE)
#Not Run
#Best Model FI
#best.modelFI=AIC.table[1,]
#best.model
#Best Model PI
#best.modelPI=AIC.table[2,]
#best.modelPI
#Not Run
#summary(GW7FI, se=TRUE)
#summary(GW7, se=TRUE)
#Not Run
```

```
#GOF for models
#ddf.gof(GW7, breaks=c(0,50,100))
#Not Run
#Density estimation across occupancy categories
#out.GW=dht(GW7, region.data, final.samples, obs, se=TRUE,
    options=list(convert.units=.01))
#Plots--Not Run
#Composite Detection Function examples
#plot(GW7, which=3, showpoints=FALSE, angle=0, density=0,
# col="black", lwd=3, main="Golden-cheeked Warbler",
# xlab="Distance (m)", las=1, cex.axis=1.25, cex.lab=1.25)
#Conditional Detection Function
#dd=expand.grid(distance=0:100,Day=(4:82)/82)
#dmat=model.matrix(~distance*Day,dd)
#dd$p=plogis(model.matrix(~distance*Day,dd)%*%coef(GW7$mr)$estimate)
#dd$Day=dd$Day*82
#with(dd[dd$Day==12,],plot(distance,p,ylim=c(0,1), las=1,
# ylab="Detection probability", xlab="Distance (m)",
# type="l",lty=1, lwd=3, bty="l", cex.axis=1.5, cex.lab=1.5))
#with(dd[dd$Day==65,],lines(distance,p,lty=2, lwd=3))
#ch=paste(bird.data$detected[bird.data$observer==1],
# bird.data$detected[bird.data$observer==2],
# sep="")
#tab=table(ch,cut(82*bird.data$Day[bird.data$observer==1],c(0,45,83)),
# cut(bird.data$distance[bird.data$observer==1],c(0,50,100)))
#tabmat=cbind(colMeans(rbind(tab[3, ,1]/colSums(tab[2:3, ,1],
#
# colMeans(rbind(tab[3,,2]/colSums(tab[2:3, ,2],
# tab[3,,2]/colSums(tab[c(1,3),,2])))))
#colnames(tabmat)=c("0-50", "51-100")
#points(c(25,75),tabmat[1,],pch=1, cex=1.5)
#points(c(25,75),tabmat[2,],pch=2, cex=1.5)
# Another alternative plot using barplot instead of points
# (this is one in paper)
#ch=paste(bird.data$detected[bird.data$observer==1],
# bird.data$detected[bird.data$observer==2],
#sep="")
#tab=table(ch,cut(82*bird.data$Day[bird.data$observer==1],c(0,45,83)),
# cut(bird.data$distance[bird.data$observer==1],c(0,50,100)))
#tabmat=cbind(colMeans(rbind(tab[3, ,1]/colSums(tab[2:3, ,1],
# tab[3,,1]/colSums(tab[c(1,3),,1])))),
#colMeans(rbind(tab[3,,2]/colSums(tab[2:3, 2],
# tab[3,,2]/colSums(tab[c(1,3), ,2])))))
#colnames(tabmat)=c("0-50", "51-100")
#par(mfrow=c(2, 1), mai=c(1,1,1,1))
#with(dd[dd$Day==12,],
# plot(distance,p,ylim=c(0,1), las=1,
# ylab="Detection probability", xlab="",
```

```
# type="l",lty=1, lwd=4, bty="l", cex.axis=1.5, cex.lab=1.5))
#segments(0, 0, .0, tabmat[1,1], lwd=3)
#segments(0, tabmat[1,1], 50, tabmat[1,1], lwd=4)
#segments(50, tabmat[1,1], 50, 0, lwd=4)
#segments(50, tabmat[1,2], 100, tabmat[1,2], lwd=4)
#segments(0, tabmat[1,1], 50, tabmat[1,1], lwd=4)
#segments(100, tabmat[1,2], 100, 0, lwd=4)
#mtext("a",line=-1, at=90)
#with(dd[dd$Day==65,],
# plot(distance,p,ylim=c(0,1), las=1, ylab="Detection probability",
# xlab="Distance", type="l",lty=1,
# lwd=4, bty="l", cex.axis=1.5, cex.lab=1.5))
#segments(0, 0, .0, tabmat[2,1], lwd=4)
#segments(0, tabmat[2,1], 50, tabmat[2,1], lwd=4)
#segments(50, tabmat[2,1], 50, 0, lwd=4)
#segments(50, tabmat[2,2], 50, tabmat[2,1], lwd=4)
#segments(50, tabmat[2,2], 100, tabmat[2,2], lwd=4)
#segments(100, tabmat[2,2], 100, 0, lwd=4)
#mtext("b",line=-1, at=90)
## End(Not run)
```

logisticbyx Logistic as a function of covariates

## Description

treats logistic as a function of covariates; for a given covariate combination it computes function at with those covariate values at a range of distances

## Usage

logisticbyx(distance, x, models, beta, point)

## Arguments

| distance | vector of distance values |
| :--- | :--- |
| x | covariate data |
| models | model list |
| beta | logistic parameters |
| point | TRUE if a point transect model |

## Value

vector of probabilities

## Author(s)

Jeff Laake
logisticbyz Logistic as a function of distance

## Description

Treats logistic as a function of distance; for a given distance it computes function at all covariate values in data.

## Usage

logisticbyz(x, distance, models, beta)

## Arguments

| x | covariate data |
| :--- | :--- |
| distance | single distance value |
| models | model list |
| beta | logistic parameters |

## Value

vector of probabilities

## Author(s)

Jeff Laake
logisticdetfct Logistic detection function

## Description

Logistic detection function

## Usage

logisticdetfct(distance, theta, w, std = FALSE)

## Arguments

| distance | perpendicular distance vector |
| :--- | :--- |
| theta | scale parameters |
| w | scale covariate matrix |
| std | if TRUE uses scale=1 |

The routine returns a vector of probabilities that the observation were detected given they were at the specified distance and assuming that $\mathrm{g}(0)=1$ (ie a standard line transect detection function).
logisticdupbyx Logistic for duplicates as a function of covariates

## Description

Treats logistic for duplicates as a function of covariate z ; for a given z it computes the function at with those covariate values at a range of distances.

## Usage

logisticdupbyx(distance, x1, x2, models, beta, point)

## Arguments

| distance | vector of distance values |
| :--- | :--- |
| x1 | covariate data for fct 1 |
| x2 | covariate data for fct 2 |
| models | model list |
| beta | logistic parameters |
| point | TRUE for point transect data |

## Value

vector of probabilities

## Author(s)

Jeff Laake
logisticdupbyx_fast Logistic for duplicates as a function of covariates (fast)

## Description

As logisticdupbyx, but faster when distance is a covariate (but no interactions with distance occur.

## Usage

logisticdupbyx_fast(distance, $x 1, \mathrm{x} 2, \mathrm{models}$, beta, point, beta_distance)

## Arguments

| distance | vector of distance values |
| :--- | :--- |
| x1 | linear predictor for 1, without distance |
| x2 | linear predictor for 2, without distance |
| models | model list |
| beta | logistic parameters |
| point | TRUE for point transect data |
| beta_distance | parameter for distance |

## Author(s)

David L Miller

| logit $\quad$ Logit function |
| :--- | :--- |

## Description

Computes logit transformation.

## Usage

logit(p)

## Arguments

$\mathrm{p} \quad$ probability

## Value

$\operatorname{logit}(p)$ returns $[\log (p /(1-p)]$

## Author(s)

Jeff Laake
logLik.ddf log-likelihood value for a fitted detection function

## Description

Extract the log-likelihood from a fitted detection function.

## Usage

\#\# S3 method for class 'ddf'
logLik(object, ...)

## Arguments

$$
\begin{array}{ll}
\text { object } & \text { a fitted detection function model object } \\
\ldots & \text { included for } \mathrm{S} 3 \text { completeness, but ignored }
\end{array}
$$

## Value

a numeric value giving the log-likelihood with two attributes: "df" the "degrees of freedom" for the model (number of parameters) and "nobs" the number of observations used to fit the model

## Author(s)

David L Miller

```
    mcds MCDS function definition
```


## Description

Creates model formula list for multiple covariate distance sampling using values supplied in call to ddf

```
Usage
    mcds(
        formula = NULL,
        key = NULL,
        adj.series = NULL,
        adj.order = c(NULL),
        adj.scale = "width",
        adj.exp = FALSE,
        shape.formula = ~1
    )
```


## Arguments

| formula | formula for scale function |
| :--- | :--- |
| key | string identifying key function (currently either "hn" (half-normal), "hr" (hazard- <br> rate), "unif" (uniform) or "gamma" (gamma distribution) |
| adj.series | string identifying adjustment functions cos (Cosine), herm (Hermite polynomi- <br> als), poly (simple polynomials) or NULL |
| adj.order | vector of order of adjustment terms to include |
| adj.scale | whether to scale the adjustment terms by "width" or "scale" |
| adj.exp | if TRUE uses exp(adj) for adjustment to keep $f(x)>0$ <br> shape.formula |
|  | formula for shape function |

Value
A formula list used to define the detection function model
fct string "mcds"
key key function string
adj.series adjustment function string
adj.order adjustment function orders
adj.scale adjustment function scale type
formula formula for scale function
shape.formula formula for shape function

## Author(s)

Jeff Laake; Dave Miller
mrds-opt Tips on optimisation issues in mrds models

## Description

Occasionally when fitting an 'mrds' model one can run into optimisation issues. In general such problems can be quite complex so these "quick fixes" may not work. If you come up against problems that are not fixed by these tips, or you feel the results are dubious please go ahead and contact the package authors.

## Debug mode

One can obtain debug output at each stage of the optimisation using the showit option. This is set via control, so adding control=list(showit=3) gives the highest level of debug output (setting showit to 1 or 2 gives less output).

## Re-scaling covariates

Sometimes convergence issues in covariate (MCDS) models are caused by values of the covariate being very large, so a rescaling of that covariate is then necessary. Simply scaling by the standard deviation of the covariate can help (e.g. dat\$size.scaled<- dat\$scale/sd(dat\$scale) for a covariate size, then including size. scaled in the model instead of size).
It is important to note that one needs to use the original covariate (size) when computing HorvitzThompson estimates of population size if the group size is used in that estimate. i.e. use the unscaled size in the numerator of the $\mathrm{H}-\mathrm{T}$ estimator.

## Factor levels

By default R will set the base factor level to be the label which comes first alphabetically. Sometimes this can be an issue when that factor level corresponds to a subset of the data with very few observations. This can lead to very large uncertainty estimates (CVs) for model parameters. One way around this is to use relevel to set the base level to a level with more observations.

## Initial values

Initial (or starting) values can be set via the initial element of the control list. initial is a list itself with elements scale, shape and adjustment, corresponding to the associated parameters. If a model has covariates then the scale or shape elements will be vectors with parameter initial values in the same order as they are specific in the model formula (using showit is a good check they are in the correct order). Adjustment starting values are in order of the order of that term (cosine order 2 is before cosine order 3 terms).
One way of obtaining starting values is to fit a simpler model first (say with fewer covariates or adjustments) and then use the starting values from this simpler model for the corresponding parameters.

Another alternative to obtain starting values is to fit the model (or some submodel) using Distance for Windows. Note that Distance reports the scale parameter (or intercept in a covariate model) on the exponential scale, so one must log this before supplying it to ddf.

## Bounds

One can change the upper and lower bounds for the parameters. These specify the largest and smallest values individual parameters can be. By placing these constraints on the parameters, it is possible to "temper" the optimisation problem, making fitting possible.
Again, one uses the control list, the elements upperbounds and lowerbounds. In this case, each of upperbounds and lowerbounds are vectors, which one can think of as each of the vectors scale, shape and adjustment from the "Initial values" section above, concatenated in that order. If one does not occur (e.g. no shape parameter) then it is simple omitted from the vector.

## Author(s)

David L. Miller [dave@ninepointeightone.net](mailto:dave@ninepointeightone.net)

NCovered Compute estimated abundance in covered (sampled) region

## Description

Generic function that computes abundance within the covered region. It calls method (class) specific functions for the computation.

## Usage

NCovered(par, model = NULL, group = TRUE)

## Arguments

| par | parameter values (used when computing derivatives wrt parameter uncertainty) if NULL parameter values in model are used |
| :---: | :---: |
| model | ddf model object |
| group | if TRUE computes group abundance and if FALSE individual abundance |

## Value

abundance estimate

## Author(s)

Jeff Laake

```
nlminb_wrapper Wrapper around nlminb
```


## Description

This is a wrapper around nlminb to use scaling, as this is not available in optimx.

## Usage

```
nlminb_wrapper(
    par,
    11,
    ugr = NULL,
    lower = NULL,
    upper = NULL,
    mcontrol,
    hess = NULL,
    ddfobj,
    data,
)
```


## Arguments

| par | starting parameters |
| :--- | :--- |
| ll | log likelihood function |
| ugr | gradient function |
| lower | lower bounds on parameters |
| upper | upper bounds on parameters |
| mcontrol | control options |
| hess | hessian function |
| ddfobj | detection function specification object |
| data | the data |
| $\ldots$ | anything else to pass to ll |

## Value

optimx object

## Author(s)

David L Miller, modified from optimx. run by JC Nash, R Varadhan, G Grothendieck.

```
p.det Double-platform detection probability
```


## Description

Computes detection probability for detection function computed from mark-recapture data with possibly different link functions.

## Usage

p.det(dpformula, dplink, dppars, dpdata)

## Arguments

dpformula formula for detection function
dplink link function ("logit","loglog","cloglog")
dppars parameter vector
dpdata double platform data

## Value

vector of predicted detection probabilities

## Author(s)

## Description

Generate a table of frequencies of probability of detection from a detection function model. This is particularly useful when employing covariates, as it can indicate if there are detections with very small detection probabilities that can be unduly influential when calculating abundance estimates.

## Usage

p.dist.table(object, bins $=\operatorname{seq}(0,1$, by $=0.1)$, proportion $=$ FALSE)
p_dist_table(object, bins $=\operatorname{seq}(0,1$, by $=0.1)$, proportion = FALSE)

## Arguments

object fitted detection function
bins how the results should be binned
proportion should proportions be returned as well as counts?

## Details

Because dht uses a Horvitz-Thompson-like estimator, abundance estimates can be sensitive to errors in the estimated probabilities. The estimator is based on $\sum 1 / \hat{P}_{a}\left(z_{i}\right)$, which means that the sensitivity is greater for smaller detection probabilities. As a rough guide, we recommend that the method be not used if more than say $5 \%$ of the $\hat{P}_{a}\left(z_{i}\right)$ are less than 0.2 , or if any are less than 0.1 . If these conditions are violated, the truncation distance w can be reduced. This causes some loss of precision relative to standard distance sampling without covariates.

## Value

a data. frame with probability bins, counts and (optionally) proportions. The object has an attribute p_range which contains the range of estimated detection probabilities

## Author(s)

David L Miller

## References

Marques, F.F.C. and S.T. Buckland. 2004. Covariate models for the detection function. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R. Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.

## Examples

\#\# Not run:
\# try out the tee data
data(book.tee.data)
egdata <- book.tee.data\$book.tee.dataframe
\# fit model with covariates
result <- ddf(dsmodel = ~mcds(key = "hn", formula = ~sex+size), data = egdata[egdata\$observer==1, ], method = "ds", meta.data $=$ list(width $=4$ ))
\# print table
p.dist.table(result)
\# with proportions
p.dist.table(result, proportion=TRUE)
\#\# End(Not run)

```
parse.optimx Parse optimx results and present a nice object
```


## Description

Take the resulting object from a call to optimx and make it into an object that mrds wants to talk to.

## Usage

parse.optimx(lt, lnl.last, par.last)

## Arguments

| lt | an optimx object |
| :--- | :--- |
| lnl.last | last value of the log likelihood |
| par.last | last value of the parameters |

## Value

lt object that can be used later on

```
pdot.dsr.integrate.logistic
```

Compute probability that a object was detected by at least one observer

## Description

Computes probability that a object was detected by at least one observer (pdot or $p_{-}$.) for a logistic detection function that contains distance.

## Usage

pdot.dsr.integrate.logistic(
right,
width,
beta,
x ,
integral.numeric,
BT,
models,
GAM = FALSE,
rem $=$ FALSE,
point $=$ FALSE
)

## Arguments

right either an integration range for binned data (vector of 2) or the rightmost value for integration (from 0 to right)
width transect width
beta parameters of logistic detection function
$x \quad$ data matrix
integral.numeric
set to TRUE unless data are binned (done in this fct) or the model is such that distance is not linear (eg distance^2), If integral.numeric is FALSE it will compute the integral analytically. It should only be FALSE if is.linear.logistic function is TRUE.
BT FALSE except for the trial configuration; BT stands for Buckland-Turnock who initially proposed a trial configuration for dual observers
models list of models including g0model
GAM Not used at present. The idea was to be able to use a GAM for $g(0)$ portion of detection function; should always be F
rem only TRUE for the removal configuration but not used and could be removed if pulled from the function calls. Originally thought the pdot integral would differ but it is the same as the io formula. The only thing that differs with removal is that $\mathrm{p}(2 \mid 1)=1$. Observer 2 sees everything seen by observer 1 ,
point TRUE for point transects

## Author(s)

Jeff Laake
plot.det.tables Observation detection tables

## Description

Plot the tables created by det.tables. Produces a series of tables for dual observer data that shows the number missed and detected for each observer within defined distance classes.

## Usage

```
## S3 method for class 'det.tables'
    plot(
        x,
        which = 1:6,
        angle = NULL,
        density = NULL,
        col1 = "white",
        col2 = "lightgrey",
        new = TRUE,
    ...
    )
```


## Arguments

x which items in $x$ to plot (vector with values in 1:6)
angle shading angle for hatching
density shading density for hatching
col1 plotting colour for total histogram bars.
col2 plotting colour for subset histogram bars.
new if TRUE new plotting window for each plot
... other graphical parameters, passed to plotting functions

## Details

Plots that are produced are as follows (controlled by the which argument):
1 Detected by either observer/Detected by observer 1
2 Detected by either observer/Detected by observer 2

3 Seen by both observers
4 Seen by either observer
5 Detected by observer 2/Detected by observer 1। 2
6 Detected by observer 1/Detected by observer 2। 1

## Value

Just plots.

## Author(s)

Jeff Laake, David L Miller

## Examples

```
data(book.tee.data)
region <- book.tee.data$book.tee.region
egdata <- book.tee.data$book.tee.dataframe
samples <- book.tee.data$book.tee.samples
obs <- book.tee.data$book.tee.obs
xx <- ddf(mrmodel=~glm(formula=~distance*observer),
    dsmodel = ~mcds(key = "hn", formula = ~sex),
    data = egdata, method = "io", meta.data = list(width = 4))
tabs <- det.tables(xx,breaks=c(0,.5,1,2,3,4))
par(mfrow=c(2,3))
plot(tabs,which=1:6,new=FALSE)
```

plot.ds

Plot fit of detection functions and histograms of data from distance sampling model

## Description

Plots the fitted detection function(s) with a histogram of the observed distances to compare visually the fitted model and data.

## Usage

```
## S3 method for class 'ds'
plot(
    x,
    which = 2,
    breaks = NULL,
    nc = NULL,
    jitter.v = rep(0, 3),
    showpoints = TRUE,
```

```
    subset = NULL,
    pl.col = "lightgrey",
    pl.den = NULL,
    pl.ang = NULL,
    main = NULL,
    pages = 0,
    pdf = FALSE,
    ylim = NULL,
    xlab = "Distance",
    ylab = NULL,
)
```


## Arguments

x
which index to specify which plots should be produced:

1 histogram of observed distances
2 histogram of observed distances with fitted line and points (default)
breaks user defined breakpoints
nc number of equal width bins for histogram
jitter.v apply jitter to points by multiplying the fitted value by a random draw from a normal distribution with mean 1 and sd jitter.v.
showpoints logical variable; if TRUE plots predicted value for each observation (conditional on its observed distance).
subset subset of data to plot.
pl.col colour for histogram bars.
pl.den shading density for histogram bars.
pl.ang shading angle for histogram bars.
main plot title.
pages the number of pages over which to spread the plots. For example, if pages=1 then all plots will be displayed on one page. Default is 0 , which prompts the user for the next plot to be displayed.
pdf plot the histogram of distances with the PDF of the probability of detection overlaid. Ignored (with warning) for line transect models.
ylim vertical axis limits.
$x l a b \quad$ horizontal axis label (defaults to "Distance").
ylab vertical axis label (default automatically set depending on plot type).
... other graphical parameters, passed to the plotting functions (plot, hist, lines, points, etc).

## Details

The structure of the histogram can be controlled by the user-defined arguments nc or breaks. The observation specific detection probabilities along with the line representing the fitted average detection probability.

It is not intended for the user to call plot.ds but its arguments are documented here. Instead the generic plot command should be used and it will call the appropriate function based on the class of the ddf object.

## Value

Just plots.

## Author(s)

Jeff Laake, Jon Bishop, David Borchers, David L Miller

## See Also

add_df_covar_line

## Examples

```
# fit a model to the tee data
data(book.tee.data)
egdata <- book.tee.data$book.tee.dataframe
xx <- ddf(dsmodel=~mcds(key="hn", formula=~sex),
    data=egdata[egdata$observer==1, ],
    method="ds", meta.data=list(width=4))
# not showing predicted probabilities
plot(xx, breaks=c(0, 0.5, 1, 2, 3, 4), showpoints=FALSE)
# two subsets
plot(xx, breaks=c(0, 0.5, 1, 2, 3, 4), subset=sex==0)
plot(xx, breaks=c(0, 0.5, 1, 2, 3, 4), subset=sex==1)
# put both plots on one page
plot(xx, breaks=c(0, 0.5, 1, 2, 3, 4), pages=1, which=1:2)
```

plot.io
Plot fit of detection functions and histograms of data from distance sampling independent observer (io) model

## Description

Plots the fitted detection functions for a distance sampling model and histograms of the distances (for unconditional detection functions) or proportion of observations detected within distance intervals (for conditional detection functions) to compare visually the fitted model and data.

```
Usage
    ## S3 method for class 'io'
    plot(
        x,
        which = 1:6,
        breaks = NULL,
        nc = NULL,
        maintitle = "",
        showlines = TRUE,
        showpoints = TRUE,
        ylim = c(0, 1),
        angle = NULL,
        density = NULL,
        col = "lightgrey",
        jitter = NULL,
        divisions = 25,
        pages = 0,
        xlab = "Distance",
        ylab = "Detection probability",
        subtitle = TRUE,
    )
```


## Arguments

x
which
fitted model from ddf
index to specify which plots should be produced.
1 Plot primary unconditional detection function
2 Plot secondary unconditional detection function
3 Plot pooled unconditional detection function
4 Plot duplicate unconditional detection function
5 Plot primary conditional detection function
6 Plot secondary conditional detection function

Note that the order of which is ignored and plots are produced in the above order.
breaks user define breakpoints
nc number of equal-width bins for histogram
maintitle main title line for each plot

| showlines | logical variable; if TRUE a line representing the average detection probability is plotted |
| :---: | :---: |
| showpoints | logical variable; if TRUE plots predicted value for each observation |
| ylim | range of vertical axis; defaults to ( 0,1 ) |
| angle | shading angle for histogram bars. |
| density | shading density for histogram bars. |
| col | colour for histogram bars. |
| jitter | scaling option for plotting points. Jitter is applied to points by multiplying the fitted value by a random draw from a normal distribution with mean 1 and sd jitter. |
| divisions | number of divisions for averaging line values; default $=25$ |
| pages | the number of pages over which to spread the plots. For example, if pages=1 then all plots will be displayed on one page. Default is 0 , which prompts the user for the next plot to be displayed. |
| xlab | label for x -axis |
| ylab | label for y -axis |
| subtitle | if TRUE, shows plot type as sub-title |
|  | other graphical parameters, passed to the plotting functions (plot, hist, lines, points, etc) |

## Details

The structure of the histogram can be controlled by the user-defined arguments nc or breaks. The observation specific detection probabilities along with the line representing the fitted average detection probability.

It is not intended for the user to call plot.io.fi but its arguments are documented here. Instead the generic plot command should be used and it will call the appropriate function based on the class of the ddf object.

## Value

Just plots

## Author(s)

Jeff Laake, Jon Bishop, David Borchers, David L Miller

## Examples

```
library(mrds)
data(book.tee.data)
egdata <- book.tee.data$book.tee.dataframe
result.io <- ddf(dsmodel=~cds(key = "hn"), mrmodel=~glm(~distance),
    data=egdata, method="io", meta.data=list(width=4))
```

```
# just plot everything
plot(result.io)
# Plot primary and secondary unconditional detection functions on one page
# and primary and secondary conditional detection functions on another
plot(result.io,which=c(1,2,5,6),pages=2)
```

plot.io.fi

Plot fit of detection functions and histograms of data from distance sampling independent observer model with full independence (io.fi)

## Description

Plots the fitted detection functions for a distance sampling model and histograms of the distances (for unconditional detection functions) or proportion of observations detected within distance intervals (for conditional detection functions) to compare visually the fitted model and data.

## Usage

```
## S3 method for class 'io.fi'
plot(
        x,
        which = 1:6,
        breaks = NULL,
        nc = NULL,
        maintitle = "",
        showlines = TRUE,
        showpoints = TRUE,
        ylim = c(0, 1),
        angle = NULL,
        density = NULL,
        col = "lightgrey",
        jitter = NULL,
        divisions = 25,
        pages = 0,
        xlab = "Distance",
        ylab = "Detection probability",
        subtitle = TRUE,
    )
```


## Arguments

which index to specify which plots should be produced.

\[\)| 1 |  Plot primary unconditional detection function  |
| :--- | :--- |
| 2 |  Plot secondary unconditional detection function  |
| 3 |  Plot pooled unconditional detection function  |
| 4 |  Plot duplicate unconditional detection function  |
| 5 |  Plot primary conditional detection function  |
| 6 |  Plot secondary conditional detection function  |

\]

| breaks | Note that the order of which is ignored and plots are produced in the above order. user define breakpoints |
| :---: | :---: |
| nc | number of equal-width bins for histogram |
| maintitle | main title line for each plot |
| showlines | logical variable; if TRUE a line representing the average detection probability is plotted |
| showpoints | logical variable; if TRUE plots predicted value for each observation |
| ylim | range of vertical axis; defaults to ( 0,1 ) |
| angle | shading angle for histogram bars. |
| density | shading density for histogram bars. |
| col | colour for histogram bars. |
| jitter | scaling option for plotting points. Jitter is applied to points by multiplying the fitted value by a random draw from a normal distribution with mean 1 and sd jitter. |
| divisions | number of divisions for averaging line values; default $=25$ |
| pages | the number of pages over which to spread the plots. For example, if pages=1 then all plots will be displayed on one page. Default is 0 , which prompts the user for the next plot to be displayed. |
| xlab | label for x -axis |
| ylab | label for y -axis |
| subtitle | if TRUE, shows plot type as sub-title |
|  | other graphical parameters, passed to the plotting functions (plot, hist, lines, points, etc) |

## Details

The structure of the histogram can be controlled by the user-defined arguments nc or breaks. The observation specific detection probabilities along with the line representing the fitted average detection probability.

It is not intended for the user to call plot.io.fi but its arguments are documented here. Instead the generic plot command should be used and it will call the appropriate function based on the class of the ddf object.

## Value

Just plots.

## Author(s)

Jeff Laake, Jon Bishop, David Borchers, David L Miller

## Examples

```
library(mrds)
data(book.tee.data)
egdata <- book.tee.data$book.tee.dataframe
result.io.fi <- ddf(mrmodel=~glm(~distance), data = egdata, method = "io.fi",
    meta.data = list(width = 4))
# just plot everything
plot(result.io.fi)
# Plot primary and secondary unconditional detection functions on one page
# and primary and secondary conditional detection functions on another
plot(result.io.fi,which=c(1, 2,5,6),pages=2)
```

plot.layout Layout for plot methods in mrds

## Description

This function does the paging, using devAskNewPage(). This means we can just call plots and R will make the prompt for us Warning, this function has side effects! It modifies devAskNewPage!

## Usage

\#\# S3 method for class 'layout' plot(which, pages)

## Arguments

$\begin{array}{ll}\text { which } & \text { which plots are to be created } \\ \text { pages } & \text { number of pages to span the plots across }\end{array}$

## Details

Code is stolen and modified from plot. R in mgcv by Simon Wood

## Author(s)

David L. Miller, based on code by Simon N. Wood

Plot fit of detection functions and histograms of data from removal distance sampling model

## Description

Plots the fitted detection functions for a distance sampling model and histograms of the distances (for unconditional detection functions) or proportion of observations detected within distance intervals (for conditional detection functions) to compare visually the fitted model and data.

## Usage

```
    ## S3 method for class 'rem'
```

    plot
        x,
        which = 1:3,
        breaks = NULL,
        nc = NULL,
        maintitle = "",
        showlines = TRUE,
        showpoints = TRUE,
        ylim \(=c(0,1)\),
        angle = NULL,
        density \(=\) NULL,
        col = "lightgrey",
        jitter = NULL,
        divisions \(=25\),
        pages \(=0\),
        xlab = "Distance",
        ylab = "Detection probability",
        subtitle = TRUE,
    )
    
## Arguments

$x$
which
fitted model from ddf
index to specify which plots should be produced.
1 Plot primary unconditional detection function
2 Plot pooled unconditional detection function
3 Plot conditional (112) detection function
breaks user define breakpoints
nc number of equal-width bins for histogram

| maintitle | main title line for each plot |
| :---: | :---: |
| showlines | logical variable; if TRUE a line representing the average detection probability is plotted |
| showpoints | logical variable; if TRUE plots predicted value for each observation |
| ylim | range of vertical axis; defaults to ( 0,1 ) |
| angle | shading angle for histogram bars. |
| density | shading density for histogram bars. |
| col | colour for histogram bars. |
| jitter | scaling option for plotting points. Jitter is applied to points by multiplying the fitted value by a random draw from a normal distribution with mean 1 and sd jitter. |
| divisions | number of divisions for averaging line values; default $=25$ |
| pages | the number of pages over which to spread the plots. For example, if pages=1 then all plots will be displayed on one page. Default is 0 , which prompts the user for the next plot to be displayed. |
| xlab | label for x -axis |
| ylab | label for y -axis |
| subtitle | if TRUE, shows plot type as sub-title |
|  | other graphical parameters, passed to the plotting functions (plot, hist, lines, points, etc) |

## Details

The structure of the histogram can be controlled by the user-defined arguments nc or breaks. The observation specific detection probabilities along with the line representing the fitted average detection probability.

It is not intended for the user to call plot.rem but its arguments are documented here. Instead the generic plot command should be used and it will call the appropriate function based on the class of the ddf object.

## Author(s)

Jeff Laake, Jon Bishop, David Borchers, David L Miller

## Description

Plots the fitted detection functions for a distance sampling model and histograms of the distances (for unconditional detection functions) or proportion of observations detected within distance intervals (for conditional detection functions) to compare visually the fitted model and data.

```
Usage
    ## S3 method for class 'rem.fi'
    plot(
        x,
        which = 1:3,
        breaks = NULL,
        nc = NULL,
        maintitle = "",
        showlines = TRUE,
        showpoints = TRUE,
        ylim = c(0, 1),
        angle = NULL,
        density = NULL,
        col = "lightgrey",
        jitter = NULL,
        divisions = 25,
        pages = 0,
        xlab = "Distance",
        ylab = "Detection probability",
        subtitle = TRUE,
    ...
    )
```


## Arguments

| which | fitted model from ddf <br> index to specify which plots should be produced. |
| :---: | :---: |
|  | 1 Plot primary unconditional detection function <br> 2 Plot pooled unconditional detection function <br> 3 Plot conditional (1\|2) detection function |
| breaks | user defined breakpoints |
| nc | number of equal-width bins for histogram |
| maintitle | main title line for each plot |
| showlines | logical variable; if TRUE a line representing the average detection probability is plotted |
| showpoints | logical variable; if TRUE plots predicted value for each observation |
| ylim | range of vertical axis; defaults to (0,1) |
| angle | shading angle for histogram bars. |
| density | shading density for histogram bars. |
| col | colour for histogram bars. |
| jitter | scaling option for plotting points. Jitter is applied to points by multiplying the fitted value by a random draw from a normal distribution with mean 1 and sd jitter |


| divisions | number of divisions for averaging line values; default $=25$ <br> pages |
| :--- | :--- |
| the number of pages over which to spread the plots. For example, if pages=1 <br> then all plots will be displayed on one page. Default is 0, which prompts the <br> user for the next plot to be displayed. |  |
| xlab | label for x-axis |
| ylab | label for y-axis |
| subtitle | if TRUE, shows plot type as sub-title |
| $\ldots$ | other graphical parameters, passed to the plotting functions (plot, hist, lines, <br> points, etc) |

## Details

The structure of the histogram can be controlled by the user-defined arguments nc or breaks. The observation specific detection probabilities along with the line representing the fitted average detection probability.
It is not intended for the user to call plot.rem.fi but its arguments are documented here. Instead the generic plot command should be used and it will call the appropriate function based on the class of the ddf object.

## Author(s)

Jeff Laake, Jon Bishop, David Borchers, David L Miller

```
plot.trial
```

Plot fit of detection functions and histograms of data from distance sampling trial observer model

## Description

Plots the fitted detection functions for a distance sampling model and histograms of the distances (for unconditional detection functions) or proportion of observations detected within distance intervals (for conditional detection functions) to compare visually the fitted model and data.

## Usage

```
## S3 method for class 'trial'
```

plot $($
x ,
which = 1:2,
breaks = NULL,
nc = NULL,
maintitle = "",
showlines = TRUE,
showpoints = TRUE,
$y \lim =c(0,1)$,

```
    angle = NULL,
    density = NULL,
    col = "lightgrey",
    jitter = NULL,
    divisions = 25,
    pages = 0,
    xlab = "Distance",
    ylab = "Detection probability",
    subtitle = TRUE,
    ...
)
```


## Arguments

x
fitted model from ddf
index to specify which plots should be produced.
1 Unconditional detection function for observer 1
2 Conditional detection function plot (1|2)
breaks user define breakpoints
nc number of equal-width bins for histogram
maintitle
main title line for each plot
showlines logical variable; if TRUE a line representing the average detection probability is plotted
showpoints logical variable; if TRUE plots predicted value for each observation
ylim range of vertical axis; defaults to $(0,1)$
angle shading angle for histogram bars.
density shading density for histogram bars.
col colour for histogram bars.
jitter scaling option for plotting points. Jitter is applied to points by multiplying the fitted value by a random draw from a normal distribution with mean 1 and sd jitter.
divisions number of divisions for averaging line values; default $=25$
pages the number of pages over which to spread the plots. For example, if pages=1 then all plots will be displayed on one page. Default is 0 , which prompts the user for the next plot to be displayed.
$x l a b \quad$ label for x -axis
ylab label for $y$-axis
subtitle if TRUE, shows plot type as sub-title
other graphical parameters, passed to the plotting functions (plot, hist, lines, points, etc)

## Details

The structure of the histogram can be controlled by the user-defined arguments nc or breaks. The observation specific detection probabilities along with the line representing the fitted average detection probability.
It is not intended for the user to call plot.io.fi but its arguments are documented here. Instead the generic plot command should be used and it will call the appropriate function based on the class of the ddf object.

## Author(s)

Jeff Laake, Jon Bishop, David Borchers
plot.trial.fi Plot fit of detection functions and histograms of data from distance sampling trial observer model

## Description

Plots the fitted detection functions for a distance sampling model and histograms of the distances (for unconditional detection functions) or proportion of observations detected within distance intervals (for conditional detection functions) to compare visually the fitted model and data.

```
Usage
    ## S3 method for class 'trial.fi'
    plot(
        x,
        which = 1:2,
        breaks = NULL,
        nc = NULL,
        maintitle = "",
        showlines = TRUE,
        showpoints = TRUE,
        ylim = c(0, 1),
        angle = NULL,
        density = NULL,
        col = "lightgrey",
        jitter = NULL,
        divisions = 25,
        pages = 0,
        xlab = "Distance",
        ylab = "Detection probability",
        subtitle = TRUE,
        )
```


## Arguments

| x | fitted model from ddf |
| :---: | :---: |
| which | index to specify which plots should be produced. |
|  | 1 Unconditional detection function for observer 1 <br> 2 Conditional detection function plot (1\|2) |
| breaks | user define breakpoints |
| nc | number of equal-width bins for histogram |
| maintitle | main title line for each plot |
| showlines | logical variable; if TRUE a line representing the average detection probability is plotted |
| showpoints | logical variable; if TRUE plots predicted value for each observation |
| ylim | range of vertical axis; defaults to ( 0,1 ) |
| angle | shading angle for histogram bars. |
| density | shading density for histogram bars. |
| col | colour for histogram bars. |
| jitter | scaling option for plotting points. Jitter is applied to points by multiplying the fitted value by a random draw from a normal distribution with mean 1 and sd jitter. |
| divisions | number of divisions for averaging line values; default $=25$ |
| pages | the number of pages over which to spread the plots. For example, if pages=1 then all plots will be displayed on one page. Default is 0 , which prompts the user for the next plot to be displayed. |
| xlab | label for x -axis |
| ylab | label for y -axis |
| subtitle | if TRUE, shows plot type as sub-title |
|  | other graphical parameters, passed to the plotting functions (plot, hist, lines, points, etc) |

## Details

The structure of the histogram can be controlled by the user-defined arguments nc or breaks. The observation specific detection probabilities along with the line representing the fitted average detection probability.
It is not intended for the user to call plot.io.fi but its arguments are documented here. Instead the generic plot command should be used and it will call the appropriate function based on the class of the ddf object.

## Author(s)

Jeff Laake, Jon Bishop, David Borchers
plot_cond
Plot conditional detection function from distance sampling model

## Description

Plot proportion of observations detected within distance intervals (for conditional detection functions) to compare visually the fitted model and data. Internal function called by plot methods.

## Usage

```
    plot_cond(
        obs,
        xmat,
        gxvalues,
        model,
        nc,
        breaks,
        finebr,
        showpoints,
        showlines,
        maintitle,
        ylim,
        angle = -45,
        density = 20,
        col = "black",
        jitter = NULL,
        xlab = "Distance",
        ylab = "Detection probability",
        subtitle = TRUE,
        ..
    )
```


## Arguments

| obs | observer code |
| :--- | :--- |
| xmat | processed data |
| gxvalues | detection function values for each observation |
| model | fitted model from ddf |
| nc | number of equal-width bins for histogram |
| breaks | user define breakpoints |
| finebr | fine break values over which line is averaged |
| showpoints | logical variable; if TRUE plots predicted value for each observation |
| showlines | logical variable; if TRUE plots average predicted value line |
| maintitle | main title line for each plot |


| ylim | range of y axis (default $\mathrm{c}(0,1)$ ) |
| :---: | :---: |
| angle | shading angle for hatching |
| density | shading density for hatching |
| col | plotting colour |
| jitter | scaling option for plotting points. Jitter is applied to points by multiplying the fitted value by a random draw from a normal distribution with mean 1 and sd jitter. |
| xlab | label for x -axis |
| ylab | label for y -axis |
| subtitle | if TRUE, shows plot type as sub-title |
|  | other graphical parameters, passed to the plotting functions (plot, hist, lines, points, etc) |

## Author(s)

Jeff Laake, Jon Bishop, David Borchers
plot_uncond

## Description

Plots unconditional detection function for observer=obs observations overlays histogram, average detection function and values for individual observations data. Internal function called by plot methods.

## Usage

```
    plot_uncond(
        model,
        obs,
        xmat,
        gxvalues,
        nc,
        finebr,
        breaks,
        showpoints,
        showlines,
        maintitle,
        ylim,
        return.lines = FALSE,
        angle = -45,
        density = 20,
        col = "black",
        jitter = NULL,
```

```
    xlab = "Distance",
    ylab = "Detection probability",
    subtitle = TRUE,
)
```


## Arguments

| model | fitted model from ddf |
| :--- | :--- |
| obs | value of observer for plot |
| xmat | processed data |
| gxvalues | detection function values for each observation |
| nc | number of equal-width bins for histogram |
| finebr | fine break values over which line is averaged |
| breaks | user define breakpoints |
| showpoints | logical variable; if TRUE plots predicted value for each observation |
| showlines | logical variable; if TRUE plots average predicted value line |
| maintitle | main title line for each plot |
| ylim | range of y axis; defaults to (0,1) |
| return.lines | if TRUE, returns values for line |
| angle | shading angle for hatching |
| density | shading density for hatching |
| col | plotting colour |
| jitter | scaling option for plotting points. Jitter is applied to points by multiplying the <br> fitted value by a random draw from a normal distribution with mean 1 and sd |
| jitter. |  |
| xlab | label for x-axis |
| ylab | label for y-axis <br> subtitle |
| if TRUE, shows plot type as sub-title |  |

Value
if return.lines==TRUE returns dataframe average. line otherwise just plots

## Author(s)

Jeff Laake, Jon Bishop, David Borchers

```
predict.ds Predictions from mrds models
```


## Description

Predict detection probabilities (or effective strip widths/effective areas of detection) from a fitted distance sampling model using either the original data (i.e. "fitted" values) or using new data.

## Usage

```
\#\# S3 method for class 'ds'
predict(object, newdata=NULL, compute=FALSE,
int.range=NULL, esw=FALSE, se.fit=FALSE, ...)
    \#\# S3 method for class 'io.fi'
    predict(object, newdata=NULL, compute=FALSE,
        int.range=NULL, integrate=FALSE, ...)
        \#\# S3 method for class 'io'
    predict(object, newdata=NULL, compute=FALSE,
            int.range=NULL, ...)
            \#\# S3 method for class 'trial'
    predict(object, newdata=NULL, compute=FALSE,
            int.range=NULL, ...)
            \#\# S3 method for class 'trial.fi'
    predict(object, newdata=NULL, compute=FALSE,
            int.range=NULL, integrate=FALSE, ...)
            \#\# S3 method for class 'rem'
    predict(object, newdata=NULL, compute=FALSE,
            int.range=NULL, ...)
            \#\# S3 method for class 'rem.fi'
predict(object, newdata=NULL, compute=FALSE,
            int.range=NULL, integrate=FALSE, ...)
```


## Arguments

object ddf model object.
newdata new data.frame for prediction, this must include a column called "distance".
compute if TRUE compute values and don't use the fitted values stored in the model object.
int.range integration range for variable range analysis; either vector or 2 column matrix.
esw if TRUE, returns effective strip half-width (or effective area of detection for point transect models) integral from 0 to the truncation distance (width) of $p(y) d y$; otherwise it returns the integral from 0 to truncation width of $p(y) \pi(y)$ where $\pi(y)=1 / w$ for lines and $\pi(y)=2 r / w^{2}$ for points.
se.fit for *.ds models only, generate standard errors on the predicted probabilities of detection (or ESW if esw=TRUE), stored in the se.fit element
... for S3 consistency
integrate for *.fi methods, see Details below.

## Details

The first 4 arguments are the same in each predict function. The latter 2 are specific to certain functions. For line transects, the effective strip half-width (esw=TRUE) is the integral of the fitted detection function over either 0 to W or the specified int. range. The predicted detection probability is the average probability which is simply the integral divided by the distance range. For point transect models, esw=TRUE calculates the effective area of detection (commonly referred to as "nu", this is the integral of $2 /$ width $^{\wedge} 2 * r g(r)$.
Fitted detection probabilities are stored in the model object and these are returned unless compute=TRUE or newdata is specified. compute=TRUE is used to estimate numerical derivatives for use in delta method approximations to the variance.
For method="io.fi" or method="trial.fi" if integrate=FALSE, predict returns the value of the conditional detection probability and if integrate=TRUE, it returns the average conditional detection probability by integrating over x (distance) with respect to a uniform distribution.
Note that the ordering of the returned results when no new data is supplied (the "fitted" values) will not necessarily be the same as the data supplied to ddf, the data (and hence results from predict) will be sorted by object ID (object) then observer ID (observer).

## Value

For all but the exceptions below, the value is a list with a single element: fitted, a vector of average detection probabilities or esw values for each observation in the original data ornewdata

For predict.ds, if se.fit=TRUE there is an additional element \$se.fit, which contains the standard errors of the probabilities of detection or ESW.

For predict.io.fi,predict.trial.fi,predict.rem.fi with integrate=TRUE, the value is a list with one element: fitted, which is a vector of integrated (average) detection probabilities for each observation in the original data or newdata.
For predict.io.fi, predict.trial.fi, or predict.rem.fi with integrate=FALSE, the value is a list with the following elements:
fitted $p(y)$ values
p1 $p_{1 \mid 2}(y)$, conditional detection probability for observer 1
p2 $p_{2 \mid 1}(y)$, conditional detection probability for observer 2
fitted $p .(y)=p_{1 \mid 2}(y)+p_{2 \mid 1}(y)-p_{1 \mid 2}(y) * p_{2 \mid 1}(y)$, conditional detection probability of being seen by either observer

## Note

Each function is called by the generic function predict for the appropriate ddf model object. They can be called directly by the user, but it is typically safest to use predict which calls the appropriate function based on the type of model.

## Author(s)

Jeff Laake, David L Miller

## See Also

ddf, summary.ds, plot.ds

```
print.ddf Simple pretty printer for distance sampling analyses
```


## Description

Simply prints out summary of the model which was fitted. For more detailed information see summary.

## Usage

\#\# S3 method for class 'ddf'
print(x, ...)

## Arguments

x
a ddf object
$\ldots$ not passed through, just for S3 compatibility.

## Author(s)

David L. Miller

```
print.ddf.gof Prints results of goodness of fit tests for detection functions
```


## Description

Provides formatted output for results of goodness of fit tests: chi-square, Kolmogorv-Smirnov and Cramer-von Mises test as appropriate.

## Usage

\#\# S3 method for class 'ddf.gof'
print(x, digits $=3, \ldots$ )

## Arguments

$x \quad$ result of call to ddf. gof
digits number of digits to round chi-squared table values to ... unused unspecified arguments for generic print
print.det.tables

## Value

None

## Author(s)

Jeff Laake

## See Also

 ddf.gof```
print.det.tables Print results of observer detection tables
```


## Description

Provides formatted output for detection tables

## Usage

\#\# S3 method for class 'det.tables' print(x, ...)

## Arguments

x result of call to ddf
... unused unspecified arguments for generic print

## Value

None

## Author(s)

Jeff Laake

## See Also

plot.det.tables

## Description

Outputs summary statistics, abundance and density by region (if any) and optionally a correlation matrix if more than one region.

## Usage

\#\# S3 method for class 'dht' print(x, cor = FALSE, bysample = FALSE, vcmatrices = FALSE, ...)

## Arguments

| $x$ | dht object that results from call to dht for a specific ddf object |
| :--- | :--- |
| cor | if TRUE outputs correlation matrix of estimates |
| bysample | if TRUE, prints results for each sample |
| vcmatrices | if TRUE, prints variance-covariance matrices |
| $\ldots$ | unspecified and unused arguments for S3 consistency |

## Value

None

## Author(s)

Jeff Laake

## See Also

dht
print.p_dist_table Print distribution of probabilities of detection

## Description

Just a pretty printer for the table of probabilities of detection.

## Usage

```
    ## S3 method for class 'p_dist_table'
```

    print(x, digits = 2, ...)
    print.summary.ds

## Arguments

x
output from p_dist_table
digits
number of significant digits to print
. . .
other arguments to be passed to print. data.frame

## Value

just prints the table and the range of ps

## Author(s)

David L Miller
print.summary.ds Print summary of distance detection function model object

## Description

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error. What is printed depends on the corresponding call to summary.

## Usage

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


## Arguments

| $x$ | a summary of ddf model object |
| :--- | :--- |
| $\ldots$ | unspecified and unused arguments for $S 3$ consistency |

## Author(s)

Jeff Laake

## See Also

summary.ds

## Description

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error. What is printed depends on the corresponding call to summary.

## Usage

\#\# S3 method for class 'summary.io' print(x, ...)

## Arguments

x
a summary of ddf model object
... unspecified and unused arguments for S3 consistency

## Author(s)

Jeff Laake

## See Also

summary.io

```
print.summary.io.fi Print summary of distance detection function model object
```


## Description

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error. What is printed depends on the corresponding call to summary.

## Usage

```
## S3 method for class 'summary.io.fi'
print(x, ...)
```


## Arguments

x
...
a summary of ddf model object
unspecified and unused arguments for S3 consistency

## Author(s)

Jeff Laake

## See Also

```
summary.io.fi
```


## Description

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error. What is printed depends on the corresponding call to summary.

## Usage

\#\# S3 method for class 'summary.rem'
print(x, ...)

## Arguments

x
a summary of ddf model object
$\ldots \quad$ unspecified and unused arguments for S3 consistency

## Author(s)

Jeff Laake

## See Also

print.summary.rem.fi Print summary of distance detection function model object

## Description

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error. What is printed depends on the corresponding call to summary.

## Usage

\#\# S3 method for class 'summary.rem.fi'
print(x, ...)

## Arguments

x
a summary of ddf model object
$\ldots \quad$ unspecified and unused arguments for S3 consistency

## Author(s)

Jeff Laake

## See Also

summary.rem.fi

```
print.summary.trial

\section*{Description}

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error. What is printed depends on the corresponding call to summary.

\section*{Usage}
```


## S3 method for class 'summary.trial'

print(x, ...)

```

\section*{Arguments}
x
...
a summary of ddf model object
unspecified and unused arguments for S 3 consistency

\section*{Author(s)}

Jeff Laake

\section*{See Also}
```

summary.trial

```
```

print.summary.trial.fi

```

Print summary of distance detection function model object

\section*{Description}

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error. What is printed depends on the corresponding call to summary.

\section*{Usage}
\#\# S3 method for class 'summary.trial.fi'
print(x, ...)

\section*{Arguments}
x a summary of ddf model object
... unspecified and unused arguments for S3 consistency

\section*{Author(s)}

Jeff Laake

\section*{See Also}
```

summary.trial.fi

```

\section*{Description}

Used in call to DeltaMethod from prob.se to get first derivatives

\section*{Usage}
prob.deriv(par, model, parfct, observer = NULL, fittedmodel = NULL)

\section*{Arguments}
par detection function parameter values
model ddf model object
parfct function of detection probabilities; currently only average (over covariates) detection probability \(p\) integrated over distance or average (over covariates) detection probability at distance \(0 ; p(0)\)
observer \(\quad 1,2,3\) for primary, secondary, or duplicates for average \(p(0)\); passed to fct
fittedmodel full fitted ddf model when trial.fi or io.fi is called from trial or io respectively

\section*{Details}

Need to add equations here as I do not think they exist in any of the texts. These should probably be checked with simulation.

\section*{Value}

Vector of values from fct at specified parameter values

\section*{Author(s)}

Jeff Laake

\section*{See Also}
prob.se
prob.se Average \(p\) and average \(p(0)\) variance

\section*{Description}

Computes components of variance for average \(p=n / N\) and average \(p(0)\) with weights based on empirical covariate distribution, if it contains covariates.

\section*{Usage}
prob.se(model, fct, vcov, observer = NULL, fittedmodel = NULL)

\section*{Arguments}
\begin{tabular}{ll} 
model & ddf model object \\
fct & \begin{tabular}{l} 
function of detection probabilities; currently only average (over covariates) de- \\
tection probability p integrated over distance or average (over covariates) detec- \\
tion probability at distance \(0 ; p(0)\)
\end{tabular} \\
vcov & \begin{tabular}{l} 
variance-covariance matrix of parameter estimates
\end{tabular} \\
observer & \begin{tabular}{l}
\(1,2,3\) for primary, secondary, or duplicates for average \(p(0)\); passed to fct \\
fittedmodel
\end{tabular} \\
\begin{tabular}{l} 
full fitted ddf model when trial.fi or io.fi is called from trial or io re- \\
spectively
\end{tabular}
\end{tabular}

\section*{Details}

Need to add equations here as I do not think they exist in any of the texts. These should probably be checked with simulation.

\section*{Value}
\begin{tabular}{ll} 
var & variance \\
partial & partial derivatives of parameters with respect to fct \\
covar & covariance of \(n\) and average \(p\) or \(p(0)\)
\end{tabular}

\section*{Author(s)}

Jeff Laake

\section*{See Also}
prob.deriv

\section*{Description}

Sets up dataframe and does some basic error checking. Adds needed fields to dataframe and to meta.data.

\section*{Usage}
process.data(data, meta.data \(=\) list(), check \(=\) TRUE)

\section*{Arguments}
\begin{tabular}{ll} 
data & dataframe object \\
meta.data & meta.data options; see ddf for a description \\
check & if TRUE check data for errors in the mrds structure; for method="ds" check=FALSE
\end{tabular}

\section*{Details}

The function does a number of error checking tasks, creating fields and adding to meta.data including:
1) If check=TRUE, check to make sure the record structure is okay for mrds data. The number of primary records (observer=1) must equal the number of secondary records (observer=2). Also, a field in the dataframe is created timesseen which counts the number of times an object was detected \(0,1,2\); if timesseen \(=0\) then the record is tossed from the analysis. Also if there are differences in the data (distance, size, covariates) for observer 1 and 2 a warning is issued that the analysis may fail. The code assumes these values are the same for both observers.
2) Based on the presence of fields distbegin and distend, a determination is made of whether the data analysis should be based on binned distances and a field binned is created, which is TRUE if the distance for the observation is binned. By assigning for each observation this allows an analysis of a mixture of binned and unbinned distances.
4) Data are restricted such that distances are not greater than width and not less than left if those values are specified in meta.data. If they are not specified then left defaults to 0 and width defaults to the largest distance measurement.
5) Determine if an integration range (int.begin and int.end has been specified for the observations. If it has, add the structure to meta.data. The integration range is typically used for aerial surveys in which the altitude varies such that the strip width (left to width) changes with a change in altitude.
6) Fields defined as factors are cleaned up such that any unused levels are eliminated.
7) If the restrictions placed on the data, eliminated all of the data, the function stops with an error message

Value
xmat processed data.frame with added fields
meta.data meta.datalist

\section*{Author(s)}

Jeff Laake
pronghorn Pronghorn aerial survey data from Wyoming

\section*{Description}

Detections of pronghorn from fixed-wing aerial surveys in Southeastern Wyoming using four angular bins defined by strut marks. Illustrates data where altitude above ground level (AGL) varies during the survey.

\section*{Format}

A data frame with 660 observations on the following 5 variables.
STRATUM a numeric vector
direction a factor with levels \(N S\) representing the survey direction
AGL height above ground level
Band a factor with levels A B C D which represent angular bands between breaks at 35.42,44.56,51.52,61.02,70.97 degrees. These angles were set based on selected distance bins based on the target AGL.
cluster number of pronghorn in the observed cluster

\section*{Details}

Each record is an observed cluster of pronghorn. The data provide the stratum for the observation, the direction of travel, the AGL at the time of the observation, the angular bin which contained the center of the pronghorn cluster(group), and the number of pronghorn in the group. The angular bins were defined by a combination of two window and five wing strut marks to define bin cutpoints for perpendicular ground distances of \(0-65,65-90,90-115,115-165\) and 165-265 meters when the plane is \(300^{\prime}\) ( 91.4 meters) above ground level. The inner band is considered a blind region due to obstruction of view beneath the plane; thus th the line is offset 65 meters from underneath the plane.

\section*{Source}

Data provided courtesy of Rich Guenzel of Wyoming Game and Fish.

\section*{References}

Laake, J., R. J. Guenzel, J. L. Bengtson, P. Boveng, M. Cameron, and M. B. Hanson. 2008. Coping with variation in aerial survey protocol for line-transect sampling. Wildlife Research 35:289-298.
ptdata.distance Single observer point count data example from Distance

\section*{Description}

Single observer point count data example from Distance

\section*{Format}

The format is 144 obs of 6 variables: distance: numeric distance from center observer: Factor w/ 2 levels "1","2": \(1212121212 \ldots\) detected: numeric \(0 / 1\) object: sequential object number Sample.Label: point label Region.Label: single region label

\section*{Examples}
```

data(ptdata.distance)
xx <- ddf(dsmodel = ~cds(key="hn", formula = ~1), data = ptdata.distance,
method = "ds", meta.data = list(point=TRUE))
summary(xx)
plot(xx,main="Distance point count data")
ddf.gof(xx)
Regions <- data.frame(Region.Label=1,Area=1)
Samples <- data.frame(Sample.Label=1:30,
Region.Label=rep(1,30),
Effort=rep(1,30))
print(dht(xx,sample.table=Samples,region.table=Regions))

```
ptdata.dual Simulated dual observer point count data

\section*{Description}

Simulated dual observer point count data with detection \(p(0)=0.8 ; h n\) sigma \(=30 ; \mathrm{w}=100\) for both observers with dependency \(y>0\), gamma \(=0.1\)

\section*{Format}

The format is 420 obs of 6 variables: distance: numeric distance from center observer: Factor w/ 2 levels "1","2": \(1212121212 \ldots\) detected: numeric \(0 / 1\) person: Factor with 2 levels A,B pair: Factor with 2 levels "AB" BA" \$ object : sequential object number

\section*{Examples}
```

data(ptdata.dual)
xx <- ddf(mrmodel=~glm(formula=~distance),
dsmodel = ~cds(key="hn", formula = ~1),
data = ptdata.dual, method = "io", meta.data = list(point=TRUE))
summary (xx)
plot(xx,main="Simulated point count data")

```
ptdata. removal Simulated removal observer point count data

\section*{Description}

Simulated removal observer point count data with detection \(p(0)=0.8\); hn sigma=30; w=100 for both observers with dependency \(y>0\), gamma \(=0.1\)

\section*{Format}

The format is 408 obs of 6 variables: distance: numeric distance from center observer: Factor w/ 2 levels "1","2": \(1212121212 \ldots\) detected: numeric \(0 / 1\) person: Factor with 2 levels A,B pair: Factor with 2 levels "AB" BA" object: sequential object number

\section*{Examples}
```

data(ptdata.removal)
xx <- ddf(mrmodel=~glm(formula=~distance),
dsmodel = ~cds(key="hn", formula = ~1),
data = ptdata.removal, method = "rem",
meta.data = list(point=TRUE))
summary (xx)
plot(xx,main="Simulated point count data")

```
    ptdata.single Simulated single observer point count data

\section*{Description}

Simulated single observer point count data with detection \(\mathrm{p}(0)=1 ; \mathrm{hn}\) sigma=30; \(\mathrm{w}=100\)

\section*{Format}

The format is 341 obs of 4 variables: .. \(\$\) distance: numeric distance from center \(\$\) observer: Factor w/ 2 levels "1","2": \(1212121212 \ldots .\). detected: numeric \(0 / 1 \$\) object : sequential object number

\section*{Examples}
```

data(ptdata.single)
xx=ddf(dsmodel = ~cds(key="hn", formula = ~1), data = ptdata.single,
method = "ds", meta.data = list(point=TRUE))
summary(xx)
plot(xx,main="Simulated point count data")

```
qqplot.ddf Quantile-quantile plot and goodness of fit tests for detection functions

\section*{Description}

Constructs a quantile-quantile (Q-Q) plot for fitted model as a graphical check of goodness of fit. Formal goodness of fit testing for detection function models using Kolmogorov-Smirnov and Cramer-von Mises tests. Both tests are based on looking at the quantile-quantile plot produced by qqplot.ddf and deviations from the line \(x=y\).

\section*{Usage}
qqplot. \(\mathrm{ddf}(\) model, \(\mathrm{plot}=\) TRUE, \(n b o o t=100, \mathrm{ks}=\) FALSE,.. )

\section*{Arguments}
model
plot the Q-Q plot be plotted or just report statistics?
nboot number of replicates to use to calculate p-values for the goodness of fit test statistics
ks perform the Kolmogorov-Smirnov test (this involves many bootstraps so can take a while)
... additional arguments passed to plot

\section*{Details}

The Kolmogorov-Smirnov test asks the question "what's the largest vertical distance between a point and the \(y=x\) line?" It uses this distance as a statistic to test the null hypothesis that the samples (EDF and CDF in our case) are from the same distribution (and hence our model fits well). If the deviation between the \(y=x\) line and the points is too large we reject the null hypothesis and say the model doesn't have a good fit.

Rather than looking at the single biggest difference between the \(y=x\) line and the points in the \(\mathrm{Q}-\mathrm{Q}\) plot, we might prefer to think about all the differences between line and points, since there may be many smaller differences that we want to take into account rather than looking for one large deviation. Its null hypothesis is the same, but the statistic it uses is the sum of the deviations from each of the point to the line.
rem.glm

\section*{Value}

A list of goodness of fit related values:
\begin{tabular}{ll} 
edf & matrix of lower and upper empirical distribution function values \\
cdf & fitted cumulative distribution function values \\
ks & list with K-S statistic (Dn) and p-value (p) \\
CvM & list with CvM statistic (W) and p-value (p)
\end{tabular}

\section*{Details}

Note that a bootstrap procedure is required to ensure that the p-values from the procedure are correct as the we are comparing the cumulative distribution function (CDF) and empirical distribution function (EDF) and we have estimated the parameters of the detection function.

\section*{Author(s)}

Jeff Laake, David L Miller

\section*{References}

Burnham, K.P., S.T. Buckland, J.L. Laake, D.L. Borchers, T.A. Marques, J.R.B. Bishop, and L. Thomas. 2004. Further topics in distance sampling. pp: 385-389. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.

\section*{See Also}
ddf.gof, cdf.ds

\section*{Description}

Detection function fitting from mark-recapture data with a removal configuration in which a secondary observer knows what the primary observer detects and detects objects missed by the primary observer. The iterative offset glm/gam uses an offset to compensate for the conditioning on the set of objects seen by either observer (eg 00 those missed by both observers are not included in the analysis. This function is similar to io.glm.

\section*{Usage}
```

    rem.glm(
        datavec,
        fitformula,
        eps \(=1 \mathrm{e}-05\),
        iterlimit = 500,
        GAM = FALSE,
        gamplot = TRUE,
        datavec2
    )
    ```

\section*{Arguments}
datavec dataframe containing records seen by either observer 1 or 2
fitformula logit link formula
eps convergence criterion
iterlimit maximum number of iterations allowed
GAM uses GAM instead of GLM for fitting
gamplot set to TRUE to get a gam plot object if GAM=TRUE
datavec2 dataframe containing all records for observer 1 and observer 2 as in io.glm form; this is used in case there is an observer(not platform effect)

\section*{Details}

The only difference between this function and io.glm is the offset and the data construction because there is only one detection function being estimated for the primary observer. The two functions could be merged.

\section*{Value}
list of class("remglm","glm","lm") or class("remglm","gam")
glmobj GLM or GAM object
offsetvalue offsetvalues from iterative fit
plotobj gam plot object (if GAM \& gamplot==TRUE, else NULL)

\section*{Note}
currently the code in this function for GAMs has been commented out until the remainder of the mrds package will work with GAMs.

\section*{Author(s)}

Jeff Laake
rescale_pars

\section*{References}

Buckland, S.T., J.M. breiwick, K.L. Cattanach, and J.L. Laake. 1993. Estimated population size of the California gray whale. Marine Mammal Science, 9:235-249.

Burnham, K.P., S.T. Buckland, J.L. Laake, D.L. Borchers, T.A. Marques, J.R.B. Bishop, and L. Thomas. 2004. Further topics in distance sampling. pp: 360-363. In: Advanced Distance Sampling, eds. S.T. Buckland, D.R.Anderson, K.P. Burnham, J.L. Laake, D.L. Borchers, and L. Thomas. Oxford University Press.
\begin{tabular}{ll} 
rescale_pars & \begin{tabular}{l} 
Calculate the parameter rescaling for parameters associated with co- \\
variates
\end{tabular}
\end{tabular} variates

\section*{Description}

This will calculate the rescaling needed when covariates to be included in the scale of the detection function are "too big". Based on code from optimx.
```

Usage
rescale_pars(initialvalues, ddfobj)

```

\section*{Arguments}
initialvalues starting values for the optimisation
ddfobj detection function object

\section*{Details}

Derivative-free methods like nlminb are sensitive to the parameters being poorly scaled. This can also cause problems for quasi-Newton methods too (at least, bad scaling won't _help_ the optimisation). So here we rescale the parameters if necessary (unless we already got scaling from control)

\section*{Author(s)}

David L Miller

\section*{Description}

Generate data from a fitted detection function and refit the model

\section*{Usage}
sample_ddf(ds.object)

\section*{Arguments}
ds.object a fitted detection function object

\section*{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.

\section*{Author(s)}

David L. Miller
setbounds Set parameter bounds

\section*{Description}

Set values of lower and upper bounds and check lengths of any user-specified values

\section*{Usage}
setbounds(lowerbounds, upperbounds, initialvalues, ddfobj, width, left)

\section*{Arguments}
\begin{tabular}{ll} 
lowerbounds & vector of lower bounds \\
upperbounds & vector of upper bounds \\
initialvalues & vector of initial parameter estimates \\
ddfobj & distance detection function object \\
width & truncation distance \\
left & left truncation distance
\end{tabular}

\section*{Value}
lower vector of lower bounds
upper vector of upper bounds
setlower logical indicating whether user set lower bounds
setupper logical indicating whether user set upper bounds

\section*{Author(s)}

Jeff Laake
```

setcov

```

\section*{Description}

This function creates a design matrix for the \(g(0)\) or scale covariates using the input model formula. It returns a list which contains 2 elements: 1 ) dim: the dimension (number of columns) of the design matrix, and 2) cov: the constructed design matrix. This function is relatively simple because it uses the built-in function model.matrix which does the majority of the work. This function handles 2 exceptions " \(\sim\).", the null model with 0 columns and " \(\sim 1\) " the intercept only model - a column of 1 s . If a model other than the 2 exceptions is provided, it calls model.matrix to construct the columns. If any of the columns of the design matrix are all 0's the column is removed. This occurs when there is no data for a particular factor.

\section*{Usage}
```

setcov(dmat, model)

```

\section*{Arguments}
```

    dmat data matrix
    model model formula
    ```

\section*{Value}
a design matrix for the specified data and model

\section*{Author(s)}

Jeff Laake

\section*{Description}

For a given detection function, it computes the initial values for the parameters including scale and shape parameters and adjustment function parameters if any. If there are user-defined initial values only the parameters not specified by the user are computed.

\section*{Usage}
setinitial.ds(ddfobj, width, initial, point, left) sethazard(ddfobj, dmat, width, left, point)

\section*{Arguments}
ddfobj distance detection function object
width half-width of transect or radius of point count
initial list of user-defined initial values with possible elements: scale, shape, adjustment
point if TRUE, point count data; otherwise, line transect data
left left truncation
dmat xmat from ddfobj

\section*{Value}
scale vector of initial scale parameter values
shape vector of initial shape parameter values
adjustment vector of initial adjustment function parameter values

\section*{Author(s)}

Jeff Laake, David L Miller

\section*{sim.mix}

Simulation of distance sampling data via mixture models Allows one to simulate line transect distance sampling data using a mixture of half-normal detection functions.

\section*{Description}

Simulation of distance sampling data via mixture models Allows one to simulate line transect distance sampling data using a mixture of half-normal detection functions.
solvecov

\section*{Usage}
sim.mix(n, sigma, mix.prop, width, means \(=0\) )

\section*{Arguments}
\begin{tabular}{ll}
n & number of samples to generate \\
sigma & vector of scale parameters \\
mix.prop & vector of mixture proportions (same length as sigma) \\
width & truncation \\
means & vector of means (used to generate wacky, non-monotonic data)
\end{tabular}

\section*{Value}
distances a vector of distances

\section*{Note}

At the moment this is TOTALLY UNSUPPORTED! Please don't use it for anything important!

\section*{Author(s)}

David Lawrence Miller
solvecov Invert of covariance matrices

\section*{Description}

Tries to invert a matrix by solve. If this fails because of singularity, an eigenvector decomposition is computed, and eigenvalues below 1 /cmax are replaced by \(1 / \mathrm{cmax}\), i.e., cmax will be the corresponding eigenvalue of the inverted matrix.

\section*{Usage}
solvecov(m, cmax \(=1 \mathrm{e}+10)\)

\section*{Arguments}
m
cmax
a numeric symmetric matrix.
a positive value, see above.

\section*{Value}

A list with the following components: inv the inverted matrix, coll TRUE if solve failed because of singularity.

Source
solvecov code was taken from package fpc: Christian Hennig

Author(s)
Christian Hennig

\section*{See Also}
solve, eigen
```

stake77 Wooden stake data from 1977 survey

```

\section*{Description}

Multiple surveys by different observers of a single 1 km transect containing 150 wooden stakes placed randomly throughout a 40 m strip ( 20 m on either side).

\section*{Format}

A data frame with 150 observations on the following 10 variables.
StakeNo unique number for each stake 1-150
PD perpendicular distance at which the stake was placed from the line
Obs1 0/1 whether missed/seen by observer 1
Obs2 \(0 / 1\) whether missed/seen by observer 2
Obs3 0/1 whether missed/seen by observer 3
Obs4 \(0 / 1\) whether missed/seen by observer 4
Obs5 \(0 / 1\) whether missed/seen by observer 5
Obs6 \(0 / 1\) whether missed/seen by observer 6
Obs7 \(0 / 1\) whether missed/seen by observer 7
Obs8 \(0 / 1\) whether missed/seen by observer 8

\section*{Source}

Laake, J. 1978. Line transect estimators robust to animal movement. M.S. Thesis. Utah State University, Logan, Utah. 55p.

\section*{References}

Burnham, K. P., D. R. Anderson, and J. L. Laake. 1980. Estimation of Density from Line Transect Sampling of Biological Populations. Wildlife Monographs:7-202.

\section*{Examples}
```

data(stake77)

# Extract functions for stake data and put in the mrds format

extract.stake <- function(stake,obs){
extract.obs <- function(obs){
example <- subset(stake,eval(parse(text=paste("Obs", obs,"==1", sep=""))),
select="PD")
example$distance <- example$PD
example$object <- 1:nrow(example)
        example$PD <- NULL
return(example)
}
if(obs!="all"){
return(extract.obs(obs=obs))
}else{
example <- NULL
for(i in 1:(ncol(stake)-2)){
df <- extract.obs(obs=i)
df$person <- i
            example <- rbind(example,df)
        }
        example$person <- factor(example$person)
        example$object <- 1:nrow(example)
return(example)
}
}
extract.stake.pairs <- function(stake,obs1,obs2,removal=FALSE){
obs1 <- paste("Obs",obs1,sep="")
obs2 <- paste("Obs",obs2,sep="")
example <- subset(stake,eval(parse(text=paste(obs1,"==1 |",obs2,"==1 ",
sep=""))), select=c("PD",obs1,obs2))
names(example) <- c("distance","obs1","obs2")
detected <- c(example$obs1,example$obs2)
example <- data.frame(object = rep(1:nrow(example),2),
distance = rep(example$distance,2),
                detected = detected,
                observer = c(rep(1,nrow(example)),
                        rep(2,nrow(example))))
    if(removal) example$detected[example\$observer==2] <- 1
return(example)
}

# extract data for observer 1 and fit a single observer model

stakes <- extract.stake(stake77,1)
ds.model <- ddf(dsmodel = ~mcds(key = "hn", formula = ~1), data = stakes,
method = "ds", meta.data = list(width = 20))
plot(ds.model,breaks=seq(0, 20, 2), showpoints=TRUE)
ddf.gof(ds.model)

# extract data from observers 1 and 3 and fit an io model

stkpairs <- extract.stake.pairs(stake77,1,3,removal=FALSE)
io.model <- ddf(dsmodel = ~mcds(key = "hn", formula=~1),

```
```

        mrmodel=~glm(formula=~distance),
        data = stkpairs, method = "io")
    summary(io.model)
par(mfrow=c(3,2))
plot(io.model,breaks=seq(0, 20,2),showpoints=TRUE,new=FALSE)
dev.new()
ddf.gof(io.model)

```
stake78 Wooden stake data from 1978 survey

\section*{Description}

Multiple surveys by different observers of a single 1 km transect containing 150 wooden stakes placed based on expected uniform distribution throughout a 40 m strip ( 20 m on either side).

\section*{Format}

A data frame with 150 observations on the following 13 variables.
StakeNo unique number for each stake 1-150
PD perpendicular distance at which the stake was placed from the line
Obs1 \(0 / 1\) whether missed/seen by observer 1
Obs2 \(0 / 1\) whether missed/seen by observer 2
Obs3 0/1 whether missed/seen by observer 3
Obs4 \(0 / 1\) whether missed/seen by observer 4
Obs5 0/1 whether missed/seen by observer 5
Obs6 \(0 / 1\) whether missed/seen by observer 6
Obs7 \(0 / 1\) whether missed/seen by observer 7
Obs8 \(0 / 1\) whether missed/seen by observer 8
Obs9 \(0 / 1\) whether missed/seen by observer 9
Obs10 0/1 whether missed/seen by observer 10
Obs11 0/1 whether missed/seen by observer 11

\section*{Details}

The 1997 survey was based on a single realization of a uniform distribution. Because it was a single transect and there was no randomization of the distances for each survey, we repeated the experiment and used distances that provided a uniform distribution but randomly sorted the positions along the line so there was no pattern obvious to the observer.

\section*{Source}

Laake, J. 1978. Line transect estimators robust to animal movement. M.S. Thesis. Utah State University, Logan, Utah. 55p.

\section*{References}

Burnham, K. P., D. R. Anderson, and J. L. Laake. 1980. Estimation of Density from Line Transect Sampling of Biological Populations. Wildlife Monographs:7-202.

\section*{Examples}
```

data(stake78)
data(stake77)

# compare distribution of distances for all stakes

hist(stake77$PD)
hist(stake78$PD)

# Extract stake data and put in the mrds format for model fitting.

extract.stake <- function(stake,obs){
extract.obs <- function(obs){
example <- subset(stake, eval(parse(text=paste("Obs",obs,"==1",sep=""))),
select="PD")
example$distance <- example$PD
example$object <- 1:nrow(example)
            example$PD <- NULL
return(example)
}
if(obs!="all"){
return(extract.obs(obs=obs))
}else{
example <- NULL
for(i in 1:(ncol(stake)-2)){
df <- extract.obs(obs=i)
df$person <- i
                example <- rbind(example,df)
            }
            example$person <- factor(example$person)
            example$object <- 1:nrow(example)
return(example)
}
}
extract.stake.pairs <- function(stake,obs1,obs2,removal=FALSE){
obs1 <- paste("Obs",obs1,sep="")
obs2 <- paste("Obs",obs2,sep="")
example <- subset(stake,eval(parse(text=paste(obs1,"==1 |",obs2,"==1 ",
sep=""))), select=c("PD",obs1,obs2))
names(example) <- c("distance","obs1","obs2")
detected <- c(example$obs1,example$obs2)
example <- data.frame(object=rep(1:nrow(example),2),
distance=rep(example$distance, 2),
                                    detected = detected,
                                    observer=c(rep(1,nrow(example)),
                                    rep(2,nrow(example))))
    if(removal) example$detected[example\$observer==2] <- 1
return(example)
}

```
```


# extract data for observer 10 and fit a single observer model

stakes <- extract.stake(stake78,10)
ds.model <- ddf(dsmodel = ~mcds(key = "hn", formula = ~1), data = stakes,
method = "ds", meta.data = list(width = 20))
plot(ds.model,breaks=seq(0,20,2),showpoints=TRUE)
ddf.gof(ds.model)

# extract data from observers 5 and 7 and fit an io model

stkpairs <- extract.stake.pairs(stake78,5,7,removal=FALSE)
io.model <- ddf(dsmodel = ~mcds(key = "hn", formula=~1),
mrmodel=~glm(formula=~distance),
data = stkpairs, method = "io")
summary(io.model)
par(mfrow=c(3,2))
plot(io.model,breaks=seq(0, 20, 2), showpoints=TRUE, new=FALSE)
ddf.gof(io.model)

```
summary.ds Summary of distance detection function model object

\section*{Description}

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error.

\section*{Usage}
\#\# S3 method for class 'ds'
summary (object, se = TRUE, \(\mathrm{N}=\) TRUE, ...)

\section*{Arguments}
object a ddf model object
se if TRUE, computes standard errors
N if TRUE, computes abundance in covered (sampled) region
... unspecified and unused arguments for S3 consistency

\section*{Details}

The argument \(N\) is used to suppress computation of abundance and average detection probability in calls to summarize the ds and either io.fi or trial.fi for summaries of io and trial objects respectively which are composed of a ds model object and a mark-recapture model object. The corresponding print function is called to print the summary results.

\section*{Value}
list of extracted and summarized objects

\section*{Note}

This function is called by the generic function summary for any ddf model object. Each function can be called directly by the user, but it is typically safest to use the generic function summary which calls the appropriate function based on the type of ddf model.

\section*{Author(s)}

Jeff Laake
summary.io Summary of distance detection function model object

\section*{Description}

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error.

\section*{Usage}
\#\# S3 method for class 'io'
summary (object, se = TRUE, ...)

\section*{Arguments}
object a ddf model object
se if TRUE, computes standard errors
... unspecified and unused arguments for S3 consistency

\section*{Details}

The argument \(N\) is used to suppress computation of abundance and average detection probability in calls to summarize the ds and either io.fi or trial.fi for summaries of io and trial objects respectively which are composed of a ds model object and a mark-recapture model object. The corresponding print function is called to print the summary results.

\section*{Value}
list of extracted and summarized objects

\section*{Note}

This function is called by the generic function summary for any ddf model object. Each function can be called directly by the user, but it is typically safest to use the generic function summary which calls the appropriate function based on the type of ddf model.

\section*{Author(s)}

Jeff Laake

\section*{Description}

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error.

\section*{Usage}
\#\# S3 method for class 'io.fi'
summary (object, se = TRUE, \(\mathrm{N}=\) TRUE, fittedmodel \(=\) NULL, ddfobj = NULL, ...)

\section*{Arguments}
\begin{tabular}{ll} 
object & a ddf model object \\
se & if TRUE, computes standard errors \\
N & if TRUE, computes abundance in covered (sampled) region \\
fittedmodel & full fitted model when called from trial or io \\
ddfobj & distance sampling object description \\
\(\ldots\) & unspecified and unused arguments for S3 consistency
\end{tabular}

\section*{Details}

The argument \(N\) is used to suppress computation of abundance and average detection probability in calls to summarize the ds and either io.fi or trial.fi for summaries of io and trial objects respectively which are composed of a ds model object and a mark-recapture model object. The corresponding print function is called to print the summary results.

\section*{Value}
list of extracted and summarized objects

\section*{Note}

This function is called by the generic function summary for any ddf model object. Each function can be called directly by the user, but it is typically safest to use the generic function summary which calls the appropriate function based on the type of ddf model.

\section*{Author(s)}

Jeff Laake

\section*{Description}

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error.

\section*{Usage}
\#\# S3 method for class 'rem'
summary (object, se = TRUE, ...)

\section*{Arguments}
```

    object a ddf model object
    se if TRUE, computes standard errors
    ... unspecified and unused arguments for S3 consistency
    ```

\section*{Details}

The argument \(N\) is used to suppress computation of abundance and average detection probability in calls to summarize the ds and either io.fi or trial.fi for summaries of io and trial objects respectively which are composed of a ds model object and a mark-recapture model object. The corresponding print function is called to print the summary results.

\section*{Value}
list of extracted and summarized objects

\section*{Note}

This function is called by the generic function summary for any ddf model object. Each function can be called directly by the user, but it is typically safest to use the generic function summary which calls the appropriate function based on the type of ddf model.

\section*{Author(s)}

Jeff Laake

\section*{Description}

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error.

\section*{Usage}
\#\# S3 method for class 'rem.fi'
summary(object, se = TRUE, \(\mathrm{N}=\) TRUE, fittedmodel = NULL, ...)

\section*{Arguments}
\begin{tabular}{ll} 
object & a ddf model object \\
se & if TRUE, computes standard errors \\
\(N\) & if TRUE, computes abundance in covered (sampled) region \\
fittedmodel & full fitted model when called from trial or io \\
\(\ldots\) & unspecified and unused arguments for S3 consistency
\end{tabular}

\section*{Details}

The argument \(N\) is used to suppress computation of abundance and average detection probability in calls to summarize the ds and either io.fi or trial.fi for summaries of io and trial objects respectively which are composed of a ds model object and a mark-recapture model object. The corresponding print function is called to print the summary results.

\section*{Value}
list of extracted and summarized objects

\section*{Note}

This function is called by the generic function summary for any ddf model object. Each function can be called directly by the user, but it is typically safest to use the generic function summary which calls the appropriate function based on the type of ddf model.

\section*{Author(s)}

Jeff Laake
```

summary.trial Summary of distance detection function model object

```

\section*{Description}

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error.

\section*{Usage}
\#\# S3 method for class 'trial'
summary (object, se = TRUE, ...)

\section*{Arguments}
```

    object a ddf model object
    se if TRUE, computes standard errors
    ... unspecified and unused arguments for S3 consistency
    ```

\section*{Details}

The argument \(N\) is used to suppress computation of abundance and average detection probability in calls to summarize the ds and either io.fi or trial.fi for summaries of io and trial objects respectively which are composed of a ds model object and a mark-recapture model object. The corresponding print function is called to print the summary results.

\section*{Value}
list of extracted and summarized objects

\section*{Note}

This function is called by the generic function summary for any ddf model object. Each function can be called directly by the user, but it is typically safest to use the generic function summary which calls the appropriate function based on the type of ddf model.

\section*{Author(s)}

Jeff Laake

\section*{Description}

Provides a brief summary of data and fitted detection probability model parameters, model selection criterion, and optionally abundance in the covered (sampled) region and its standard error.

\section*{Usage}
\#\# S3 method for class 'trial.fi'
summary (object, se \(=\) TRUE, \(N=\) TRUE, fittedmodel = NULL, ...)

\section*{Arguments}
\begin{tabular}{ll} 
object & a ddf model object \\
se & if TRUE, computes standard errors \\
N & if TRUE, computes abundance in covered (sampled) region \\
fittedmodel & full fitted model when called from trial or io \\
\(\ldots\) & unspecified and unused arguments for S3 consistency
\end{tabular}

\section*{Details}

The argument \(N\) is used to suppress computation of abundance and average detection probability in calls to summarize the ds and either io.fi or trial.fi for summaries of io and trial objects respectively which are composed of a ds model object and a mark-recapture model object. The corresponding print function is called to print the summary results.

\section*{Value}
list of extracted and summarized objects

\section*{Note}

This function is called by the generic function summary for any ddf model object. Each function can be called directly by the user, but it is typically safest to use the generic function summary which calls the appropriate function based on the type of ddf model.

\section*{Author(s)}

Jeff Laake
survey.region.dht Extrapolate Horvitz-Thompson abundance estimates to entire surveyed region

\section*{Description}

Extrapolate Horvitz-Thompson abundance estimates to entire surveyed region

\section*{Usage}
survey.region.dht(Nhat.by.sample, samples, width, left, point, areas.supplied)

\section*{Arguments}

Nhat. by. sample dataframe of abundance by sample
samples samples table
width truncation width
left left truncation if any
point if TRUE point count otherwise line transect
areas. supplied if TRUE, covered area is extracted from the CoveredArea column of Nhat. by. sample

\section*{Value}

Revised Nhat.by.sample dataframe containing estimates extrapolated to survey region

\section*{Note}

Internal function called by dht and related functions.

\section*{Author(s)}

Jeff Laake and David L Miller
```

test.breaks Test validity for histogram breaks(cutpoints)

```

\section*{Description}

Determines whether user specified breaks for histograms are properly ordered and match the left and right truncation.

\section*{Usage}
test.breaks(breaks, left, width)

\section*{Arguments}
breaks vector of cutpoints (breaks) for distance histogram
left left truncation value
width right truncation value; either radius of point count or half-width of transect

\section*{Value}
vector of breaks modified to be valid if necessary

\section*{Author(s)}

Jeff Laake
```

varn
Compute empirical variance of encounter rate

```

\section*{Description}

Computes one of a series of possible variance estimates for the observed encounter rate for a set of sample measurements (e.g., line lengths) and number of observations per sample.

\section*{Usage}
varn(lvec, nvec, type)
covn(lvec, groups1, groups2, type)

\section*{Arguments}
lvec \(\quad\) vector of sample measurements (e.g., line lengths)
nvec vector of number observed
type choice of variance estimator to use for encounter rate
groups1 vector of number of groups observed
groups2 vector of number of individuals observed

\section*{Details}

The choice of type follows the notation of Fewster et al. (2009) in that there are 8 choices of encounter rate variance that can be computed for lines and one for points:

R2 random line placement with unequal line lengths (design-assisted estimator)
R3 random line placement, model-assisted estimator, based on true contagion process
R4 random line placement, model-assisted estimator, based on apparent contagion process
S1 systematic line placement, post-stratification with no strata overlap

S2 systematic line placement, post-stratification with no strata overlap, variances weighted by line length per stratum
01 systematic line placement, post-stratification with overlapping strata (akin to S1)
02 systematic line placement, post-stratification with overlapping strata (weighted by line length per stratum, akin to S2)
03 systematic line placement, post-stratification with overlapping strata, model-assisted estimator with trend in encounter rate with line length

P2 random point placement, potentially unequal number of visits per point, design-based estimator P3 random point placement, potentially unequal number of visits per point, model-based estimator

Default value is "R2", shown in Fewster et al. (2009) to have good performance for completely random designs for lines. For systematic parallel line transect designs, Fewster et al. recommend " 02 ". For point transects the default is "P3" (but "P2" is also available).
For the systematic estimators, pairs are assigned in the order they are given in the lengths and groups vectors.

\section*{Value}

Variance of encounter rate as defined by arguments

\section*{Note}

This function is also used with different calling arguments to compute Innes et al variance of the estimated abundances/length rather than observation encounter rate. The function covn is probably only valid for R3 and R2. Currently, the R2 form is used for all types other than R3.

\section*{Author(s)}

Jeff Laake, David L Miller

\section*{References}

Fewster, R.M., S.T. Buckland, K.P. Burnham, D.L. Borchers, P.E. Jupp, J.L. Laake and L. Thomas. 2009. Estimating the encounter rate variance in distance sampling. Biometrics 65: 225-236.

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