

# Package ‘ICAOD’

January 14, 2019

**Title** Optimal Designs for Nonlinear Models

**Version** 0.9.8

**Description** Finds optimal designs for nonlinear models using a metaheuristic algorithm called imperialist competitive algorithm ICA. See, for details, Masoudi et al. (2017) <doi:10.1016/j.csda.2016.06.014>.

**Depends** R (>= 3.1.3)

**License** GPL (>= 2)

**LazyData** true

**LinkingTo** Rcpp, RcppEigen,

**Imports** Rcpp, nloptr, stats, graphics, grDevices, cubature, sn,  
mnormt, methods, mvQuad

**Suggests** rgl, lattice, knitr, rmarkdown

**BugReports** <https://github.com/ehsan66/ICAOD/issues>

**URL** <https://github.com/ehsan66/ICAOD>

**RoxygenNote** 6.1.0

**Encoding** UTF-8

**NeedsCompilation** yes

**Author** Ehsan Masoudi [aut, cre],  
Heinz Holling [aut],  
Weng Kee Wong [aut]

**Maintainer** Ehsan Masoudi <ehsan.masoudi@wwu.de>

**Repository** CRAN

**Date/Publication** 2019-01-14 16:30:03 UTC

## R topics documented:

bayes	3
bayescomp	12
beff	16

crt.bayes.control . . . . .	22
crt.minimax.control . . . . .	23
FIM_2par_exp_censor1 . . . . .	24
FIM_2par_exp_censor2 . . . . .	25
FIM_3par_exp_censor1 . . . . .	26
FIM_3par_exp_censor2 . . . . .	26
FIM_exp_2par . . . . .	27
FIM_kinetics_alcohol . . . . .	28
FIM_logistic . . . . .	28
FIM_logistic_2pred . . . . .	29
FIM_logistic_4par . . . . .	30
FIM_loglin . . . . .	31
FIM_mixed_inhibition . . . . .	31
FIM_power_logistic . . . . .	32
FIM_sig_emax . . . . .	33
ICA.control . . . . .	33
ICAOD . . . . .	35
iterate . . . . .	37
iterate.bayes . . . . .	37
iterate.minimax . . . . .	38
leff . . . . .	38
locally . . . . .	40
locallycomp . . . . .	45
minimax . . . . .	49
multiple . . . . .	57
normal . . . . .	61
plot.bayes . . . . .	62
plot.minimax . . . . .	64
print.bayes . . . . .	65
print.minimax . . . . .	66
print.sensbayes . . . . .	66
print.sensminimax . . . . .	67
robust . . . . .	67
sens.bayes.control . . . . .	70
sens.minimax.control . . . . .	72
sensbayes . . . . .	74
sensbayescomp . . . . .	79
senslocally . . . . .	82
senslocallycomp . . . . .	85
sensminimax . . . . .	88
sensmultiple . . . . .	94
sensrobust . . . . .	97
skewnormal . . . . .	100
student . . . . .	101
uniform . . . . .	102

**Description**

Finds (pseudo) Bayesian D-optimal designs for nonlinear models. It should be used when the user assumes a (truncated) prior distribution for the unknown model parameters. If you have a discrete prior, please use the function [robust](#).

**Usage**

```
bayes(formula, predvars, parvars, family = gaussian(), prior, lx, ux,
      iter, k, fimfunc = NULL, ICA.control = list(),
      crt.bayes.control = list(), sens.bayes.control = list(),
      crt_method = c("cubature", "quadrature"), sens_method = c("cubature",
      "quadrature"), initial = NULL, npar = NULL, plot_3d = c("lattice",
      "rgl"), x = NULL)
```

**Arguments**

formula	A nonlinear model <a href="#">formula</a> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <a href="#">formula</a> if necessary.
predvars	A vector of characters. Denotes the predictors in the <a href="#">formula</a> .
parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .
family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <a href="#">gaussian()</a> is applied.
prior	An object of class <code>cprior</code> . User can also use one of the functions <a href="#">uniform</a> , <a href="#">normal</a> , <a href="#">skewnormal</a> or <a href="#">student</a> to create the prior. See 'Details' of <a href="#">bayes</a> .
lx	Vector of lower bounds for the predictors. Should be in the same order as <code>predvars</code> .
ux	Vector of upper bounds for the predictors. Should be in the same order as <code>predvars</code> .
iter	Maximum number of iterations.
k	Number of design points. Must be at least equal to the number of model parameters to avoid singularity of the FIM.
fimfunc	A function. Returns the FIM as a matrix. Required when <code>formula</code> is missing. See 'Details' of <a href="#">minimax</a> .
ICA.control	ICA control parameters. For details, see <a href="#">ICA.control</a> .

<code>crt.bayes.control</code>	A list. Control parameters to approximate the integral in the Bayesian criterion at a given design over the parameter space. For details, see <a href="#">crt.bayes.control</a> .
<code>sens.bayes.control</code>	A list. Control parameters to verify the general equivalence theorem. For details, see <a href="#">sens.bayes.control</a> .
<code>crt_method</code>	A character denotes which method to be used to approximate the integrals in Bayesian criteria. "cubature" corresponds to the adaptive multivariate integration method using the <a href="#">hcubature</a> algorithm (default). "quadrature" corresponds the traditional quadrature formulas and calls the function <a href="#">createNIGrid</a> . The tuning parameters are adjusted by <code>crt.bayes.control</code> . Default is set to "cubature".
<code>sens_method</code>	Similar to <code>crt_method</code> , but for approximating the integrals in the sensitivity (derivative) function. The tuning parameters are adjusted by <code>sens.bayes.control</code> . Default is set to "cubature".
<code>initial</code>	A matrix of the initial design points and weights that will be inserted into the initial solutions (countries) of the algorithm. Every row is a design, i.e. a concatenation of <code>x</code> and <code>w</code> . Will be coerced to a matrix if necessary. See 'Details' of <a href="#">minimax</a> .
<code>npar</code>	Number of model parameters. Used when <code>fimfunc</code> is given instead of <code>formula</code> to specify the number of model parameters. If not specified correctly, the sensitivity (derivative) plot may be shifted below the y-axis. When NULL (default), it will be set to <code>length(parvars)</code> or <code>prior\$npar</code> when missing( <code>formula</code> ).
<code>plot_3d</code>	Which package should be used to plot the sensitivity (derivative) function for two-dimensional design space. Defaults to "lattice".
<code>x</code>	A vector of candidate design (support) points. When is not set to NULL (default), the algorithm only finds the optimal weights for the candidate points in <code>x</code> . Should be set when the user has a finite number of candidate design points and the purpose is to find the optimal weight for each of them (when zero, they will be excluded from the design). For design points with more than one dimension, see 'Details' of <a href="#">sensminimax</a> .

## Details

Let  $\Xi$  be the space of all approximate designs with  $k$  design points (support points) at  $x_1, x_2, \dots, x_k$  from design space  $\chi$  with corresponding weights  $w_1, \dots, w_k$ . Let  $M(\xi, \theta)$  be the Fisher information matrix (FIM) of a  $k$ -point design  $\xi$  and  $\pi(\theta)$  is a user-given prior distribution for the vector of unknown parameters  $\theta$ . A Bayesian D-optimal design  $\xi^*$  minimizes over  $\Xi$

$$\int_{\theta \in \Theta} -\log |M(\xi, \theta)| \pi(\theta) d\theta.$$

An object of class `cprior` is a list with the following components:

- `fn`: Prior distribution as an R function with argument `param` that is the vector of the unknown parameters. See below.
- `npar`: Number of unknown parameters and is equal to the length of `param`.

- lower: Argument lower. It has the same length as param.
- upper: Argument upper. It has the same length as param.

A cprior object will be passed to the argument prior of the function `bayes`. The argument param in fn has the same order as the argument parvars when the model is specified by a formula. Otherwise, it is the same as the argument param in the function `fimfunc`.

The user can use the implemented priors that are `uniform`, `normal`, `skewnormal` and `student` to create a cprior object.

To verify the equivalence theorem of the output design, use `plot` function or change the value of the `checkfreq` in the argument `ICA.control`.

To increase the speed of the algorithm, change the value of the tuning parameters `tol` and `maxEval` via the argument `crt.bayes.control` when `crt_method = "cubature"`. Similarly, this applies when `crt_method = "quadrature"`. In general, if the CPU time matters, the user should find an appropriate speed-accuracy trade-off for her/his own problem. See 'Examples' for more details.

If some of the parameters are known and fixed, they should be set to their values via the argument `paravars` when the model is given by formula. In this case, the user must provide the number of parameters via the argument `npar` for verifying the general equivalence theorem. See 'Examples', Section 'Weibull', 'Richards' and 'Exponential' model.

## Value

an object of class `bayes` that is a list including three sub-lists:

`arg` A list of design and algorithm parameters.

`evol` A list of length equal to the number of iterations that stores the information about the best design (design with the minimum criterion value) of each iteration as follows: `evol[[iter]]` contains:

<code>iter</code>	Iteration number.
<code>x</code>	Design points.
<code>w</code>	Design weights.
<code>min_cost</code>	Value of the criterion for the best imperialist (design).
<code>mean_cost</code>	Mean of the criterion values of all the imperialists.
<code>sens</code>	An object of class 'sensbayes'. See below.

`empires` A list of all the empires of the last iteration.

`alg` A list with following information:

<code>nfeval</code>	Number of function evaluations. It does not count the function evaluations from checking the general equivalence theorem.
<code>nlocal</code>	Number of successful local searches.
<code>nrevol</code>	Number of successful revolutions.
<code>nimprove</code>	Number of successful movements toward the imperialists in the assimilation step.
<code>convergence</code>	Stopped by 'maxiter' or 'equivalence'?

The list `sens` contains information about the design verification by the general equivalence theorem. See `sensbayes` for more Details. It is only given every `ICA.control$checkfreq` iterations and also the last iteration if `ICA.control$checkfreq >= 0`. Otherwise, `NULL`.

## See Also

[sensbayes](#)

## Examples

```
#####
# Two parameter logistic model: uniform prior
#####
# set the uniform prior
uni <- uniform(lower = c(-3, .1), upper = c(3, 2))
# set the logistic model with formula
res1 <- bayes(formula = ~1/(1 + exp(-b *(x - a))),
              predvars = "x", parvars = c("a", "b"),
              family = binomial(), lx = -3, ux = 3,
              k = 5, iter = 1, prior = uni,
              ICA.control = list(rseed = 1366))

## Not run:
res1 <- iterate(res1, 500)
plot(res1)

## End(Not run)
# You can also use your Fisher information matrix (FIM) if you think it is faster!
## Not run:
bayes(fimfunc = FIM_logistic, lx = -3, ux = 3, k = 5, iter = 500,
      prior = uni, ICA.control = list(rseed = 1366))

## End(Not run)

# with fixed x
## Not run:
res1.1 <- bayes(formula = ~1/(1 + exp(-b *(x - a))),
                predvars = "x", parvars = c("a", "b"),
                family = binomial(), lx = -3, ux = 3,
                k = 5, iter = 100, prior = uni,
                x = c(-3, -1.5, 0, 1.5, 3),
                ICA.control = list(rseed = 1366))

plot(res1.1)
# not optimal

## End(Not run)

# with quadrature formula
## Not run:
res1.2 <- bayes(formula = ~1/(1 + exp(-b *(x - a))),
                predvars = "x", parvars = c("a", "b"),
                family = binomial(), lx = -3, ux = 3,
```

```

        k = 5, iter = 1, prior = uni,
        crt_method = "quadrature",
        ICA.control = list(rseed = 1366))
res1.2 <- iterate(res1.2, 500)
plot(res1.2) # not optimal
# compare it with res1 that was found by automatic integration
plot(res1)

# we increase the number of quadrature nodes
res1.3 <- bayes(formula = ~1/(1 + exp(-b *(x - a))),
  predvars = "x", parvars = c("a", "b"),
  family = binomial(), lx = -3, ux = 3,
  k = 5, iter = 1, prior = uni,
  crt_method = "quadrature",
  crt.bayes.control = list(quadrature = list(level = 9)),
  ICA.control = list(rseed = 1366))
res1.3 <- iterate(res1.3, 500)
plot(res1.3)
# by automatic integration (crt_method = "cubature"),
# we did not need to worry about the number of nodes.

## End(Not run)
#####
# Two parameter logistic model: normal prior #1
#####
# defining the normal prior #1
norm1 <- normal(mu = c(0, 1),
  sigma = matrix(c(1, -0.17, -0.17, .5), nrow = 2),
  lower = c(-3, .1), upper = c(3, 2))

## Not run:
# initializing
res2 <- bayes(formula = ~1/(1 + exp(-b *(x - a))), predvars = "x", parvars = c("a", "b"),
  family = binomial(), lx = -3, ux = 3, k = 4, iter = 1, prior = norm1,
  ICA.control = list(rseed = 1366))
res2 <- iterate(res2, 500)
plot(res2)

## End(Not run)

#####
# Two parameter logistic model: normal prior #2
#####
# defining the normal prior #1
norm2 <- normal(mu = c(0, 1),
  sigma = matrix(c(1, 0, 0, .5), nrow = 2),
  lower = c(-3, .1), upper = c(3, 2))

## Not run:
# initializing
res3 <- bayes(formula = ~1/(1 + exp(-b *(x - a))), predvars = "x", parvars = c("a", "b"),
  family = binomial(), lx = -3, ux = 3, k = 4, iter = 1, prior = norm2,
  ICA.control = list(rseed = 1366))

res3 <- iterate(res3, 700)

```

```

plot(res3, sens.bayes.control = list(cubature = list(maxEval = 3000, tol = 1e-4),
                                     optslist = list(maxeval = 3000)))

## End(Not run)

#####
# Two parameter logistic model: skewed normal prior #1
#####
skew1 <- skewnormal(xi = c(0, 1),
                   Omega = matrix(c(1, -0.17, -0.17, .5), nrow = 2),
                   alpha = c(1, 0), lower = c(-3, .1), upper = c(3, 2))

## Not run:
res4 <- bayes(formula = ~1/(1 + exp(-b *(x - a))), predvars = "x", parvars = c("a", "b"),
              family = binomial(), lx = -3, ux = 3, k = 4, iter = 700, prior = skew1,
              ICA.control = list(rseed = 1366, ncount = 60))
plot(res4, sens.bayes.control = list(cubature = list(maxEval = 3000, tol = 1e-4),
                                     optslist = list(maxeval = 3000)))

## End(Not run)

#####
# Two parameter logistic model: skewed normal prior #2
#####
skew2 <- skewnormal(xi = c(0, 1),
                   Omega = matrix(c(1, -0.17, -0.17, .5), nrow = 2),
                   alpha = c(-1, 0), lower = c(-3, .1), upper = c(3, 2))

## Not run:
res5 <- bayes(formula = ~1/(1 + exp(-b *(x - a))), predvars = "x", parvars = c("a", "b"),
              family = binomial(), lx = -3, ux = 3, k = 4, iter = 700, prior = skew2,
              ICA.control = list(rseed = 1366, ncount = 60))
plot(res5, sens.bayes.control = list(cubature = list(maxEval = 3000, tol = 1e-4),
                                     optslist = list(maxeval = 3000)))

## End(Not run)

#####
# Two parameter logistic model: t student prior
#####
# set the prior
stud <- student(mean = c(0, 1), S = matrix(c(1, -0.17, -0.17, .5), nrow = 2),
                df = 3, lower = c(-3, .1), upper = c(3, 2))

## Not run:
res6 <- bayes(formula = ~1/(1 + exp(-b *(x - a))), predvars = "x", parvars = c("a", "b"),
              family = binomial(), lx = -3, ux = 3, k = 5, iter = 500, prior = stud,
              ICA.control = list(ncount = 50, rseed = 1366))

plot(res6)

## End(Not run)
# not bad, but to find a very accurate designs we increase
# the ncount to 200 and repeat the optimization
## Not run:

```

```

res6 <- bayes(formula = ~1/(1 + exp(-b *(x - a))),
  predvars = "x", parvars = c("a", "b"),
  family = binomial(), lx = -3, ux = 3, k = 5, iter = 1000, prior = stud,
  ICA.control = list(ncount = 200, rseed = 1366))

plot(res6)

## End(Not run)

#####
# 4-parameter sigmoid Emax model: uniform prior
#####
lb <- c(4, 11, 100, 5)
ub <- c(8, 15, 130, 9)
## Not run:
res7 <- bayes(formula = ~ theta1 + (theta2 - theta1)*(x^theta4)/(x^theta4 + theta3^theta4),
  predvars = c("x"), parvars = c("theta1", "theta2", "theta3", "theta4"),
  lx = .001, ux = 500, k = 5, iter = 200, prior = uniform(lb, ub),
  ICA.control = list(rseed = 1366, ncount = 60))
plot(res7, sens.bayes.control = list(cubature = list(maxEval = 500, tol = 1e-3),
  optslist = list(maxeval = 500)))

## End(Not run)

#####
# 2-parameter Cox Proportional-Hazards Model for type one censored data
#####
# The Fisher information matrix is available here with name FIM_2par_exp_censor1
# However, we should reparameterize the function to match the standard of the argument 'fimfunc'
myfim <- function(x, w, param)
  FIM_2par_exp_censor1(x = x, w = w, param = param, tcensor = 30)
## Not run:
res8 <- bayes(fimfunc = myfim, lx = 0, ux = 1, k = 4,
  iter = 1, prior = uniform(c(-11, -11), c(11, 11)),
  ICA.control = list(rseed = 1366))

res8 <- iterate(res8, 200)
plot(res8, sens.bayes.control = list(cubature = list(maxEval = 500, tol = 1e-3),
  optslist = list(maxeval = 500)))

## End(Not run)

#####
# 2-parameter Cox Proportional-Hazards Model for random censored data
#####
# The Fisher information matrix is available here with name FIM_2par_exp_censor2
# However, we should reparameterize the function to match the standard of the argument 'fimfunc'
myfim <- function(x, w, param)
  FIM_2par_exp_censor2(x = x, w = w, param = param, tcensor = 30)
## Not run:
res9 <- bayes(fimfunc = myfim, lx = 0, ux = 1, k = 2,
  iter = 200, prior = uniform(c(-11, -11), c(11, 11)),

```

```

ICA.control = list(rseed = 1366))
plot(res9, sens.bayes.control = list(cubature = list(maxEval = 100, tol = 1e-3),
                                     optslist = list(maxeval = 100)))

## End(Not run)

#####
# Weibull model: Uniform prior
#####
# see Dette, H., & Pepelyshev, A. (2008).
# Efficient experimental designs for sigmoidal growth models.
# Journal of statistical planning and inference, 138(1), 2-17.

## See how we fixed a some parameters in Bayesian designs
## Not run:
res10 <- bayes(formula = ~a - b * exp(-lambda * t ^h),
               predvars = c("t"),
               parvars = c("a=1", "b=1", "lambda", "h=1"),
               lx = .00001, ux = 20,
               prior = uniform(.5, 2.5), k = 5, iter = 400,
               ICA.control = list(rseed = 1366))

plot(res10)

## End(Not run)

#####
# Weibull model: Normal prior
#####
norm3 <- normal(mu = 1, sigma = .1, lower = .5, upper = 2.5)
res11 <- bayes(formula = ~a - b * exp(-lambda * t ^h),
               predvars = c("t"),
               parvars = c("a=1", "b=1", "lambda", "h=1"),
               lx = .00001, ux = 20, prior = norm3, k = 4, iter = 1,
               ICA.control = list(rseed = 1366))

## Not run:
res11 <- iterate(res11, 400)
plot(res11)

## End(Not run)

#####
# Richards model: Normal prior
#####
norm4 <- normal(mu = c(1, 1), sigma = matrix(c(.2, 0.1, 0.1, .4), 2, 2),
               lower = c(.4, .4), upper = c(1.6, 1.6))

## Not run:
res12 <- bayes(formula = ~a/(1 + b * exp(-lambda*t))^h,
               predvars = c("t"),
               parvars = c("a=1", "b", "lambda", "h=1"),
               lx = .00001, ux = 10,
               prior = norm4,
               k = 5, iter = 400,

```

```

ICA.control = list(rseed = 1366))
plot(res12, sens.bayes.control = list(cubature = list(maxEval = 1000, tol = 1e-3),
                                     optslist = list(maxeval = 1000)))
## or we can use the quadrature formula to plot the derivative function
plot(res12, sens_method = "quadrature",
      sens.bayes.control = list(optslist = list(maxeval = 1000)))

## End(Not run)

#####
# Exponential model: Uniform prior
#####
res13 <- bayes(formula = ~a + exp(-b*x), predvars = "x",
              parvars = c("a = 1", "b"),
              lx = 0.0001, ux = 1,
              prior = uniform(lower = 1, upper = 20),
              iter = 1, k = 3,
              ICA.control= list(rseed = 100))

## Not run:
res13 <- iterate(res13, 300)
plot(res13)

## End(Not run)

#####
# Power logistic model
#####
# See, Duarte, B. P., & Wong, W. K. (2014).
# A Semidefinite Programming based approach for finding
# Bayesian optimal designs for nonlinear models
uni1 <- uniform(lower = c(-.3, 6, .5), upper = c(.3, 8, 1))
## Not run:
res14 <- bayes(formula = ~1/(1 + exp(-b *(x - a)))^s, predvars = "x",
              parvars = c("a", "b", "s"),
              lx = -1, ux = 1, prior = uni1, k = 5, iter = 1)
res14 <- iterate(res14, 300)
plot(res14)

## End(Not run)

#####
# A two-variable generalized linear model with a gamma distributed response
#####
lb <- c(.5, 0, 0, 0, 0, 0)
ub <- c(2, 1, 1, 1, 1, 1)
myformula1 <- ~beta0+beta1*x1+beta2*x2+beta3*x1^2+beta4*x2^2+beta5*x1*x2
## Not run:
res15 <- bayes(formula = myformula1,
              predvars = c("x1", "x2"), parvars = paste("beta", 0:5, sep = ""),
              family = Gamma(),
              lx = rep(0, 2), ux = rep(1, 2),
              prior = uniform(lower = lb, upper = ub),

```

```

      k = 7, iter = 1, ICA.control = list(rseed = 1366))
res14 <- iterate(res14, 500)
plot(res14, sens.bayes.control = list(cubature = list(maxEval = 5000, tol = 1e-4),
                                     optslist = list(maxeval = 3000)))

## End(Not run)

#####
# Three parameter logistic model
#####

sigma1 <- matrix(-0.1, nrow = 3, ncol = 3)
diag(sigma1) <- c(.5, .4, .1)
norm5 <- normal(mu = c(0, 1, .2), sigma = sigma1,
               lower = c(-3, .1, 0), upper = c(3, 2, .7))
## Not run:
res16 <- bayes(formula = ~ c + (1-c)/(1 + exp(-b *(x - a))), predvars = "x",
               parvars = c("a", "b", "c"),
               family = binomial(), lx = -3, ux = 3,
               k = 4, iter = 500, prior = norm5,
               ICA.control = list(rseed = 1366, ncount = 50),
               crt.bayes.control = list(cubature = list(maxEval = 2500, tol = 1e-4)))
plot(res16, sens.bayes.control = list(cubature = list(maxEval = 3000, tol = 1e-4),
                                     optslist = list(maxeval = 3000)))

# took 925 second on my system

## End(Not run)

```

---

 bayescomp

*Bayesian Compound DP-Optimal Designs*


---

## Description

Finds compound Bayesian DP-optimal designs that meet the dual goal of parameter estimation and increasing the probability of a particular outcome in a binary response model. A compound Bayesian DP-optimal design maximizes the product of the Bayesian efficiencies of a design  $\xi$  with respect to D- and average P-optimality, weighted by a pre-defined mixing constant  $0 \leq \alpha \leq 1$ .

## Usage

```

bayescomp(formula, predvars, parvars, family = binomial(), prior, alpha,
           prob, lx, ux, iter, k, fimfunc = NULL, ICA.control = list(),
           crt.bayes.control = list(), sens.bayes.control = list(),
           crt_method = c("cubature", "quadrature"), sens_method = c("cubature",
                               "quadrature"), initial = NULL, npar = NULL, plot_3d = c("lattice",
                               "rgl"))

```

**Arguments**

formula	A nonlinear model <a href="#">formula</a> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <a href="#">formula</a> if necessary.
predvars	A vector of characters. Denotes the predictors in the <a href="#">formula</a> .
parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .
family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <code>gaussian()</code> is applied.
prior	An object of class <code>cprior</code> . User can also use one of the functions <a href="#">uniform</a> , <a href="#">normal</a> , <a href="#">skewnormal</a> or <a href="#">student</a> to create the prior. See 'Details' of <a href="#">bayes</a> .
alpha	A value between 0 and 1. Compound or combined DP-criterion is the product of the efficiencies of a design with respect to D- and average P- optimality, weighted by alpha.
prob	Either <a href="#">formula</a> or a function. When function, its argument are <code>x</code> and <code>param</code> , and they are the same as the arguments in <code>fimfunc</code> . <code>prob</code> as a function takes the design points and vector of parameters and returns the probability of success at each design points. See 'Examples'.
lx	Vector of lower bounds for the predictors. Should be in the same order as <code>predvars</code> .
ux	Vector of upper bounds for the predictors. Should be in the same order as <code>predvars</code> .
iter	Maximum number of iterations.
k	Number of design points. Must be at least equal to the number of model parameters to avoid singularity of the FIM.
fimfunc	A function. Returns the FIM as a matrix. Required when <code>formula</code> is missing. See 'Details' of <a href="#">minimax</a> .
ICA.control	ICA control parameters. For details, see <a href="#">ICA.control</a> .
crt.bayes.control	A list. Control parameters to approximate the integral in the Bayesian criterion at a given design over the parameter space. For details, see <a href="#">crt.bayes.control</a> .
sens.bayes.control	A list. Control parameters to verify the general equivalence theorem. For details, see <a href="#">sens.bayes.control</a> .
crt_method	A character denotes which method to be used to approximate the integrals in Bayesian criteria. "cubature" corresponds to the adaptive multivariate integration method using the <a href="#">hcubature</a> algorithm (default). "quadrature" corresponds the traditional quadrature formulas and calls the function <a href="#">createNIGrid</a> . The tuning parameters are adjusted by <code>crt.bayes.control</code> . Default is set to "cubature".

<code>sens_method</code>	Similar to <code>crt_method</code> , but for approximating the integrals in the sensitivity (derivative) function. The tuning parameters are adjusted by <code>sens.bayes.control</code> . Default is set to "cubature".
<code>initial</code>	A matrix of the initial design points and weights that will be inserted into the initial solutions (countries) of the algorithm. Every row is a design, i.e. a concatenation of <code>x</code> and <code>w</code> . Will be coerced to a matrix if necessary. See 'Details' of <a href="#">minimax</a> .
<code>npar</code>	Number of model parameters. Used when <code>fimfunc</code> is given instead of <code>formula</code> to specify the number of model parameters. If not specified correctly, the sensitivity (derivative) plot may be shifted below the y-axis. When NULL (default), it will be set to <code>length(parvars)</code> or <code>prior\$npar</code> when missing( <code>formula</code> ).
<code>plot_3d</code>	Which package should be used to plot the sensitivity (derivative) function for two-dimensional design space. Defaults to "lattice".

### Details

Let  $\Xi$  be the space of all approximate designs with  $k$  design points (support points) at  $x_1, x_2, \dots, x_k$  from design space  $\chi$  with corresponding weights  $w_1, \dots, w_k$ . Let  $M(\xi, \theta)$  be the Fisher information matrix (FIM) of a  $k$ -point design  $\xi$ ,  $\pi(\theta)$  is a user-given prior distribution for the vector of unknown parameters  $\theta$  and  $p(x_i, \theta)$  is the  $i$ th probability of success given by  $x_i$  in a binary response model. A compound Bayesian DP-optimal design maximizes over  $\Xi$

$$\int_{\theta \in \Theta} \frac{\alpha}{q} \log |M(\xi, \theta)| + (1 - \alpha) \log \left( \sum_{i=1}^k w_i p(x_i, \theta) \right) \pi(\theta) d\theta.$$

To verify the equivalence theorem of the output design, use `plot` function or change the value of the `checkfreq` in the argument `ICA.control`.

To increase the speed of the algorithm, change the value of the tuning parameters `tol` and `maxEval` via the argument `crt.bayes.control` when `crt_method = "cubature"`. In general, if the CPU time matters, the user should find an appropriate speed-accuracy trade-off for her/his own problem. See 'Examples' for more details.

### Value

an object of class `bayes` that is a list including three sub-lists:

`arg` A list of design and algorithm parameters.

`evol` A list of length equal to the number of iterations that stores the information about the best design (design with the minimum criterion value) of each iteration as follows: `evol[[iter]]` contains:

<code>iter</code>	Iteration number.
<code>x</code>	Design points.
<code>w</code>	Design weights.
<code>min_cost</code>	Value of the criterion for the best imperialist (design).
<code>mean_cost</code>	Mean of the criterion values of all the imperialists.
<code>sens</code>	An object of class 'sensbayes'. See below.

empires A list of all the empires of the last iteration.

alg A list with following information:

nfeval	Number of function evaluations. It does not count the function evaluations from checking the general equivalence theorem.
nlocal	Number of successful local searches.
nrevol	Number of successful revolutions.
nimprove	Number of successful movements toward the imperialists in the assimilation step.
convergence	Stopped by 'maxiter' or 'equivalence'?

The list sens contains information about the design verification by the general equivalence theorem. See sensbayes for more Details. It is only given every ICA.control\$checkfreq iterations and also the last iteration if ICA.control\$checkfreq >= 0. Otherwise, NULL.

## References

McGree, J. M., Eccleston, J. A., and Duffull, S. B. (2008). Compound optimal design criteria for nonlinear models. *Journal of Biopharmaceutical Statistics*, 18(4), 646-661.

## See Also

[sensbayescomp](#)

## Examples

```
#####
# DP-optimal design for a logitic model with two predictors: with formula
#####
p <- c(1, -2, 1, -1)
myprior <- uniform(p -1.5, p + 1.5)
myformula1 <- ~exp(b0+b1*x1+b2*x2+b3*x1*x2)/(1+exp(b0+b1*x1+b2*x2+b3*x1*x2))
res1 <- bayescomp(formula = myformula1,
                  predvars = c("x1", "x2"),
                  parvars = c("b0", "b1", "b2", "b3"),
                  family = binomial(),
                  lx = c(-1, -1), ux = c(1, 1),
                  prior = myprior, iter = 1, k = 7,
                  prob = ~1-1/(1+exp(b0 + b1 * x1 + b2 * x2 + b3 * x1 * x2)),
                  alpha = .5, ICA.control = list(rseed = 1366))

## Not run:
res1 <- iterate(res1, 1000)
plot(res1, sens.bayes.control = list(cubature = list(tol = 1e-3, maxEval = 1000)))
# or use quadrature method
plot(res1, sens_method = "quadrature")

## End(Not run)

#####
# DP-optimal design for a logitic model with two predictors: with fimfunc
```

```
#####
# The function of the Fisher information matrix for this model is 'FIM_logistic_2pred'
# We should reparameterize it to match the standard of the argument 'fimfunc'
myfim <- function(x, w, param){
  npoint <- length(x)/2
  x1 <- x[1:npoint]
  x2 <- x[(npoint+1):(npoint*2)]
  FIM_logistic_2pred(x1 = x1,x2 = x2, w = w, param = param)
}

## The following function is equivalent to the function created
# by the formula:  $\sim 1-1/(1+\exp(b_0 + b_1 * x_1 + b_2 * x_2 + b_3 * x_1 * x_2))$ 
# It returns probability of success given x and param
# x = c(x1, x2) and param = c()

myprob <- function(x, param){
  npoint <- length(x)/2
  x1 <- x[1:npoint]
  x2 <- x[(npoint+1):(npoint*2)]
  b0 <- param[1]
  b1 <- param[2]
  b2 <- param[3]
  b3 <- param[4]
  out <- 1-1/(1+exp(b0 + b1 * x1 + b2 * x2 + b3 * x1 * x2))
  return(out)
}
## Not run:
res2 <- bayescomp(fimfunc = myfim,
                  lx = c(-1, -1), ux = c(1, 1),
                  prior = myprior, iter = 1000, k = 7,
                  prob = myprob, alpha = .5,
                  ICA.control = list(rseed = 1366))
plot(res2, sens.bayes.control = list(cubature = list(maxEval = 1000, tol = 1e-4)))
# quadrature with 6 nodes (default)
plot(res2, sens_method = "quadrature")

## End(Not run)
```

**Description**

Given a prior distribution for the parameters, this function calculates the Bayesian D- and PA- efficiency of a design  $\xi_1$  with respect to a design  $\xi_2$ . Usually,  $\xi_2$  is an optimal design. This function is especially useful for investigating the robustness of Bayesian optimal designs under different prior distributions (See 'Examples').

**Usage**

```
beff(formula, predvars, parvars, family = gaussian(), prior,
     fimfunc = NULL, xopt, wopt, x, w, crt.bayes.control = list(),
     npar = NULL, type = c("D", "PA"), prob = NULL)
```

**Arguments**

formula	A nonlinear model <a href="#">formula</a> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <a href="#">formula</a> if necessary.
predvars	A vector of characters. Denotes the predictors in the <a href="#">formula</a> .
parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .
family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <code>gaussian()</code> is applied.
prior	An object of class <code>cprior</code> . User can also use one of the functions <a href="#">uniform</a> , <a href="#">normal</a> , <a href="#">skewnormal</a> or <a href="#">student</a> to create the prior. See 'Details' of <a href="#">bayes</a> .
fimfunc	A function. Returns the FIM as a matrix. Required when <code>formula</code> is missing. See 'Details' of <a href="#">minimax</a> .
xopt	Vector of design (support) points of the optimal design ( $\xi_2$ ). Similar to <code>x</code> .
wopt	Vector of corresponding design weights for <code>xopt</code> .
x	Vector of design (support) points of $\xi_1$ . See 'Details' of <a href="#">leff</a> .
w	Vector of corresponding design weights for <code>x</code> .
crt.bayes.control	A list. Control parameters to approximate the integral in the Bayesian criterion at a given design over the parameter space. For details, see <a href="#">crt.bayes.control</a> .
npar	Number of model parameters. Used when <code>fimfunc</code> is given instead of <code>formula</code> to specify the number of model parameters. If not given, the sensitivity plot may be shifted below the y-axis. When <code>NULL</code> , it will be set here to <code>length(inipars)</code> .
type	A character. "D" denotes the D-efficiency and "PA" denotes the average P-efficiency.
prob	Either <code>formula</code> or a function. When function, its argument are <code>x</code> and <code>param</code> , and they are the same as the arguments in <code>fimfunc</code> . <code>prob</code> as a function takes the design points and vector of parameters and returns the probability of success at each design points. See 'Examples'.

**Details**

See Masoudi et al. (2018) for formula details (the paper is under review and will be updated as soon as accepted).

The argument  $x$  is the vector of design points. For design points with more than one dimension (the models with more than one predictors), it is a concatenation of the design points, but **dimension-wise**. For example, let the model has three predictors  $(I, S, Z)$ . Then, a two-point optimal design has the following points:  $\{\text{point1} = (I_1, S_1, Z_1), \text{point2} = (I_2, S_2, Z_2)\}$ . Then, the argument  $x$  is equal to  $x = c(I1, I2, S1, S2, Z1, Z2)$ .

## Examples

```
#####
# 2PL model
#####
formula4.1 <- ~ 1/(1 + exp(-b *(x - a)))
predvars4.1 <- "x"
parvars4.1 <- c("a", "b")

# des4.1 is a list of Bayesian optimal designs with corresponding priors.

des4.1 <- vector("list", 6)
des4.1[[1]]$x <- c(-3, -1.20829, 0, 1.20814, 3)
des4.1[[1]]$w <- c(.24701, .18305, .13988, .18309, .24702)
des4.1[[1]]$prior <- uniform(lower = c(-3, .1), upper = c(3, 2))

des4.1[[2]]$x <- c(-2.41692, -1.16676, .04386, 1.18506, 2.40631)
des4.1[[2]]$w <- c(.26304, .18231, .14205, .16846, .24414)
des4.1[[2]]$prior <- student(mean = c(0, 1), S = matrix(c(1, -0.17, -0.17, .5), nrow = 2),
                             df = 3, lower = c(-3, .1), upper = c(3, 2))

des4.1[[3]]$x <- c(-2.25540, -.76318, .54628, 2.16045)
des4.1[[3]]$w <- c(.31762, .18225, .18159, .31853)
des4.1[[3]]$prior <- normal(mu = c(0, 1),
                             sigma = matrix(c(1, -0.17, -0.17, .5), nrow = 2),
                             lower = c(-3, .1), upper = c(3, 2))

des4.1[[4]]$x <- c(-2.23013, -.66995, .67182, 2.23055)
des4.1[[4]]$w <- c(.31420, .18595, .18581, .31404)
des4.1[[4]]$prior <- normal(mu = c(0, 1),
                             sigma = matrix(c(1, 0, 0, .5), nrow = 2),
                             lower = c(-3, .1), upper = c(3, 2))

des4.1[[5]]$x <- c(-1.51175, .12043, 1.05272, 2.59691)
des4.1[[5]]$w <- c(.37679, .14078, .12676, .35567)
des4.1[[5]]$prior <- skewnormal(xi = c(0, 1),
                                 Omega = matrix(c(1, -0.17, -0.17, .5), nrow = 2),
                                 alpha = c(1, 0), lower = c(-3, .1), upper = c(3, 2))

des4.1[[6]]$x <- c(-2.50914, -1.16780, -.36904, 1.29227)
des4.1[[6]]$w <- c(.35767, .11032, .15621, .37580)
des4.1[[6]]$prior <- skewnormal(xi = c(0, 1),
                                 Omega = matrix(c(1, -0.17, -0.17, .5), nrow = 2),
                                 alpha = c(-1, 0), lower = c(-3, .1), upper = c(3, 2))
```

```

## now we want to find the relative efficiency of
## all Bayesian optimal designs assuming different priors (6 * 6)
eff4.1 <- matrix(NA, 6, 6)
colnames(eff4.1) <- c("uni", "t", "norm1", "norm2", "skew1", "skew2")
rownames(eff4.1) <- colnames(eff4.1)
## Not run:
for (i in 1:6)
  for(j in 1:6)
    eff4.1[i, j] <- beff(formula = formula4.1,
                        predvars = predvars4.1,
                        parvars = parvars4.1,
                        family = binomial(),
                        prior = des4.1[[i]]$prior,
                        xopt = des4.1[[i]]$x,
                        wopt = des4.1[[i]]$w,
                        x = des4.1[[j]]$x,
                        w = des4.1[[j]]$w)

# For example the first row represents Bayesian D-efficiencies of different
# Bayesian optimal design found assuming different priors with respect to
# the Bayesian D-optimal design found under uniform prior distribution.

## End(Not run)

#####
# Relative efficiency for the DP-Compound criterion
#####
p <- c(1, -2, 1, -1)
prior4.4 <- uniform(p -1.5, p + 1.5)
formula4.4 <- ~exp(b0+b1*x1+b2*x2+b3*x1*x2)/(1+exp(b0+b1*x1+b2*x2+b3*x1*x2))
prob4.4 <- ~1-1/(1+exp(b0 + b1 * x1 + b2 * x2 + b3 * x1 * x2))
predvars4.4 <- c("x1", "x2")
parvars4.4 <- c("b0", "b1", "b2", "b3")
lb <- c(-1, -1)
ub <- c(1, 1)

## des4.4 is a list of DP-optimal designs found using different values for alpha
des4.4 <- vector("list", 5)
des4.4[[1]]$x <- c(-1, 1)
des4.4[[1]]$w <- c(1)
des4.4[[1]]$alpha <- 0

des4.4[[2]]$x <- c(1, -.62534, .11405, -1, 1, .28175, -1, -1, 1, -1, -1, 1, 1, .09359)
des4.4[[2]]$w <- c(.08503, .43128, .01169, .14546, .05945, .08996, .17713)
des4.4[[2]]$alpha <- .25

des4.4[[3]]$x <- c(-1, .30193, 1, 1, .07411, -1, -.31952, -.08251, 1, -1, 1, -1, -1, 1)
des4.4[[3]]$w <- c(.09162, .10288, .15615, .13123, .01993, .22366, .27454)
des4.4[[3]]$alpha <- .5

```

```

des4.4[[4]]$x <- c(1, -1, .28274, 1, -1, -.19674, .03288, 1, -1, 1, -1, -.16751, 1, -1)
des4.4[[4]]$w <- c(.19040, .24015, .10011, .20527, .0388, .20075, .02452)
des4.4[[4]]$alpha <- .75

des4.4[[5]]$x <- c(1, -1, .26606, -.13370, 1, -.00887, -1, 1, -.2052, 1, 1, -1, -1, -1)
des4.4[[5]]$w <- c(.23020, .01612, .09546, .16197, .23675, .02701, .2325)
des4.4[[5]]$alpha <- 1

# D-efficiency of the DP-optimal designs:
# des4.4[[5]]$x and des4.4[[5]]$w is the D-optimal design

beff(formula = formula4.4,
      predvars = predvars4.4,
      parvars = parvars4.4,
      family = binomial(),
      prior = prior4.4,
      xopt = des4.4[[5]]$x,
      wopt = des4.4[[5]]$w,
      x = des4.4[[2]]$x,
      w = des4.4[[2]]$w)

beff(formula = formula4.4,
      predvars = predvars4.4,
      parvars = parvars4.4,
      family = binomial(),
      prior = prior4.4,
      xopt = des4.4[[5]]$x,
      wopt = des4.4[[5]]$w,
      x = des4.4[[3]]$x,
      w = des4.4[[3]]$w)

beff(formula = formula4.4,
      predvars = predvars4.4,
      parvars = parvars4.4,
      family = binomial(),
      prior = prior4.4,
      xopt = des4.4[[5]]$x,
      wopt = des4.4[[5]]$w,
      x = des4.4[[4]]$x,
      w = des4.4[[4]]$w)

# must be one!
beff(formula = formula4.4,
      predvars = predvars4.4,
      parvars = parvars4.4,
      family = binomial(),
      prior = prior4.4,
      prob = prob4.4,
      type = "PA",
      xopt = des4.4[[5]]$x,
      wopt = des4.4[[5]]$w,
      x = des4.4[[5]]$x,

```

```
w = des4.4[[5]]$w)

## P-efficiency
# reported in Table 4 as eff_P
# des4.4[[1]]$x and des4.4[[1]]$w is the P-optimal design
beff(formula = formula4.4,
      predvars = predvars4.4,
      parvars = parvars4.4,
      family = binomial(),
      prior = prior4.4,
      prob = prob4.4,
      type = "PA",
      xopt = des4.4[[1]]$x,
      wopt = des4.4[[1]]$w,
      x = des4.4[[2]]$x,
      w = des4.4[[2]]$w)

beff(formula = formula4.4,
      predvars = predvars4.4,
      parvars = parvars4.4,
      family = binomial(),
      prior = prior4.4,
      prob = prob4.4,
      type = "PA",
      xopt = des4.4[[1]]$x,
      wopt = des4.4[[1]]$w,
      x = des4.4[[3]]$x,
      w = des4.4[[3]]$w)

beff(formula = formula4.4,
      predvars = predvars4.4,
      parvars = parvars4.4,
      family = binomial(),
      prior = prior4.4,
      prob = prob4.4,
      type = "PA",
      xopt = des4.4[[1]]$x,
      wopt = des4.4[[1]]$w,
      x = des4.4[[4]]$x,
      w = des4.4[[4]]$w)

beff(formula = formula4.4,
      predvars = predvars4.4,
      parvars = parvars4.4,
      family = binomial(),
      prior = prior4.4,
      prob = prob4.4,
      type = "PA",
      xopt = des4.4[[1]]$x,
      wopt = des4.4[[1]]$w,
      x = des4.4[[5]]$x,
      w = des4.4[[5]]$w)
```

---

crt.bayes.control      *Control Parameters for Approximating Bayesian Criteria*

---

### Description

This function returns two lists each corresponds to an implemented integration method for approximating the integrals in Bayesian criteria. The first list is named `cubature` and contains the `hcubature` control parameters to approximate the integrals with an adaptive multivariate integration method over hypercubes. The second list is named `quadrature` and contains the `createNIGrid` tuning parameters to approximate the integrals with the quadrature methods.

### Usage

```
crt.bayes.control(cubature = list(tol = 1e-05, maxEval = 50000, absError
  = 0), quadrature = list(type = "GLe", level = 6, ndConstruction =
  "product", level.trans = FALSE))
```

### Arguments

- |                         |  |
|-------------------------|--|
| <code>cubature</code>   | A list that will be passed to the arguments of the function <code>hcubature</code> for the adaptive multivariate integration. It is required and used when <code>crt_method = "cubature"</code> in the parent function, e.g. <code>bayes</code> . See 'Details'. |
| <code>quadrature</code> | A list that will be passed to the arguments of the function <code>createNIGrid</code> for the quadrature-based integration. It is required and used when <code>crt_method = "quadrature"</code> in the parent function, e.g. <code>bayes</code> . See 'Details'. |

### Details

`cubature` is a list that its components will be passed to the function `hcubature` and they are:

`tol` The maximum tolerance. Defaults to `1e-5`.

`maxEval` The maximum number of function evaluations needed. Note that the actual number of function evaluations performed is only approximately guaranteed not to exceed this number. Defaults to `5000`.

`absError` The maximum absolute error tolerated. Defaults to `0`.

One can specify a maximum number of function evaluations. Otherwise, the integration stops when the estimated error is less than the absolute error requested, or when the estimated error is less than `tol` times the absolute value of the integral, or when the maximum number of iterations is reached, whichever is earlier. `cubature` is activated when `crt_method = "cubature"` in any of the parent functions (for example, `bayes`).

`quadrature` is a list that its components will be passed to the function `createNIGrid` and they are:

**type** Quadrature rule (see Details of `createNIGrid`) Defaults to "GLe".  
**level** Accuracy level (typically number of grid points for the underlying 1D quadrature rule). Defaults to 6.  
**ndConstruction** Character vector which denotes the construction rule for multidimensional grids. "product" for product rule, returns a full grid (default). "sparse" for combination technique, leads to a regular sparse grid.  
**level.trans** Logical variable denotes either to take the levels as number of grid points (FALSE = default) or to transform in that manner that number of grid points =  $2^{(\text{levels}-1)}$  (TRUE). See, `createNIGrid`, for details.  
 quadrature is activated when `crt_method = "quadrature"` in any of the parent functions (for example, `bayes`).

### Value

A list of two lists each contains the control parameters for `hcubature` and `createNIGrid`, respectively.

### Examples

```

crt.bayes.control()
crt.bayes.control(cubature = list(tol = 1e-4))
crt.bayes.control(quadrature = list(level = 4))
  
```

---

`crt.minimax.control`     *Control Parameters for Optimizing Minimax Criteria Over The Parameter Space*

---

### Description

The function `crt.minimax.control` returns a list of `nloptr` control parameters for optimizing the minimax criterion over the parameter space.

The key tuning parameter for our application is `maxeval`. Its value should be increased when either the dimension or the size of the parameter space becomes larger to avoid pre-mature convergence in the inner optimization problem over the parameter space. If the CPU time matters, the user should find an appropriate speed-accuracy trade-off for her/his own design problem.

### Usage

```

crt.minimax.control(x0 = NULL, optslist = list(stopval = -Inf,
  algorithm = "NLOPT_GN_DIRECT_L", xtol_rel = 1e-05, ftol_rel = 1e-08,
  maxeval = 1000), ...)
  
```

**Arguments**

- `x0` Vector of the starting values for the optimization problem (must be from the parameter space).
- `optslist` A list. It will be passed to the argument `opts` of the function `nloptr`. See 'Details'.
- `...` Further arguments will be passed to `nl.opts` from package `nloptr`.

**Details**

Argument `optslist` will be passed to the argument `opts` of the function `nloptr`:

`stopval` Stop minimization when an objective value  $\leq$  `stopval` is found. Setting `stopval` to `-Inf` disables this stopping criterion (default).

`algorithm` Defaults to `NLOPT_GN_DIRECT_L`. `DIRECT-L` is a deterministic-search algorithm based on systematic division of the search domain into smaller and smaller hyperrectangles.

`xtol_rel` Stop when an optimization step (or an estimate of the optimum) changes every parameter by less than `xtol_rel` multiplied by the absolute value of the parameter. Criterion is disabled if `xtol_rel` is non-positive. Defaults to `1e-5`.

`ftol_rel` Stop when an optimization step (or an estimate of the optimum) changes the objective function value by less than `ftol_rel` multiplied by the absolute value of the function value. Criterion is disabled if `ftol_rel` is non-positive. Defaults to `1e-8`.

`maxeval` Stop when the number of function evaluations exceeds `maxeval`. Criterion is disabled if `maxeval` is non-positive. Defaults to `1000`. See below.

A detailed explanation of all the options is shown by `nloptr.print.options()` in package `nloptr`.

**Value**

A list of control parameters for the function `nloptr`.

**Examples**

```
crt.minimax.control(optslist = list(maxeval = 2000))
```

---

FIM\_2par\_exp\_censor1 *Fisher Information Matrix for a 2-Parameter Cox Proportional-Hazards Model for Type One Censored Data*

---

**Description**

It provides the `cpp` function for the FIM introduced in Eq. (3.1) of Schmidt and Schwabe (2015) for type one censored data.

**Usage**

```
FIM_2par_exp_censor1(x, w, param, tcensor)
```

**Arguments**

x	Vector of design points.
w	Vector of design weight. Its length must be equal to the length of x and $\text{sum}(w) = 1$ .
param	Vector of values for the model parameters $c(\beta_0, \beta_1)$ .
tcensor	The experiment is terminated at the fixed time point tcensor.

**Value**

Fisher information matrix.

**References**

Schmidt, D., & Schwabe, R. (2015). On optimal designs for censored data. *Metrika*, 78(3), 237-257.

---

FIM\_2par\_exp\_censor2 *Fisher Information Matrix for a 2-Parameter Cox Proportional-Hazards Model for Random Censored Data*

---

**Description**

It provides the cpp function for the FIM introduced in Eq. (3.1) of Schmidt and Schwabe (2015) for random censored data (type two censored data).

**Usage**

```
FIM_2par_exp_censor2(x, w, param, tcensor)
```

**Arguments**

x	Vector of design points.
w	Vector of design weight. Its length must be equal to the length of x and $\text{sum}(w) = 1$ .
param	Vector of values for the model parameters $c(\beta_0, \beta_1)$ .
tcensor	The experiment is terminated at the fixed time point tcensor.

**Value**

Fisher information matrix.

**References**

Schmidt, D., & Schwabe, R. (2015). On optimal designs for censored data. *Metrika*, 78(3), 237-257.

---

FIM\_3par\_exp\_censor1 *Fisher Information Matrix for a 3-Parameter Cox Proportional-Hazards Model for Type One Censored Data*

---

### Description

It provides the cpp function for the FIM introduced in Page 247 of Schmidt and Schwabe (2015) for type one censored data.

### Usage

```
FIM_3par_exp_censor1(x, w, param, tcensor)
```

### Arguments

x	Vector of design points.
w	Vector of design weight. Its length must be equal to the length of x and $\text{sum}(w) = 1$ .
param	Vector of values for the model parameters $c(\beta_0, \beta_1, \beta_2)$ .
tcensor	The experiment is terminated at the fixed time point tcensor.

### Value

Fisher information matrix.

### References

Schmidt, D., & Schwabe, R. (2015). On optimal designs for censored data. *Metrika*, 78(3), 237-257.

---

FIM\_3par\_exp\_censor2 *Fisher Information Matrix for a 3-Parameter Cox Proportional-Hazards Model for Random Censored Data*

---

### Description

It provides the cpp function for the FIM introduced in Page 247 of Schmidt and Schwabe (2015) for random censored data (type two censored data).

### Usage

```
FIM_3par_exp_censor2(x, w, param, tcensor)
```

**Arguments**

x	Vector of design points.
w	Vector of design weight. Its length must be equal to the length of x and $\text{sum}(w) = 1$ .
param	Vector of values for the model parameters $(\beta_0, \beta_1, \beta_2)$ .
tcensor	The experiment is terminated at the fixed time point tcensor.

**Value**

Fisher information matrix.

**References**

Schmidt, D., & Schwabe, R. (2015). On optimal designs for censored data. *Metrika*, 78(3), 237-257.

---

FIM\_exp\_2par

*Fisher Information Matrix for the 2-Parameter Exponential Model*


---

**Description**

It provides the cpp function for FIM for the model  $\sim a + \exp(-b \cdot x)$ .

**Usage**

```
FIM_exp_2par(x, w, param)
```

**Arguments**

x	Vector of design points.
w	Vector of design weight. Its length must be equal to the length of x and $\text{sum}(w) = 1$ .
param	Vector of values for the model parameters $c(a, b)$ .

**Details**

The FIM does not depend on the value of a.

**Value**

Fisher information matrix.

**References**

Dette, H., & Neugebauer, H. M. (1997). Bayesian D-optimal designs for exponential regression models. *Journal of Statistical Planning and Inference*, 60(2), 331-349.

**Examples**

```
FIM_exp_2par(x = c(1, 2), w = c(.5, .5), param = c(3, 4))
```

---

FIM\_kinetics\_alcohol *Fisher Information Matrix for the Alcohol-Kinetics Model*

---

### Description

It provides the cpp function for FIM for the model  $\sim(b_3 * x_1)/(1 + b_1 * x_1 + b_2 * x_2)$

### Usage

FIM\_kinetics\_alcohol(x1, x2, w, param)

### Arguments

x1	Vector of design points (first dimension).
x2	Vector of design points (second dimension).
w	Vector of design weight. Its length must be equal to the length of x and $\text{sum}(w) = 1$ .
param	Vector of values for the model parameters $c(b_1, b_2, b_3)$ .

### Value

Fisher information matrix.

---

FIM\_logistic *Fisher Information Matrix for the 2-Parameter Logistic (2PL) Model*

---

### Description

It provides the cpp function for FIM for the model  $\sim 1/(1 + \exp(-b * (x - a)))$ . In item response theory (IRT),  $a$  is the item difficulty parameter,  $b$  is the item discrimination parameter and  $x$  is the person ability parameter.

### Usage

FIM\_logistic(x, w, param)

### Arguments

x	Vector of design points.
w	Vector of design weight. Its length must be equal to the length of x and $\text{sum}(w) = 1$ .
param	Vector of values for the model parameters $c(a, b)$ .

**Details**

It can be shown that minimax and standardized D-optimal designs for the 2PL model is symmetric around point  $a_M = (a^L + a^U)/2$  where  $a^L$  and  $a^U$  are the lower bound and upper bound for parameter  $a$ , respectively. In `ICA.control`, arguments `sym` and `sym_point` can be used to specify  $a_M$  and find accurate symmetric optimal designs.

**Value**

Fisher information matrix.

**Examples**

```
FIM_logistic(x = c(1, 2), w = c(.5, .5), param = c(2, 1))
```

---

FIM_logistic_2pred	<i>Fisher Information Matrix for the Logistic Model with Two Predictors</i>
--------------------	---

---

**Description**

It provides the `cpp` function for FIM for the following model:  
 $\sim \exp(b_0 + b_1 * x_1 + b_2 * x_2 + b_3 * x_1 * x_2) / (1 + \exp(b_0 + b_1 * x_1 + b_2 * x_2 + b_3 * x_1 * x_2))$ .

**Usage**

```
FIM_logistic_2pred(x1, x2, w, param)
```

**Arguments**

<code>x1</code>	Vector of design points (for first predictor).
<code>x2</code>	Vector of design points (for second predictor).
<code>w</code>	Vector of design weight. Its length must be equal to the length of <code>x</code> and $\text{sum}(w) = 1$ .
<code>param</code>	Vector of values for the model parameters $c(b_0, b_1, b_2, b_3)$ .

**Value**

Fisher information matrix.



---

FIM\_loglin

*Fisher Information Matrix for the Mixed Inhibition Model*


---

**Description**

It provides the cpp function for the FIM for the model  $\sim \text{theta0} + \text{theta1} * \log(x + \text{theta2})$ .

**Usage**

FIM\_loglin(x, w, param)

**Arguments**

x                      Vector of design points.  
w                        Vector of design weight. Its length must be equal to the length of x and  $\text{sum}(w) = 1$ .  
param                    Vector of values for the model parameters  $c(\text{theta0}, \text{theta1}, \text{theta2})$ .

**Details**

The FIM of this model does not depend on the parameter  $\text{theta0}$ .

**Value**

Fisher information matrix.

**References**

Dette, H., Kiss, C., Bevanda, M., & Bretz, F. (2010). Optimal designs for the EMAX, log-linear and exponential models. *Biometrika*, 97(2), 513-518.

---

FIM\_mixed\_inhibition

*Fisher Information Matrix for the Mixed Inhibition Model.*


---

**Description**

It provides the cpp function for FIM for the model  $\sim V * S / (K_m * (1 + I / K_{ic}) + S * (1 + I / K_{iu}))$

**Usage**

FIM\_mixed\_inhibition(S, I, w, param)

**Arguments**

S	Vector of S component of design points. S is the substrate concentration.
I	Vector of I component of design points. I is the inhibitor concentration.
w	Vector of design weight. Its length must be equal to the length of S and I, besides $\text{sum}(w) = 1$ .
param	Vector of values for the model parameters $c(V, K_m, K_{ic}, K_{iu})$ .

**Details**

The optimal design does not depend on parameter  $V$ .

**Value**

Fisher information matrix of design.

**References**

Bogacka, B., Patan, M., Johnson, P. J., Youdim, K., & Atkinson, A. C. (2011). Optimum design of experiments for enzyme inhibition kinetic models. *Journal of biopharmaceutical statistics*, 21(3), 555-572.

**Examples**

```
FIM_mixed_inhibition(S = c(30, 3.86, 30, 4.60),
                    I = c(0, 0, 5.11, 4.16), w = rep(.25, 4),
                    param = c(1.5, 5.2, 3.4, 5.6))
```

---

FIM\_power\_logistic      *Fisher Information Matrix for the Power Logistic Model*

---

**Description**

It provides the cpp function for FIM for the model  $\sim 1/(1 + \exp(-b * (x - a)))^s$ , but when s is fixed (a two by two matrix).

**Usage**

```
FIM_power_logistic(x, w, param, s)
```

**Arguments**

x	Vector of design points.
w	Vector of design weight. Its length must be equal to the length of x and $\text{sum}(w) = 1$ .
param	Vector of values for the model parameters $c(a, b)$ .
s	parameter s.

**Value**

Fisher information matrix.

**Note**

This matrix is a two by two matrix and not equal to the Fisher information matrix for the power logistic model when the derivative is taken with respect to all the three parameters. This matrix is only given to be used in some illustrative examples.

---

 FIM\_sig\_emax

*Fisher Information Matrix for the Sigmoid Emax Model*


---

**Description**

It provides the cpp function for FIM for the model  $\sim b_1 + (b_2 - b_1) * (x^{b_4}) / (x^{b_4} + b_3^{b_4})$

**Usage**

```
FIM_sig_emax(x, w, param)
```

**Arguments**

x	Vector of design points.
w	Vector of design weight. Its length must be equal to the length of x and $\text{sum}(w) = 1$ .
param	Vector of values for the model parameters $c(b_1, b_2, b_3, b_4)$ . The mean of response variable is .

**Value**

Fisher information matrix.

---

 ICA.control

*ICA Control Optimization Parameters*


---

**Description**

The function ICA.control returns a list of ICA control parameters.

**Usage**

```
ICA.control(ncount = 40, nimp = ncount/10, assim_coeff = 4,
  revol_rate = 0.3, damp = 0.99, uniting_threshold = 0.02,
  equal_weight = FALSE, sym = FALSE, sym_point = NULL,
  stop_rule = c("maxiter", "equivalence"), stoptol = 0.99,
  checkfreq = 0, plot_cost = TRUE, plot_sens = TRUE,
  plot_3d = c("lattice", "rgl"), trace = TRUE, rseed = NULL)
```

**Arguments**

ncount	Number of countries. Defaults to 40.
nimp	Number of imperialists. Defaults to 10 percent of ncount.
assim_coeff	Assimilation coefficient. Defaults to 4.
revol_rate	Revolution rate. Defaults to 0.3.
damp	Damp ratio for revolution rate. <code>revol_rate</code> is decreased in every iteration by a factor of <code>damp</code> ( <code>revol_rate * damp</code> ). Defaults to 0.99.
uniting_threshold	If the distance between two imperialists is less than the product of the uniting threshold by the largest distance in the search space, ICA unites the empires. Defaults to 0.02.
equal_weight	Should the weights of design points assumed to be equal? Defaults to FALSE. If TRUE, it reduces the dimension of the search space and produces a design that gives equal weight to all of its support points.
sym	Should the design points be symmetric around <code>sym_point</code> ? Defaults to FALSE. When TRUE, <code>sym_point</code> must be given.
sym_point	If <code>sym = TRUE</code> , the design points will be symmetric around <code>sym_point</code> . See 'Details'.
stop_rule	Either 'maxiter' or 'equivalence'. Denotes the type of stopping rule. See 'Details'. Defaults to 'maxiter'.
stoptol	If <code>stop_rule = 'equivalence'</code> , algorithm stops when ELB is larger than <code>stoptol</code> . Defaults to 0.99.
checkfreq	The algorithm verifies the general equivalence theorem in every <code>checkfreq</code> iterations. When <code>checkfreq = 0</code> , no verification will be done. When <code>checkfreq = Inf</code> , only the output design will be verified. Defaults to 0.
plot_cost	Plot the iterations (evolution) of algorithm? Defaults to TRUE.
plot_sens	Plot the sensitivity (derivative) function at every <code>checkfreq</code> . Defaults to TRUE.
plot_3d	Character. Which package should be used to plot the sensitivity plot for models with two explanatory variables?
trace	Print the information in every iteration? Defaults to TRUE.
rseed	Random seed. Defaults to NULL.

**Details**

If `stop_rule = 'maxiter'`, the algorithm iterates until maximum number of iterations.

If `stop_rule = 'equivalence'`, the algorithm stops when either ELB is greater than `stoptol` or it reaches `maxiter`. In this case, you must specify the check frequency by `checkfreq`. Note that checking equivalence theorem is a very time consuming process, especially for Bayesian and minimax design problems. We advise using this option only for locally, multiple objective and robust optimal designs.

What to follows shows how `sym_point` and `sym` may be useful?

Assume the 2PL model of the form  $P(Y = 1) = \frac{1}{1 + \exp(-b(x-a))}$  and let the parameters  $a$  and  $b$  belong to  $[a_L, a_U]$  and  $[b_L, b_U]$ , respectively. It can be shown that the optimal design for this model

is symmetric around  $a_M = \frac{a_L + a_U}{2}$ . For this model, to find accurate symmetric designs, one can set `sym = TRUE` and provide the value of the  $a_M$  via `sym_point`. In this case, the output design will be symmetric around the point `sym_point`. The length of `sym_point` must be equal to the number of model predictors, here, is equal to 1.

### Value

A list of ICA control parameters.

### Examples

```
ICA.control(ncount = 100)
```

---

ICAOD

*ICAOD: Finding Optimal Designs for Nonlinear Models*

---

### Description

Different functions are available to find optimal designs for nonlinear models. The user should choose either of them based on her/his strategy to deal with the unknown model parameters: :

- **locally**: finds locally D-optimal designs. A vector of initial estimates or guess is available for the vector of model parameters.
- **robust**: finds robust optimal designs or optimal designs in average. Some (weighted) vector of initial estimates are available for the vector of unknown model parameters (discrete prior).
- **bayes**: finds Bayesian D-optimal designs. A continuous prior is available for the vector of unknown model parameters.
- **minimax**: finds minimax and standardized maximin D-optimal designs. Each of the unknown model parameters belong to an user-specified interval. The purpose is to find a design that protects the user against the worst scenario over the parameter space. Standardized designs should be used when locally optimal design of the model of interest has analytical solution.

There are also some functions to find optimal designs for special problems or models:

- **multiple**: finds locally multiple objective optimal designs for the 4-parameter Hill model. It uses the same strategy as the function `locally` to deal with the unknown model parameters.
- **bayescomp**: finds a design that that meets the dual goal of the parameter estimation and increasing the probability of a particular outcome in a binary response model. It uses the same strategy as the function `bayes` to deal with the unknown mode parameters.

### Details

The optimization of the outer problem (over the design space) is done using a metaheuristic algorithm called imperialist competitive algorithm (ICA). The user can always adjust its tuning parameters through the function `ICA.control`. Based on our experience, the most important parameter here is the number of countries, equivalent to the number of particles in the PSO algorithm and that can be regulated via the argument `ncount`.

Depending on the strategy dealing with the unknown model parameters, each type of optimal design problem may have an inner problem. Given a known design (support points and weights), the inner problems is:

- **locally**: equivalent to a simple function (criterion) evaluation at the vector of initial estimates of the vector of model parameters.
- **robust**: equivalent to a weighted some of function (criterion) evaluations at the vectors of initial estimates of the model parameters.
- **bayes**: equivalent to an integral approximation over the multiple prior distribution of the unknown model parameters (by function `hcubature`). The integration tuning parameters can be adjusted by the function `crt.bayes.control`. The most important ones are `maxEval` and `tol`.
- **bayes**: equivalent to an optimization problem over the parameter space that is the cartesian product of some intervals for the model unknown parameters (by function `nloptr`). The optimization tuning parameters can be regulated by the function `crt.minimax.control`. The most important tuning parameter is `maxeval`. The convergence of the algorithm for minimax and standardized maximin designs requires solving global multi-optima optimization problems over the parameter space. Therefore, the parameter `maxeval` plays a very important role to avoid getting trapped in the local optima of the inner problem.

The functions `locally` and `robust` are very easy to be applied and they are usually fast. The speed of the functions `bayes` and `minimax` considerably depends on the value of the tuning parameters.

Because the output design in this package is always an **approximate** or **continuous** design, the general equivalence theorem may be used to assess the proximity of the design to the true-optimal design based on derivative plot and efficiency lower bound.

For each type of design, the user can use the following functions to verify the optimality of her/his output design:

- `senslocally`
- `sensrobust`
- `sensbayes`
- `sensminimax`
- `sensmultiple`
- `sensbayescomp`

Functions for computing different relative efficiencies are also available.

## References

Masoudi E, Holling H, Wong W.K. (2017). Application of Imperialist Competitive Algorithm to Find Minimax and Standardized Maximin Optimal Designs. *Computational Statistics and Data Analysis*, 113, 330-345. <doi:10.1016/j.csda.2016.06.014>

---

iterate	<i>Updating an Object of Class 'bayes' or 'minimax'</i>
---------	---

---

**Description**

Runs ICA for more number of iterations.

**Usage**

```
iterate(object, iter)
```

**Arguments**

object	An object of class 'minimax' or 'bayes'.
iter	Number of iterations.

**Value**

An (updated) object of class 'bayes' or 'minimax'.

**See Also**

[iterate.minimax](#) and [iterate.bayes](#).

---

iterate.bayes	<i>Updating an Object of Class bayes</i>
---------------	--

---

**Description**

Runs the ICA optimization algorithm on an object of class bayes for more number of iterations and updates the results.

**Usage**

```
## S3 method for class 'bayes'  
iterate(object, iter)
```

**Arguments**

object	An object of class bayes.
iter	Number of iterations.

**See Also**

[bayes](#)

---

iterate.minimax	<i>Updating an Object of Class minimax</i>
-----------------	--

---

**Description**

Runs the ICA optimization algorithm on an object of class `minimax` for more number of iterations and updates the results.

**Usage**

```
## S3 method for class 'minimax'
iterate(object, iter)
```

**Arguments**

object	An object of class <code>minimax</code> .
iter	Number of iterations.

**See Also**

[minimax](#)

---

leff	<i>Calculates Relative Efficiency for Locally Optimal Designs</i>
------	---

---

**Description**

Given a vector of initial estimates for the parameters, this function calculates the D-and PA- efficiency of a design  $\xi_1$  with respect to a design  $\xi_2$ . Usually,  $\xi_2$  is an optimal design.

**Usage**

```
leff(formula, predvars, parvars, family = gaussian(), inipars,
     type = c("D", "PA"), fimfunc = NULL, xopt, wopt, x, w,
     npar = length(inipars), prob = NULL)
```

**Arguments**

formula	A nonlinear model <a href="#">formula</a> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <a href="#">formula</a> if necessary.
predvars	A vector of characters. Denotes the predictors in the <a href="#">formula</a> .
parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .

family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <code>gaussian()</code> is applied.
inipars	Vector. Initial values for the unknown parameters. It will be passed to the information matrix and also probability function.
type	A character. "D" denotes the D-efficiency and "PA" denotes the average P-efficiency.
fimfunc	A function. Returns the FIM as a matrix. Required when formula is missing. See 'Details' of <a href="#">minimax</a> .
xopt	Vector of design (support) points of the optimal design ( $\xi_2$ ). Similar to x.
wopt	Vector of corresponding design weights for xopt.
x	Vector of design (support) points of $\xi_1$ . See 'Details' of <a href="#">leff</a> .
w	Vector of corresponding design weights for x.
npar	Number of model parameters. Used when fimfunc is given instead of formula to specify the number of model parameters. If not given, the sensitivity plot may be shifted below the y-axis. When NULL, it will be set here to <code>length(inipars)</code> .
prob	Either formula or a function. When function, its argument are x and param, and they are the same as the arguments in fimfunc. prob as a function takes the design points and vector of parameters and returns the probability of success at each design points. See 'Examples'.

### Details

For a known  $\theta_0$ , relative D-efficiency is

$$\exp\left(\frac{\log|M(\xi, \theta_0)| - \log|M(\xi^*, \theta_0)|}{npar}\right)$$

The relative P-efficiency is

$$\exp\left(\log\left(\sum_{i=1}^k w_i p(x_i, \theta_0)\right) - \log\left(\sum_{i=1}^k w_i^* p(x_i^*, \theta_0)\right)\right)$$

where  $x^*$  and  $w^*$  are the support points and the corresponding weights of the optimal design, respectively.

The argument x is the vector of design points. For design points with more than one dimension (the models with more than one predictors), it is a concatenation of the design points, but **dimension-wise**. For example, let the model has three predictors ( $I, S, Z$ ). Then, a two-point optimal design has the following points:  $\{\text{point1} = (I_1, S_1, Z_1), \text{point2} = (I_2, S_2, Z_2)\}$ . Then, the argument x is equal to  $x = c(I1, I2, S1, S2, Z1, Z2)$ .

### Value

A value between 0 and 1.

## References

McGree, J. M., Eccleston, J. A., and Duffull, S. B. (2008). Compound optimal design criteria for nonlinear models. *Journal of Biopharmaceutical Statistics*, 18(4), 646-661.

## Examples

```
p <- c(1, -2, 1, -1)
prior4.4 <- uniform(p -1.5, p + 1.5)
formula4.4 <- ~exp(b0+b1*x1+b2*x2+b3*x1*x2)/(1+exp(b0+b1*x1+b2*x2+b3*x1*x2))
prob4.4 <- ~1-1/(1+exp(b0 + b1 * x1 + b2 * x2 + b3 * x1 * x2))
predvars4.4 <- c("x1", "x2")
parvars4.4 <- c("b0", "b1", "b2", "b3")

# Locally D-optimal design is as follows:
## weight and point of D-optimal design
# Point1    Point2    Point3    Point4
# /1.00000 \ /-1.00000 \ /0.06801 \ /1.00000 \
# \-1.00000/ \-1.00000/ \1.00000 / \1.00000 /
# Weight1    Weight2    Weight3    Weight4
# 0.250      0.250      0.250      0.250

xopt_D <- c(1, -1, .0680, 1, -1, -1, 1, 1)
wopt_D <- rep(.25, 4)

# Let see if we use only three of the design points, what is the relative efficiency.
leff(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4, family = binomial(),
      x = c(1, -1, .0680, -1, -1, 1), w = c(.33, .33, .33),
      inipars = p,
      xopt = xopt_D, wopt = wopt_D)
# Wow, it heavily drops!

# Locally P-optimal design has only one support point and is -1 and 1
xopt_P <- c(-1, 1)
wopt_P <- 1

# What is the relative P-efficiency of the D-optimal design with respect to P-optimal design?
leff(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4, family = binomial(),
      x = xopt_D, w = wopt_D,
      inipars = p,
      type = "PA",
      prob = prob4.4,
      xopt = xopt_P, wopt = wopt_P)
# .535
```

## Description

Finds locally D-optimal designs for nonlinear models. It should be used when a vector of initial estimates is available for the unknown model parameters. Locally optimal designs may not be efficient when the initial estimates are far away from the true values of the parameters.

## Usage

```
locally(formula, predvars, parvars, family = gaussian(), lx, ux, iter, k,
        inipars, fimfunc = NULL, ICA.control = list(),
        sens.minimax.control = list(), initial = NULL,
        npar = length(inipars), plot_3d = c("lattice", "rgl"), x = NULL)
```

## Arguments

formula	A nonlinear model <a href="#">formula</a> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <a href="#">formula</a> if necessary.
predvars	A vector of characters. Denotes the predictors in the <a href="#">formula</a> .
parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .
family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <a href="#">gaussian()</a> is applied.
lx	Vector of lower bounds for the predictors. Should be in the same order as <a href="#">predvars</a> .
ux	Vector of upper bounds for the predictors. Should be in the same order as <a href="#">predvars</a> .
iter	Maximum number of iterations.
k	Number of design points. Must be at least equal to the number of model parameters to avoid singularity of the FIM.
inipars	A vector of initial estimates for the unknown parameters. It must match <a href="#">parvars</a> or the argument <a href="#">param</a> of the function <a href="#">fimfunc</a> , when provided.
fimfunc	A function. Returns the FIM as a matrix. Required when <a href="#">formula</a> is missing. See 'Details' of <a href="#">minimax</a> .
ICA.control	ICA control parameters. For details, see <a href="#">ICA.control</a> .
sens.minimax.control	Control parameters to verify the general equivalence theorem. For details, see the function <a href="#">sens.minimax.control</a> .
initial	A matrix of the initial design points and weights that will be inserted into the initial solutions (countries) of the algorithm. Every row is a design, i.e. a concatenation of $x$ and $w$ . Will be coerced to a matrix if necessary. See 'Details' of <a href="#">minimax</a> .

npar	Number of model parameters. Used when <code>fimfunc</code> is given instead of <code>formula</code> to specify the number of model parameters. If not given, the sensitivity plot may be shifted below the y-axis. When <code>NULL</code> , it is set to <code>length(inipars)</code> .
plot_3d	Which package should be used to plot the sensitivity (derivative) function for two-dimensional design space. Defaults to "lattice".
x	A vector of candidate design (support) points. When is not set to <code>NULL</code> (default), the algorithm only finds the optimal weights for the candidate points in <code>x</code> . Should be set when the user has a finite number of candidate design points and the purpose is to find the optimal weight for each of them (when zero, they will be excluded from the design). For design points with more than one dimension, see 'Details' of <a href="#">sensminimax</a> .

### Details

Let  $M(\xi, \theta_0)$  be the Fisher information matrix (FIM) of a  $k$ -point design  $\xi$  and  $\theta_0$  be the vector of the initial estimates for the unknown parameters. A locally D-optimal design  $\xi^*$  minimizes over  $\Xi$

$$-\log |M(\xi, \theta_0)|.$$

One can adjust the tuning parameters in [ICA.control](#) to set a stopping rule based on the general equivalence theorem. See "Examples" below.

### Value

an object of class `minimax` that is a list including three sub-lists:

`arg` A list of design and algorithm parameters.

`evol` A list of length equal to the number of iterations that stores the information about the best design (design with least criterion value) of each iteration. `evol[[iter]]` contains:

<code>iter</code>	Iteration number.
<code>x</code>	Design points.
<code>w</code>	Design weights.
<code>min_cost</code>	Value of the criterion for the best imperialist (design).
<code>mean_cost</code>	Mean of the criterion values of all the imperialists.
<code>sens</code>	An object of class 'sensminimax'. See below.
<code>param</code>	Vector of parameters.

`empires` A list of all the empires of the last iteration.

`alg` A list with following information:

<code>nfeval</code>	Number of function evaluations. It does not count the function evaluations from checking the general equivalence theorem.
<code>nlocal</code>	Number of successful local searches.
<code>nrevol</code>	Number of successful revolutions.
<code>nimprove</code>	Number of successful movements toward the imperialists in the assimilation step.
<code>convergence</code>	Stopped by 'maxiter' or 'equivalence'?

The list `sens` contains information about the design verification by the general equivalence theorem. See `sensminimax` for more details. It is given every `ICA.control$checkfreq` iterations and also the last iteration if `ICA.control$checkfreq >= 0`. Otherwise, `NULL`.

`param` is a vector of parameters that is the global minimum of the minimax criterion or the global maximum of the standardized maximin criterion over the parameter space, given the current `x`, `w`.

## References

Masoudi E, Holling H, Wong W.K. (2017). Application of Imperialist Competitive Algorithm to Find Minimax and Standardized Maximin Optimal Designs. *Computational Statistics and Data Analysis*, 113, 330-345.

## See Also

[senslocally](#)

## Examples

```
#####
# Exponential growth model
#####
# See how we set stopping rule by adjusting 'stop_rule', 'checkfreq' and 'stoptol'
# It calls the 'senslocally' function every checkfreq = 50 iterations to
# calculate the ELB. if ELB is greater than stoptol = .95, then the algorithm stops.

# initializing by one iteration
res1 <- locally(formula = ~a + exp(-b*x), predvars = "x", parvars = c("a", "b"),
               lx = 0, ux = 1, inipars = c(1, 10),
               iter = 1, k = 2,
               ICA.control= ICA.control(rseed = 100,
                                       stop_rule = "equivalence",
                                       checkfreq = 20, stoptol = .95))

## Not run:
# update the algorithm
res1 <- iterate(res1, 150)
#stops at iteration 21 because ELB is greater than .95

## End(Not run)

### fixed x, lx and ux are only required for equivalence theorem
## Not run:
res1.1 <- locally(formula = ~a + exp(-b*x), predvars = "x", parvars = c("a", "b"),
                 lx = 0, ux = 1, inipars = c(1, 10),
                 iter = 100, k = 3,
                 x = c(.25, .5, .75),
                 ICA.control= ICA.control(rseed = 100))

plot(res1.1)
# we can not have an optimal design using this x

## End(Not run)
```

```
#####
## two parameter logistic model
#####
res2 <- locally(formula = ~1/(1 + exp(-b *(x - a))),
                predvars = "x", parvars = c("a", "b"),
                family = binomial(), lx = -3, ux = 3,
                inipars = c(1, 3), iter = 1, k = 2,
                ICA.control= list(rseed = 100, stop_rule = "equivalence",
                                checkfreq = 50, stoptol = .95))

## Not run:
  res2 <- iterate(res2, 100)
  # stops at iteration 51

## End(Not run)

#####
# A model with two predictors
#####
# mixed inhibition model
## Not run:
  res3 <- locally(formula = ~ V*S/(Km * (1 + I/Kic)+ S * (1 + I/Kiu)),
                predvars = c("S", "I"),
                parvars = c("V", "Km", "Kic", "Kiu"),
                family = gaussian(),
                lx = c(0, 0), ux = c(30, 60),
                k = 4,
                iter = 300,
                inipars = c(1.5, 5.2, 3.4, 5.6),
                ICA.control= list(rseed = 100, stop_rule = "equivalence",
                                checkfreq = 50, stoptol = .95))

  # stops at iteration 100

## End(Not run)

## Not run:
  # fixed x
  res3.1 <- locally(formula = ~ V*S/(Km * (1 + I/Kic)+ S * (1 + I/Kiu)),
                predvars = c("S", "I"),
                parvars = c("V", "Km", "Kic", "Kiu"),
                family = gaussian(),
                lx = c(0, 0), ux = c(30, 60),
                k = 5,
                iter = 100,
                x = c(20, 4, 20, 4, 10, 0, 0, 30, 3, 2),
                inipars = c(1.5, 5.2, 3.4, 5.6),
                ICA.control= list(rseed = 100))

## End(Not run)
```

**Description**

Finds compound locally DP-optimal designs that meet the dual goal of parameter estimation and increasing the probability of a particular outcome in a binary response model. A compound locally DP-optimal design maximizes the product of the efficiencies of a design  $\xi$  with respect to D- and average P-optimality, weighted by a pre-defined mixing constant  $0 \leq \alpha \leq 1$ .

**Usage**

```
locallycomp(formula, predvars, parvars, family = gaussian(), lx, ux,
  alpha, prob, iter, k, inipars, fimfunc = NULL, ICA.control = list(),
  sens.minimax.control = list(), initial = NULL,
  npar = length(inipars), plot_3d = c("lattice", "rgl"))
```

**Arguments**

formula	A nonlinear model <a href="#">formula</a> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <a href="#">formula</a> if necessary.
predvars	A vector of characters. Denotes the predictors in the <a href="#">formula</a> .
parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .
family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <code>gaussian()</code> is applied.
lx	Vector of lower bounds for the predictors. Should be in the same order as <code>predvars</code> .
ux	Vector of upper bounds for the predictors. Should be in the same order as <code>predvars</code> .
alpha	A value between 0 and 1. Compound or combined DP-criterion is the product of the efficiencies of a design with respect to D- and average P- optimality, weighted by alpha.
prob	Either <a href="#">formula</a> or a function. When function, its argument are <code>x</code> and <code>param</code> , and they are the same as the arguments in <code>fimfunc</code> . <code>prob</code> as a function takes the design points and vector of parameters and returns the probability of success at each design points. See 'Examples'.
iter	Maximum number of iterations.
k	Number of design points. When <code>alpha = 0</code> , then <code>k</code> can be less than the number of parameters.

<code>inipars</code>	Vector. Initial values for the unknown parameters. It will be passed to the information matrix and also probability function.
<code>fimfunc</code>	A function. Returns the FIM as a matrix. Required when formula is missing. See 'Details' of <code>minimax</code> .
<code>ICA.control</code>	ICA control parameters. For details, see <code>ICA.control</code> .
<code>sens.minimax.control</code>	Control parameters to verify the general equivalence theorem. For details, see the function <code>sens.minimax.control</code> .
<code>initial</code>	A matrix of the initial design points and weights that will be inserted into the initial solutions (countries) of the algorithm. Every row is a design, i.e. a concatenation of <code>x</code> and <code>w</code> . Will be coerced to a matrix if necessary. See 'Details' of <code>minimax</code> .
<code>npar</code>	Number of model parameters. Used when <code>fimfunc</code> is given instead of formula to specify the number of model parameters. If not given, the sensitivity plot may be shifted below the y-axis. When NULL, it will be set here to <code>length(inipars)</code> .
<code>plot_3d</code>	Which package should be used to plot the sensitivity (derivative) function for two-dimensional design space. Defaults to "lattice".

## Details

Let  $\Xi$  be the space of all approximate designs with  $k$  design points (support points) at  $x_1, x_2, \dots, x_k$  from design space  $\chi$  with corresponding weights  $w_1, \dots, w_k$ . Let  $M(\xi, \theta)$  be the Fisher information matrix (FIM) of a  $k$ -point design  $\xi$ ,  $\theta_0$  is a user-given vector of initial estimates for the unknown parameters  $\theta$  and  $p(x_i, \theta)$  is the  $i$ th probability of success given by  $x_i$  in a binary response model. A compound locally DP-optimal design maximizes over  $\Xi$

$$\frac{\alpha}{q} \log |M(\xi, \theta_0)| + (1 - \alpha) \log \left( \sum_{i=1}^k w_i p(x_i, \theta_0) \right).$$

Use `plot` function to verify the general equivalence theorem for the output design or change `checkfreq` in `ICA.control`.

One can adjust the tuning parameters in `ICA.control` to set a stopping rule based on the general equivalence theorem. See "Examples" in `locally`.

## Value

an object of class `minimax` that is a list including three sub-lists:

`arg` A list of design and algorithm parameters.

`evol` A list of length equal to the number of iterations that stores the information about the best design (design with least criterion value) of each iteration. `evol[[iter]]` contains:

<code>iter</code>	Iteration number.
<code>x</code>	Design points.
<code>w</code>	Design weights.
<code>min_cost</code>	Value of the criterion for the best imperialist (design).
<code>mean_cost</code>	Mean of the criterion values of all the imperialists.

sens An object of class 'sensminimax'. See below.  
 param Vector of parameters.

empires A list of all the empires of the last iteration.

alg A list with following information:

nfeval Number of function evaluations. It does not count the function evaluations from checking the general equivalence theorem.  
 nlocal Number of successful local searches.  
 nrevol Number of successful revolutions.  
 nimprove Number of successful movements toward the imperialists in the assimilation step.  
 convergence Stopped by 'maxiter' or 'equivalence'?

The list sens contains information about the design verification by the general equivalence theorem. See sensminimax for more details. It is given every ICA.control\$checkfreq iterations and also the last iteration if ICA.control\$checkfreq >= 0. Otherwise, NULL.

param is a vector of parameters that is the global minimum of the minimax criterion or the global maximum of the standardized maximin criterion over the parameter space, given the current x, w.

## References

McGree, J. M., Eccleston, J. A., and Duffull, S. B. (2008). Compound optimal design criteria for nonlinear models. *Journal of Biopharmaceutical Statistics*, 18(4), 646-661.

## Examples

```
## Here we produce the results of Table 2 in in McGree and Eccleston (2008)
# For D- and P-efficiency see, ?leff and ?peff

p <- c(1, -2, 1, -1)
prior4.4 <- uniform(p -1.5, p + 1.5)
formula4.4 <- ~exp(b0+b1*x1+b2*x2+b3*x1*x2)/(1+exp(b0+b1*x1+b2*x2+b3*x1*x2))
prob4.4 <- ~1-1/(1+exp(b0 + b1 * x1 + b2 * x2 + b3 * x1 * x2))
predvars4.4 <- c("x1", "x2")
parvars4.4 <- c("b0", "b1", "b2", "b3")
lb <- c(-1, -1)
ub <- c(1, 1)

# set checkfreq = Inf to ask for equivalence theorem at final step.
res.0 <- locallycomp(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4,
  family = binomial(), prob = prob4.4, lx = lb, ux = ub,
  alpha = 0, k = 1, inipars = p, iter = 10,
  ICA.control = ICA.control(checkfreq = Inf))

## Not run:
res.25 <- locallycomp(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4,
  family = binomial(), prob = prob4.4, lx = lb, ux = ub,
  alpha = .25, k = 4, inipars = p, iter = 350,
```

```

ICA.control = ICA.control(checkfreq = Inf))

res.5 <- locallycomp(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4,
  family = binomial(), prob = prob4.4, lx = lb, ux = ub,
  alpha = .5, k = 4, inipars = p, iter = 350,
  ICA.control = ICA.control(checkfreq = Inf))
res.75 <- locallycomp(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4,
  family = binomial(), prob = prob4.4, lx = lb, ux = ub,
  alpha = .75, k = 4, inipars = p, iter = 350,
  ICA.control = ICA.control(checkfreq = Inf))

res.1 <- locallycomp(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4,
  family = binomial(), prob = prob4.4, lx = lb, ux = ub,
  alpha = 1, k = 4, inipars = p, iter = 350,
  ICA.control = ICA.control(checkfreq = Inf))

#### computing the D-efficiency
# locally D-optimal design is locally DP-optimal design when alpha = 1.

leff(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4, family = binomial(),
  x = res.0$evol[[10]]$x, w = res.0$evol[[10]]$w,
  inipars = p,
  xopt = res.1$evol[[350]]$x, wopt = res.1$evol[[350]]$w)

leff(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4, family = binomial(),
  x = res.25$evol[[350]]$x, w = res.25$evol[[350]]$w,
  inipars = p,
  xopt = res.1$evol[[350]]$x, wopt = res.1$evol[[350]]$w)

leff(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4, family = binomial(),
  x = res.5$evol[[350]]$x, w = res.5$evol[[350]]$w,
  inipars = p,
  xopt = res.1$evol[[350]]$x, wopt = res.1$evol[[350]]$w)

leff(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4, family = binomial(),
  x = res.75$evol[[350]]$x, w = res.75$evol[[350]]$w,
  inipars = p,
  xopt = res.1$evol[[350]]$x, wopt = res.1$evol[[350]]$w)

#### computing the P-efficiency
# locally p-optimal design is locally DP-optimal design when alpha = 0.

leff(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4, family = binomial(),
  xopt = res.0$evol[[10]]$x, wopt = res.0$evol[[10]]$w,
  prob = prob4.4,
  type = "PA",
  inipars = p,
  x = res.25$evol[[350]]$x, w = res.25$evol[[350]]$w)

leff(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4, family = binomial(),

```

```

xopt = res.0$evol[[10]]$x, wopt = res.0$evol[[10]]$w,
prob = prob4.4,
inipars = p,
type = "PA",
x = res.5$evol[[350]]$x, w = res.5$evol[[350]]$w)

leff(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4, family = binomial(),
xopt = res.0$evol[[10]]$x, wopt = res.0$evol[[10]]$w,
prob = prob4.4,
inipars = p,
type = "PA",
x = res.75$evol[[350]]$x, w = res.75$evol[[350]]$w)

leff(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4, family = binomial(),
xopt = res.0$evol[[10]]$x, wopt = res.1$evol[[10]]$w,
prob = prob4.4,
type = "PA",
inipars = p,
x = res.1$evol[[350]]$x, w = res.1$evol[[350]]$w)

## End(Not run)

```

---

minimax

---

*Minimax and Standardized Maximin D-Optimal Designs*


---

## Description

Finds minimax and standardized maximin D-optimal designs for nonlinear models. It should be used when the user assumes the unknown parameters belong to a parameter region  $\Theta$ , which is called “region of uncertainty”, where the purpose is to protect the experiment from the worst case scenario over  $\Theta$ .

## Usage

```

minimax(formula, predvars, parvars, family = gaussian(), lx, ux, lp, up,
iter, k, n.grid = 0, fimfunc = NULL, ICA.control = list(),
sens.minimax.control = list(), crt.minimax.control = list(),
standardized = FALSE, initial = NULL, localdes = NULL,
npar = length(lp), plot_3d = c("lattice", "rgl"), x = NULL)

```

## Arguments

formula	A nonlinear model <code>formula</code> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <code>formula</code> if necessary.
predvars	A vector of characters. Denotes the predictors in the <code>formula</code> .

parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .
family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <code>gaussian()</code> is applied.
lx	Vector of lower bounds for the predictors. Should be in the same order as <code>predvars</code> .
ux	Vector of upper bounds for the predictors. Should be in the same order as <code>predvars</code> .
lp	Vector of lower bounds for the model parameters. Should be in the same order as <code>parvars</code> or <code>param</code> in the argument <code>fimfunc</code> .
up	Vector of upper bounds for the model parameters. Should be in the same order as <code>parvars</code> or <code>param</code> in the argument <code>fimfunc</code> . When a parameter is known (has a fixed value), its associated lower and upper bound values in <code>lp</code> and <code>up</code> must be set equal.
iter	Maximum number of iterations.
k	Number of design points. Must be at least equal to the number of model parameters to avoid singularity of the FIM.
n.grid	Only required when the parameter space is going to be discretized. The total number of grid points from the parameter space is <code>n.grid^p</code> . When <code>n.grid &gt; 0</code> , optimal design protects the experimenter against the worst case scenario only over the grid points, and not over the continuous parameter space. The resulting designs may not be globally optimal. In some literature, this type of designs has been used as a compromise to the minimax type designs to avoid continuous optimization problem over the parameter space and simplify the minimax design problems. Especially when the design criterion is convex with respect to the given parameter space at every given design from the design space, the obtained design may also be globally optimal (because the maximum of a convex function is attained on the bounds, and here, are included in the grid points). See 'Details' of <a href="#">minimax</a> .
fimfunc	A function. Returns the FIM as a matrix. Required when <code>formula</code> is missing. See 'Details' of <a href="#">minimax</a> .
ICA.control	ICA control parameters. For details, see <a href="#">ICA.control</a> .
sens.minimax.control	Control parameters to verify the general equivalence theorem. For details, see the function <a href="#">sens.minimax.control</a> .
crt.minimax.control	Control parameters to optimize the minimax or standardized maximin criterion at a given design over a <b>continuous</b> parameter space (when <code>n.grid = 0</code> ). For details, see the function <a href="#">crt.minimax.control</a> .
standardized	Maximin standardized design? When <code>standardized = TRUE</code> , the argument <code>localdes</code> must be given. Defaults to FALSE. See 'Details' of <a href="#">minimax</a> .

initial	A matrix of the initial design points and weights that will be inserted into the initial solutions (countries) of the algorithm. Every row is a design, i.e. a concatenation of $x$ and $w$ . Will be coerced to a matrix if necessary. See 'Details' of <a href="#">minimax</a> .
localdes	A function that takes the parameter values as inputs and returns the design points and weights of the locally optimal design. Required when <code>standardized = "TRUE"</code> . See 'Details' of <a href="#">minimax</a> .
npar	Number of model parameters. Used when <code>fimfunc</code> is given instead of <code>formula</code> to specify the number of model parameters. If not specified truly, the sensitivity (derivative) plot may be shifted below the y-axis. When NULL (default), it is set to <code>length(lp)</code> .
plot_3d	Which package should be used to plot the sensitivity (derivative) function for two-dimensional design space. Defaults to "lattice".
x	A vector of candidate design (support) points. When is not set to NULL (default), the algorithm only finds the optimal weights for the candidate points in $x$ . Should be set when the user has a finite number of candidate design points and the purpose is to find the optimal weight for each of them (when zero, they will be excluded from the design). For design points with more than one dimension, see 'Details' of <a href="#">sensminimax</a> .

## Details

Let  $\Xi$  be the space of all approximate designs with  $k$  design points (support points) at  $x_1, x_2, \dots, x_k$  from the design space  $\chi$  with corresponding weights  $w_1, \dots, w_k$ . Let  $M(\xi, \theta)$  be the Fisher information matrix (FIM) of a  $k$ -point design  $\xi$  and  $\theta$  be the vector of unknown parameters. A minimax D-optimal design  $\xi^*$  minimizes over  $\Xi$

$$\max_{\theta \in \Theta} -\log |M(\xi, \theta)|.$$

A standardized maximin D-optimal design  $\xi^*$  maximizes over  $\Xi$

$$\inf_{\theta \in \Theta} \left[ \left( \frac{|M(\xi, \theta)|}{|M(\xi_\theta, \theta)|} \right)^{\frac{1}{p}} \right],$$

where  $p$  is the number of model parameters and  $\xi_\theta$  is the locally D-optimal design with respect to  $\theta$ .

A minimax criterion (cost function or objective function) is evaluated at each design (decision variables) by maximizing the criterion over the parameter space. We call the optimization problem over the parameter space as *inner optimization problem*. Two different strategies may be applied to solve the inner problem at a given design (design points and weights):

1. **Continuous inner problem:** we optimize the criterion over a continuous parameter space using the function `nloptr`. In this case, the tuning parameters can be regulated via the argument `crt.minimax.control`, when the most influential one is `maxeval`.
2. **Discrete inner problem:** we map the parameter space to the grid points and optimize the criterion over a discrete parameter space. In this case, the number of grid points can be regulated via `n.grid`. This strategy is quite efficient (and fast) when the maxima most likely attain

the vertices of the continuous parameter space at any given design. The output design here protects the experiment from the worst scenario over the grid points.

The formula is used to automatically create the Fisher information matrix (FIM) for a nonlinear model provided that the distribution of the response variable belongs to the natural exponential family. Function `minimax` also provides an option to assign a user-defined FIM directly via the argument `fimfunc`. In this case, the argument `fimfunc` takes a function that has three arguments as follows:

1. `x` a vector of design points. For design points with more than one dimension, it is a concatenation of the design points, but dimension-wise. For example, let the model has three predictors  $(I, S, Z)$ . Then, a two-point design is of the format  $\{\text{point1} = (I_1, S_1, Z_1), \text{point2} = (I_2, S_2, Z_2)\}$ . and the argument `x` is equivalent to `x = c(I1, I2, S1, S2, Z1, Z2)`.
2. `w` a vector that includes the design weights associated with `x`.
3. `param` a vector of parameter values associated with `lp` and `up`.

The output must be the Fisher information matrix with number of rows equal to `length(param)`. See 'Examples'.

Minimax optimal designs can have very different criterion values depending on the nominal set of parameter values. Accordingly, it is desirable to standardize the criterion and control for the potentially widely varying magnitude of the criterion (Dette, 1997). Evaluating a standardized maximin criterion requires knowing locally optimal designs. We strongly advise setting `standardized = 'TRUE'` only when analytical solutions for the locally D-optimal designs is available. In this case, for any initial estimate of the unknown parameters, an analytical solution for the locally optimal design, i.e, the support points `x` and the corresponding weights `w`, must be provided via the argument `localdes`. Therefore, depending on how the model is specified, `localdes` is a function with the following arguments (input).

- If `formula` is given (`!missing(formula)`):
  - The parameter names given by `parvars` in the same order.
- If FIM is given via the argument `fimfunc` (`missing(formula)`):
  - `param`: A vector of the parameters equal to the argument `param` in `fimfunc`.

This function must return a list with the components `x` and `w` (they match the same arguments in the function `fimfunc`). See 'Examples'.

The standardized D-criterion is equal to the D-efficiency and it must be between 0 and 1. However, in practice, when running the algorithm, it may be the case that the criterion takes a value larger than one. This may happen because the user-function that is given via `localdes` does not return the true (accurate) locally optimal designs for some requested initial estimates of the parameters from  $\Theta$ . In this case, the function `minimax` throw an error where the error message helps the user to debug her/his function.

Each row of `initial` is one design, i.e. a concatenation of values for design (support) points and the associated design weights. Let `x0` and `w0` be the vector of initial values with exactly the same length and order as `x` and `w` (the arguments of `fimfunc`). As an example, the first row of the matrix `initial` is equal to `initial[1, ] = c(x0, w0)`. For models with more than one predictors, `x0` is a concatenation of the initial values for design points, but **dimension-wise**. See the details of the argument `fimfunc`, above.

To verify the optimality of the output design by the general equivalence theorem, the user can either plot the results or set `checkfreq` in `ICA.control` to `Inf`. In either way, the function `sensminimax` is called for verification. Note that the function `sensminimax` always verifies the optimality of a design assuming a continuous parameter space. See 'Examples'.

### Value

an object of class `minimax` that is a list including three sub-lists:

`arg` A list of design and algorithm parameters.

`evol` A list of length equal to the number of iterations that stores the information about the best design (design with least criterion value) of each iteration. `evol[[iter]]` contains:

<code>iter</code>	Iteration number.
<code>x</code>	Design points.
<code>w</code>	Design weights.
<code>min_cost</code>	Value of the criterion for the best imperialist (design).
<code>mean_cost</code>	Mean of the criterion values of all the imperialists.
<code>sens</code>	An object of class 'sensminimax'. See below.
<code>param</code>	Vector of parameters.

`empires` A list of all the empires of the last iteration.

`alg` A list with following information:

<code>nfeval</code>	Number of function evaluations. It does not count the function evaluations from checking the general equivalence theorem.
<code>nlocal</code>	Number of successful local searches.
<code>nrevol</code>	Number of successful revolutions.
<code>nimprove</code>	Number of successful movements toward the imperialists in the assimilation step.
<code>convergence</code>	Stopped by 'maxiter' or 'equivalence'?

The list `sens` contains information about the design verification by the general equivalence theorem. See `sensminimax` for more details. It is given every `ICA.control$checkfreq` iterations and also the last iteration if `ICA.control$checkfreq >= 0`. Otherwise, `NULL`.

`param` is a vector of parameters that is the global minimum of the minimax criterion or the global maximum of the standardized maximin criterion over the parameter space, given the current `x`, `w`.

### Note

For larger parameter space or model with more number of unknown parameters, it is always important to increase the value of `ncount` in `ICA.control` and `optslst$maxeval` in `crt.minimax.control` to produce very accurate designs.

Although standardized criteria have been preferred theoretically, in practice, they should be applied only when an analytical solution for the locally D-optimal designs is available for the model of interest. Otherwise, we encounter a three-level nested-optimization algorithm.

## References

- Masoudi E, Holling H, Wong W.K. (2017). Application of Imperialist Competitive Algorithm to Find Minimax and Standardized Maximin Optimal Designs. *Computational Statistics and Data Analysis*, 113, 330-345.
- Dette, H. (1997). Designing experiments with respect to 'standardized' optimality criteria. *Journal of the Royal Statistical Society: Series B (Statistical Methodology)*, 59(1), 97-110.

## See Also

[sensminimax](#)

## Examples

```
#####
# Two-parameter exponential growth model
#####
res1 <- minimax (formula = ~a + exp(-b*x), predvars = "x", parvars = c("a", "b"),
                lx = 0, ux = 1, lp = c(1, 1), up = c(1, 10),
                iter = 1, k = 4, ICA.control= ICA.control(rseed = 100))
# the optimal design has 3 points, but we set k = 4 for illustration purpose to
# show how the algorithm modify the design by assigning weights and locating
# the points.
## Not run:
  res1 <- iterate(res1, 150)
  # iterating the algorithm up to 150 more iterations

## End(Not run)

res1 # print method
plot(res1) # verifying the general equivalence theorem

## Not run:
  ## fixed x
res1.1 <- minimax (formula = ~a + exp(-b*x), predvars = "x", parvars = c("a", "b"),
                  lx = 0, ux = 1, lp = c(1, 1), up = c(1, 10),
                  x = c(0, .5, 1),
                  iter = 150, k = 3, ICA.control= ICA.control(rseed = 100))
# not optimal

## End(Not run)

#####
# Two-parameter logistic model.
#####
# A little bit tickling with the tuning parameters
# reducing the value of maxeval to 200 to increase the speed
cont1 <- crt.minimax.control(optslist = list(maxeval = 500))
cont2 <- ICA.control(rseed = 100, checkfreq = Inf, ncount = 60)

## Not run:
```

```

res2 <- minimax (formula = ~1/(1 + exp(-b *(x - a))), predvars = "x",
                parvars = c("a", "b"),
                family = binomial(), lx = -3, ux = 3,
                lp = c(0, 1), up = c(1, 2.5), iter = 200, k = 3,
                ICA.control= cont2, crt.minimax.control = cont1)

print(res2)
plot(res2)

## End(Not run)

#####
# An example of a model with two predictors
#####
# Mixed inhibition model
lower <- c(1, 4, 2, 4)
upper <- c(1, 5, 3, 5)
cont <- crt.minimax.control(optslist = list(maxeval = 100)) # to be faster
## Not run:
res3 <- minimax(formula = ~ V*S/(Km * (1 + I/Kic)+ S * (1 + I/Kiu)),
                predvars = c("S", "I"),
                parvars = c("V", "Km", "Kic", "Kiu"),
                lx = c(0, 0), ux = c(30, 60), k = 4,
                iter = 100, lp = lower, up = upper,
                ICA.control= list(rseed = 100),
                crt.minimax.control = cont)

res3 <- iterate(res3, 100)
print(res3)
plot(res3) # sensitivity plot
res3$arg$time

## End(Not run)

# Now consider grid points instead of assuming continuous parameter space
# set n.grid to 5
## Not run:
res4 <- minimax(formula = ~ V*S/(Km * (1 + I/Kic)+ S * (1 + I/Kiu)),
                predvars = c("S", "I"),
                parvars = c("V", "Km", "Kic", "Kiu"),
                lx = c(0, 0), ux = c(30, 60),
                k = 4, iter = 130, n.grid = 5, lp = lower, up = upper,
                ICA.control= list(rseed = 100, checkfreq = Inf),
                crt.minimax.control = cont)

print(res4)
plot(res4) # sensitivity plot

## End(Not run)

#####
# Standardized maximin D-optimal designs
#####
# Now assume the purpose is finding STANDARDIZED designs
# We know from the literature that the locally D-optimal design (LDOD)

```

```

# for this model has analytical solution.
# The following function takes the parameter as input and returns
# the design points and weights of LDOD.
# x and w are exactly similar to the arguments of 'fimfunc'.
# x is a vector and returns the design points 'dimension-wise'.
# see explanation of the arguments of 'fimfunc' in 'Details'.

LDOD <- function(V, Km, Kic, Kiu){
  #first dimension is for S and the second one is for I.
  S_min <- 0
  S_max <- 30
  I_min <- 0
  I_max <- 60
  s2 <- max(S_min, S_max*Km*Kiu*(Kic+I_min)/
            (S_max*Kic*I_min+S_max*Kic*Kiu+2*Km*Kiu*I_min+2*Km*Kiu*Kic))
  i3 <- min((2*S_max*Kic*I_min + S_max*Kic*Kiu+2*Km*Kiu*I_min+Km*Kiu*Kic)/
            (Km*Kiu+S_max*Kic), I_max)
  i4 <- min(I_min + (sqrt((Kic+I_min)*(Km*Kic*Kiu+Km*Kiu*I_min+
                               S_max*Kic*Kiu+S_max*Kic*I_min)/
                               (Km*Kiu+S_max*Kic))), I_max )
  s4 <- max(-Km*Kiu*(Kic+2*I_min-i4)/(Kic*(Kiu+2*I_min-i4)), S_min)
  x <- c(S_max, s2, S_max, s4, I_min, I_min, i3, i4)
  return(list(x = x, w =rep(1/4, 4)))
}
formalArgs(LDOD)
## Not run:
  minimax(formula = ~ V*S/(Km * (1 + I/Kic)+ S * (1 + I/Kiu)),
          predvars = c("S", "I"),
          parvars = c("V", "Km", "Kic", "Kiu"),
          lx = c(0, 0), ux = c(30, 60),
          k = 4, iter = 300,
          lp = lower, up = upper,
          ICA.control= list(rseed = 100, checkfreq = Inf),
          crt.minimax.control = cont,
          standardized = TRUE,
          localdes = LDOD)

## End(Not run)

#####
# Not necessary!
# The rest of the examples here are only for professional uses.
#####
# Imagine you have written your own FIM, say in Rcpp that is faster than
# the FIM created by the formula interface above.

#####
# An example of a model with two predictors
#####
# For example, the cpp FIM function for the mixed inhibition model is named:
formalArgs(FIM_mixed_inhibition)

```

```

# We should reparameterize the arguments to match the standard of the
# argument 'fimfunc' (see 'Details').
myfim <- function(x, w, param){
  npoint <- length(x)/2
  S <- x[1:npoint]
  I <- x[(npoint+1):(npoint*2)]
  out <- FIM_mixed_inhibition(S = S, I = I, w = w, param = param)
  return(out)
}
formalArgs(myfim)

# Finds minimax optimal design, exactly as before, but NOT using the
# formula interface.
## Not run:
res5 <- minimax(fimfunc = myfim,
                lx = c(0, 0), ux = c(30, 60), k = 4,
                iter = 100, lp = lower, up = upper,
                ICA.control= list(rseed = 100),
                crt.minimax.control = cont)

print(res5)
plot(res5) # sensitivity plot

## End(Not run)
#####
# Standardized maximin D-optimal designs
#####
# To match the argument 'localdes' when no formula interface is used,
# we should reparameterize LDOD.
# The input must be 'param' same as the argument of 'fimfunc'
LDOD2 <- function(param)
  LDOD(V = param[1], Km = param[2], Kic = param[3], Kiu = param[4])

# compare these two:
formalArgs(LDOD)
formalArgs(LDOD2)
## Not run:
res6 <- minimax(fimfunc = myfim,
                lx = c(0, 0), ux = c(30, 60), k = 4,
                iter = 300, lp = lower, up = upper,
                ICA.control= list(rseed = 100, checkfreq = Inf),
                crt.minimax.control = cont,
                standardized = TRUE,
                localdes = LDOD2)

res6
plot(res6)

## End(Not run)

```

---

multiple

*Locally Multiple Objective Optimal Designs for the 4-Parameter Hill Model***Description**

The 4-parameter Hill model is of the form

$$f(D) = c + \frac{(d-c)\left(\frac{D}{a}\right)^b}{1 + \left(\frac{D}{a}\right)^b} + \epsilon,$$

where  $\epsilon \sim N(0, \sigma^2)$ ,  $D$  is the dose level and the predictor,  $a$  is the ED50,  $d$  is the upper limit of response,  $c$  is the lower limit of response and  $b$  denotes the Hill constant that control the flexibility in the slope of the response curve.

Sometimes, the Hill model is re-parameterized and written as

$$f(x) = \frac{\theta_1}{1 + \exp(\theta_2 x + \theta_3)} + \theta_4,$$

where  $\theta_1 = d - c$ ,  $\theta_2 = -b$ ,  $\theta_3 = b \log(a)$ ,  $\theta_4 = c$ ,  $\theta_1 > 0$ ,  $\theta_2 \neq 0$ , and  $-\infty < ED50 < \infty$ , where  $x = \log(D) \in [-M, M]$  for some sufficiently large value of  $M$ . The new form is sometimes called 4-parameter logistic model.

The function `multiple` finds locally multiple-objective optimal designs for estimating the model parameters, the ED50, and the MED, simultaneously. For more details, see Hyun and Wong (2015).

**Usage**

```
multiple(minDose, maxDose, iter, k, inipars, Hill_par = FALSE, delta,
         lambda, fimfunc = NULL, ICA.control = list(),
         sens.minimax.control = list(), initial = NULL,
         tol = sqrt(.Machine$double.xmin))
```

**Arguments**

<code>minDose</code>	Minimum dose $D$ . For the 4-parameter logistic model, i.e. when <code>Hill_par = FALSE</code> , it is the minimum of $\log(D)$ .
<code>maxDose</code>	Maximum dose $D$ . For the 4-parameter logistic model, i.e. when <code>Hill_par = FALSE</code> , it is the maximum of $\log(D)$ .
<code>iter</code>	Maximum number of iterations.
<code>k</code>	Number of design points. Must be at least equal to the number of model parameters to avoid singularity of the FIM.
<code>inipars</code>	A vector of initial estimates for the vector of parameters $(a, b, c, d)$ . For the 4-parameter logistic model, i.e. when <code>Hill_par = FALSE</code> , it is a vector of initial estimates for $(\theta_1, \theta_2, \theta_3, \theta_4)$ .
<code>Hill_par</code>	Hill model parameterization? Defaults to TRUE.
<code>delta</code>	Predetermined meaningful value of the minimum effective dose MED. When $\delta < 0$ , then $\theta_2 > 0$ or when $\delta > 0$ , then $\theta_2 < 0$ .

<code>lambda</code>	A vector of relative importance of each of the three criteria, i.e. $\lambda = (\lambda_1, \lambda_2, \lambda_3)$ . Here $0 < \lambda_i < 1$ and $\sum \lambda_i = 1$ .
<code>fimfunc</code>	A function. Returns the FIM as a matrix. Required when formula is missing. See 'Details' of <code>minimax</code> .
<code>ICA.control</code>	ICA control parameters. For details, see <code>ICA.control</code> .
<code>sens.minimax.control</code>	Control parameters to verify the general equivalence theorem. For details, see the function <code>sens.minimax.control</code> .
<code>initial</code>	A matrix of the initial design points and weights that will be inserted into the initial solutions (countries) of the algorithm. Every row is a design, i.e. a concatenation of <code>x</code> and <code>w</code> . Will be coerced to a matrix if necessary. See 'Details' of <code>minimax</code> .
<code>tol</code>	Tolerance for finding the general inverse of the Fisher information matrix. Defaults to <code>.Machine\$double.xmin</code> .

### Details

When  $\lambda_1 > 0$ , then the number of support points `k` must at least be four to avoid singularity of the Fisher information matrix.

One can adjust the tuning parameters in `ICA.control` to set a stopping rule based on the general equivalence theorem. See 'Examples' below.

### Value

an object of class `minimax` that is a list including three sub-lists:

`arg` A list of design and algorithm parameters.

`evol` A list of length equal to the number of iterations that stores the information about the best design (design with least criterion value) of each iteration. `evol[[iter]]` contains:

<code>iter</code>	Iteration number.
<code>x</code>	Design points.
<code>w</code>	Design weights.
<code>min_cost</code>	Value of the criterion for the best imperialist (design).
<code>mean_cost</code>	Mean of the criterion values of all the imperialists.
<code>sens</code>	An object of class 'sensminimax'. See below.
<code>param</code>	Vector of parameters.

`empires` A list of all the empires of the last iteration.

`alg` A list with following information:

<code>nfeval</code>	Number of function evaluations. It does not count the function evaluations from checking the general equivalence theorem.
<code>nlocal</code>	Number of successful local searches.
<code>nrevol</code>	Number of successful revolutions.
<code>nimprove</code>	Number of successful movements toward the imperialists in the assimilation step.
<code>convergence</code>	Stopped by 'maxiter' or 'equivalence'?

The list `sens` contains information about the design verification by the general equivalence theorem. See `sensminimax` for more details. It is given every `ICA.control$checkfreq` iterations and also the last iteration if `ICA.control$checkfreq >= 0`. Otherwise, `NULL`.

`param` is a vector of parameters that is the global minimum of the minimax criterion or the global maximum of the standardized maximin criterion over the parameter space, given the current `x`, `w`.

### Note

This function is NOT appropriate for finding c-optimal designs for estimating 'MED' or 'ED50' (single objective optimal designs) and the results may not be stable. The reason is that for the c-optimal criterion the generalized inverse of the Fisher information matrix is not stable and depends on the tolerance value (`tol`).

### References

Hyun, S. W., and Wong, W. K. (2015). Multiple-Objective Optimal Designs for Studying the Dose Response Function and Interesting Dose Levels. The international journal of biostatistics, 11(2), 253-271.

### See Also

[sensmultiple](#)

### Examples

```
# All th examples here are available in Hyun and Wong (2015)

#####
# 4-parameter logistic model
# Example 1, Table 3
#####
lam <- c(0.05, 0.05, .90)
# Initial estimates are derived from Table 1
# See how the stopping rules are set via 'stop_rul', 'checkfreq' and 'stoptol'
Theta1 <- c(1.563, 1.790, 8.442, 0.137)
res1 <- multiple(minDose = log(.001), maxDose = log(1000),
                inipars = Theta1, k = 4, lambda = lam, delta = -1,
                iter = 1,
                ICA.control = list(rseed = 1366, ncount = 100,
                                   stop_rule = "equivalence",
                                   checkfreq = 100, stoptol = .95))

## Not run:
res1 <- iterate(res1, 1000)
# stops at iteration 101

## End(Not run)

#####
# 4-parameter Hill model
```

```
#####
## initial estimates for the parameters of Hill model:
a <- 0.008949 # ED50
b <- -1.79 # Hill constant
c <- 0.137 # lower limit
d <- 1.7 # upper limit
# D belongs to c(.001, 1000) ## dose in mg
## the vector of Hill parameters are now c(a, b, c, d)
## Not run:
res2 <- multiple(minDose = .001, maxDose = 1000,
                 inipars = c(a, b, c, d),
                 Hill_par = TRUE, k = 4, lambda = lam,
                 delta = -1, iter = 1000,
                 ICA.control = list(rseed = 1366, ncount = 100,
                                    stop_rule = "equivalence",
                                    checkfreq = 100, stoptol = .95))

# stops at iteration 100

## End(Not run)
```

---

normal	<i>Assumes A Multivariate Normal Prior Distribution for The Model Parameters</i>
--------	--

---

### Description

Creates a multivariate normal prior distribution for the unknown parameters as an object of class `cprior`.

### Usage

```
normal(mu, sigma, lower, upper)
```

### Arguments

mu	A vector representing the mean values.
sigma	A symmetric positive-definite matrix representing the variance-covariance matrix of the distribution.
lower	A vector of lower bounds for the model parameters.
upper	A vector of upper bounds for the model parameters.

**Value**

An object of class `cprior` that is a list with the following components:

- `fn`: prior distribution as an R function with argument `param` that is the vector of the unknown parameters. See below.
- `npar`: Number of unknown parameters and is equal to the length of `param`.
- `lower`: Argument `lower`. It has the same length as `param`.
- `upper`: Argument `lower`. It has the same length as `param`.

The list will be passed to the argument `prior` of the function `bayes`. The order of the argument `param` in `fn` has the same order as the argument `parvars` when the model is specified by a formula. Otherwise, it is equal to the argument `param` in the function `fimfunc`.

**See Also**

[bayes](#) [sensbayes](#)

**Examples**

```
normal(mu = c(0, 1), sigma = matrix(c(1, -0.17, -0.17, .5), nrow = 2),
      lower = c(-3, .1), upper = c(3, 2))
```

---

plot.bayes

*Plotting bayes Objects*

---

**Description**

This function plots the evolution of the ICA algorithm (iteration vs the best (minimum) criterion value at each iteration) and also verifies the optimality of the last obtained design using the general equivalence theorem. It plots the sensitivity function and calculates the ELB for the best design generated at iteration number `iter`.

**Usage**

```
## S3 method for class 'bayes'
plot(x, iter = NULL, sensitivity = TRUE,
     calculate_criterion = FALSE, crt_method = NULL, sens_method = NULL,
     sens.bayes.control = list(), crt.bayes.control = list(),
     silent = FALSE, plot_3d = c("lattice", "rgl"), evolution = FALSE,
     ...)
```

**Arguments**

<code>x</code>	An object of class <code>bayes</code> .
<code>iter</code>	Iteration number. if NULL (default), it will be set to the last iteration.
<code>sensitivity</code>	Logical. If TRUE (default), the general equivalence theorem is used to check the optimality if the best design in iteration number <code>iter</code> and the sensitivity function will be plotted.
<code>calculate_criterion</code>	Logical. Re-calculate the Bayesian criterion value (maybe with a set of new tuning parameters to be sure of the accuracy of the integral approximation)? Defaults to FALSE. See 'Details'.
<code>crt_method</code>	A character denotes which method to be used to approximate the integrals in Bayesian criteria. "cubature" corresponds to the adaptive multivariate integration method using the <a href="#">hcubature</a> algorithm (default). "quadrature" corresponds the traditional quadrature formulas and calls the function <a href="#">createNIGrid</a> . The tuning parameters are adjusted by <code>crt.bayes.control</code> . If NULL (default), it will use the method already applied to create the object <code>x</code> .
<code>sens_method</code>	Similar to <code>crt_method</code> , but for approximating the integrals in the sensitivity (derivative) function. The tuning parameters are adjusted by <code>sens.bayes.control</code> . If NULL (default), it will use the method already applied to create the object <code>x</code> .
<code>sens.bayes.control</code>	Control parameters to verify general equivalence theorem. For details, see <a href="#">sens.bayes.control</a> . If NULL (default), it will be set to the tuning parameters used to create object <code>x</code> .
<code>crt.bayes.control</code>	Control parameters to approximate the integration in the Bayesian criterion at a given design. For details, see <a href="#">crt.bayes.control</a> . If NULL (default), it will be set to the tuning parameters used to create object <code>x</code> .
<code>silent</code>	Do not print anything? Defaults to FALSE.
<code>plot_3d</code>	Which package should be used to plot the sensitivity function for two-dimensional design space. Defaults to <code>plot_3d = "lattice"</code> . Only applicable when <code>sensitivity = TRUE</code> .
<code>evolution</code>	Plot Evolution? Defaults to FALSE.
<code>...</code>	Argument with no further use.

**Details**

In addition to verifying the general equivalence theorem, this function makes it possible to re-calculate the criterion value for the output designs using a new (say, more conservative) set of tuning parameters. This is useful for Bayesian optimal designs to assess the robustness of the criterion value with respect to different values of the tuning parameters. To put it simple, for Bayesian generated designs, the user can re-calculate the criterion value for the output design (best design generated in `iter`) with different values for `maxEval` and `tol` in [crt.bayes.control](#) to verify that the function `hcubature` approximates the integrals to an user-acceptable accuracy. The same also applies for the quadrature methods using different number of nodes.

**See Also**

[bayes](#), [bayescomp](#)

---

plot.minimax

*Plotting minimax Objects*

---

**Description**

This function plots the evolution of the ICA algorithm (iteration vs the best (minimum) criterion value at each iteration) and also verifies the optimality of the last obtained design using the general equivalence theorem. It plots the sensitivity function and calculates the ELB for the best design generated at iteration number `iter`.

**Usage**

```
## S3 method for class 'minimax'
plot(x, iter = NULL, sensitivity = TRUE,
     calculate_criterion = FALSE, sens.minimax.control = list(),
     crt.minimax.control = list(), silent = FALSE,
     plot_3d = c("lattice", "rgl"), evolution = FALSE, ...)
```

**Arguments**

<code>x</code>	An object of class <code>minimax</code> .
<code>iter</code>	Iteration number. if <code>NULL</code> (default), it will be set to the last iteration.
<code>sensitivity</code>	Logical. If <code>TRUE</code> (default), the general equivalence theorem is used to check the optimality if the best design in iteration number <code>iter</code> and the sensitivity function will be plotted.
<code>calculate_criterion</code>	Logical. Re-calculate the criterion value (maybe with a set of new tuning parameters to be sure of the globality of the maximum over the parameter space given the design)? It only assumes a continuous parameter space for the minimax and standardized maximin designs. Defaults to <code>FALSE</code> . See 'Details'.
<code>sens.minimax.control</code>	Control parameters to verify general equivalence theorem. For details, see <a href="#">sens.minimax.control</a> . If <code>NULL</code> (default), it will be set to the tuning parameters used to create object <code>x</code> .
<code>crt.minimax.control</code>	Control parameters to optimize the minimax or standardized maximin criterion at a given design over a <b>continuous</b> parameter space. For details, see <a href="#">crt.minimax.control</a> . If <code>NULL</code> (default), it will be set to the tuning parameters used to create object <code>x</code> .
<code>silent</code>	Do not print anything? Defaults to <code>FALSE</code> .
<code>plot_3d</code>	Which package should be used to plot the sensitivity function for two-dimensional design space. Defaults to <code>plot_3d = "lattice"</code> . Only applicable when <code>sensitivity = TRUE</code> .
<code>evolution</code>	Plot Evolution? Defaults to <code>FALSE</code> .
<code>...</code>	Argument with no further use.

## Details

In addition to verifying the general equivalence theorem, this function makes it possible to re-calculated the criterion value for the output designs using a new set of tuning parameters, especially, a large value for `maxeval` in the function `crt.minimax.control`. This is useful for minimax and standardized maximin optimal designs to assess the robustness of the criterion value with respect to different values of `maxeval`. To put it simple, for these designs, the user can re-calculate the criterion value (finds the global maximum over the parameter space given an output design in a minimax problem) with larger values for `maxeval` in `crt.minimax.control` to be sure that the function `nloptr` finds global optima of the inner optimization problem over the parameter space using the default value (or the user-given value) of `maxeval`. If increasing the value of `maxeval` returns different criterion values, then the results can not be trusted and the algorithm should be repeated with a higher value for `maxeval`.

## See Also

[minimax](#), [locally](#), [robust](#)

---

print.bayes

*Printing bayes Objects*

---

## Description

Print method for an object of class bayes.

## Usage

```
## S3 method for class 'bayes'  
print(x, iter = NULL, ...)
```

## Arguments

<code>x</code>	an object of class bayes.
<code>iter</code>	Iteration number. if NULL, it will be set to the last iteration.
<code>...</code>	Argument with no further use.

## See Also

[bayes](#)

print.minimax

*Printing minimax Objects*

---

**Description**

Print method for an object of class minimax.

**Usage**

```
## S3 method for class 'minimax'  
print(x, iter = NULL, ...)
```

**Arguments**

x	An object of class minimax.
iter	Iteration number. if NULL, will be set equal to the last iteration.
...	Argument with no further use.

**See Also**

[minimax](#), [locally](#), [robust](#)

---

print.sensbayes

*Printing sensbayes Objects*

---

**Description**

Print method for an object of class sensbayes.

**Usage**

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

**Arguments**

x	An object of class sensbayes.
...	Argument with no further use.

**See Also**

[sensbayes](#), [sensbayescomp](#)

---

print.sensminimax	<i>Printing sensminimax Objects</i>
-------------------	-------------------------------------

---

**Description**

Print method for an object of class sensminimax.

**Usage**

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

**Arguments**

x	An object of class sensminimax.
...	Argument with no further use.

**See Also**

[sensminimax](#), [senslocally](#), [sensrobust](#)

---

robust	<i>Robust D-Optimal Designs</i>
--------	---------------------------------

---

**Description**

Finds Robust designs or optimal in-average designs for nonlinear models. It is useful when a set of different vectors of initial estimates along with a discrete probability measure are available for the unknown model parameters. It is a discrete version of [bayes](#).

**Usage**

```
robust(formula, predvars, parvars, family = gaussian(), lx, ux, iter, k,
       prob, parset, fimfunc = NULL, ICA.control = list(),
       sens.minimax.control = list(), initial = NULL,
       npar = dim(parset)[2], plot_3d = c("lattice", "rgl"), x = NULL)
```

**Arguments**

formula	A nonlinear model <a href="#">formula</a> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <a href="#">formula</a> if necessary.
predvars	A vector of characters. Denotes the predictors in the <a href="#">formula</a> .
parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .

family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <code>gaussian()</code> is applied.
lx	Vector of lower bounds for the predictors. Should be in the same order as <code>predvars</code> .
ux	Vector of upper bounds for the predictors. Should be in the same order as <code>predvars</code> .
iter	Maximum number of iterations.
k	Number of design points. Must be at least equal to the number of model parameters to avoid singularity of the FIM.
prob	A vector of the probability measure $\pi$ associated with each row of <code>parset</code> .
parset	A matrix that provides the vector of initial estimates for the model parameters, i.e. support of $\pi$ . Every row is one vector ( <code>nrow(parset) == length(prob)</code> ). See 'Details'.
fimfunc	A function. Returns the FIM as a matrix. Required when <code>formula</code> is missing. See 'Details' of <a href="#">minimax</a> .
ICA.control	ICA control parameters. For details, see <a href="#">ICA.control</a> .
sens.minimax.control	Control parameters to verify the general equivalence theorem. For details, see the function <a href="#">sens.minimax.control</a> .
initial	A matrix of the initial design points and weights that will be inserted into the initial solutions (countries) of the algorithm. Every row is a design, i.e. a concatenation of <code>x</code> and <code>w</code> . Will be coerced to a matrix if necessary. See 'Details' of <a href="#">minimax</a> .
npar	Number of model parameters. Used when <code>fimfunc</code> is given instead of <code>formula</code> to specify the number of model parameters. If not given, the sensitivity plot may be shifted below the y-axis. When <code>NULL</code> , it is set to <code>dim(parset)[2]</code> .
plot_3d	Which package should be used to plot the sensitivity (derivative) function for two-dimensional design space. Defaults to "lattice".
x	A vector of candidate design (support) points. When is not set to <code>NULL</code> (default), the algorithm only finds the optimal weights for the candidate points in <code>x</code> . Should be set when the user has a finite number of candidate design points and the purpose is to find the optimal weight for each of them (when zero, they will be excluded from the design). For design points with more than one dimension, see 'Details' of <a href="#">sensminimax</a> .

## Details

Let  $\Theta$  be a set of initial estimates for the unknown parameters. A robust criterion is evaluated at the elements of  $\Theta$  weighted by a probability measure  $\pi$  as follows:

$$B(\xi, \pi) = \int_{\Theta} |M(\xi, \theta)| \pi(\theta) d\theta.$$

A robust design  $\xi^*$  maximizes  $B(\xi, \pi)$  over the space of all designs.

When the model is given via formula, columns of `parset` must match the parameters introduced in `parvars`. Otherwise, when the model is introduced via `fimfunc`, columns of `parset` must match the argument `param` in `fimfunc`.

To verify the optimality of the output design by the general equivalence theorem, the user can either plot the results or set `checkfreq` in `ICA.control` to `Inf`. In either way, the function `sensrobust` is called for verification. One can also adjust the tuning parameters in `ICA.control` to set a stopping rule based on the general equivalence theorem. See 'Examples' below.

## Value

an object of class `minimax` that is a list including three sub-lists:

`arg` A list of design and algorithm parameters.

`evol` A list of length equal to the number of iterations that stores the information about the best design (design with least criterion value) of each iteration. `evol[[iter]]` contains:

<code>iter</code>	Iteration number.
<code>x</code>	Design points.
<code>w</code>	Design weights.
<code>min_cost</code>	Value of the criterion for the best imperialist (design).
<code>mean_cost</code>	Mean of the criterion values of all the imperialists.
<code>sens</code>	An object of class 'sensminimax'. See below.
<code>param</code>	Vector of parameters.

`empires` A list of all the empires of the last iteration.

`alg` A list with following information:

<code>nfeval</code>	Number of function evaluations. It does not count the function evaluations from checking the general equivalence theorem.
<code>nlocal</code>	Number of successful local searches.
<code>nrevol</code>	Number of successful revolutions.
<code>nimprove</code>	Number of successful movements toward the imperialists in the assimilation step.
<code>convergence</code>	Stopped by 'maxiter' or 'equivalence'?

The list `sens` contains information about the design verification by the general equivalence theorem. See `sensminimax` for more details. It is given every `ICA.control$checkfreq` iterations and also the last iteration if `ICA.control$checkfreq`  $\geq 0$ . Otherwise, `NULL`.

`param` is a vector of parameters that is the global minimum of the minimax criterion or the global maximum of the standardized maximin criterion over the parameter space, given the current `x`, `w`.

## Note

When a continuous prior distribution for the unknown model parameters is available, use `bayes`. When only one initial estimates of the unknown model parameters is available ( $\Theta$  has only one element), use `locally`.

**See Also**[bayes sensrobust](#)**Examples**

```

# Finding a robust design for the two-parameter logistic model
# See how we set a stopping rule.
# The ELB is computed every checkfreq = 30 iterations
# The optimization stops when the ELB is larger than stoptol = .95
res1 <- robust(formula = ~1/(1 + exp(-b *(x - a))),
               predvars = c("x"), parvars = c("a", "b"),
               family = binomial(),
               lx = -5, ux = 5, prob = rep(1/4, 4),
               parset = matrix(c(0.5, 1.5, 0.5, 1.5, 4.0, 4.0, 5.0, 5.0), 4, 2),
               iter = 1, k =3,
               ICA.control = list(stop_rule = "equivalence",
                                  stoptol = .95, checkfreq = 30))

## Not run:
res1 <- iterate(res1, 100)
# stops at iteration 51

## End(Not run)

## Not run:
res1.1 <- robust(formula = ~1/(1 + exp(-b *(x - a))),
                 predvars = c("x"), parvars = c("a", "b"),
                 family = binomial(),
                 lx = -5, ux = 5, prob = rep(1/4, 4),
                 parset = matrix(c(0.5, 1.5, 0.5, 1.5, 4.0, 4.0, 5.0, 5.0), 4, 2),
                 x = c(-3, 0, 3),
                 iter = 150, k =3)

plot(res1.1)
# not optimal

## End(Not run)

```

---

sens.bayes.control      *Control Parameters for Verifying General Equivalence Theorem for Bayesian Designs*

---

**Description**

This function returns two lists each corresponds to an implemented integration method for approximating the integrals in the sensitivity (derivative) functions for Bayesian criteria. Moreover, it contains a list of `nloptr` control parameters to find maximum of the sensitivity (derivative) function over the design space, used to calculate the efficiency lower bound (ELB).

**Usage**

```
sens.bayes.control(cubature = list(tol = 1e-05, maxEval = 50000, absError
= 0), quadrature = list(type = NULL, level = NULL, ndConstruction =
"product", level.trans = FALSE), x0 = NULL, optslist = list(stopval =
-Inf, algorithm = "NLOPT_GN_DIRECT_L", xtol_rel = 1e-08, ftol_rel =
1e-10, maxeval = 2000), ...)
```

**Arguments**

cubature	A list that will be passed to the arguments of the <a href="#">hcubature</a> function. See 'Details' of <a href="#">crt.bayes.control</a> .
quadrature	A list that will be passed to the arguments of the <a href="#">createNIGrid</a> function. See 'Details' of <a href="#">crt.bayes.control</a> .
x0	Vector of starting values for maximizing the sensitivity (derivative) function over the design space $x$ . It will be passed to the optimization function <a href="#">nloptr</a> .
optslist	A list will be passed to opts argument of the function <a href="#">nloptr</a> to find the maximum of the sensitivity function over the design space. See 'Details'.
...	Further arguments will be passed to <a href="#">nl.opts</a> from package <a href="#">nloptr</a> .

**Details**

ELB is a measure of proximity of a design to the optimal design without knowing the latter. Given a design, let  $\epsilon$  be the global maximum of the sensitivity (derivative) function with respect the vector of the model predictors  $x$  over the design space. ELB is given by

$$ELB = p/(p + \epsilon),$$

where  $p$  is the number of model parameters. Obviously, calculating ELB requires finding  $\epsilon$  and therefore, a maximization problem to be solved. The function [nloptr](#) is used here to solve this maximization problem. The arguments  $x0$  and [optslist](#) will be passed to this function as follows:

Argument  $x0$  provides the user initial values for this maximization problem and will be passed to the argument with the same name in the function [nloptr](#).

Argument [optslist](#) is a list and it will be passed to the argument [opts](#) of the function [nloptr](#). The most important components of [optslist](#) are:

[stopval](#) Stop minimization when an objective value  $\leq$  [stopval](#) is found. Setting [stopval](#) to  $-\text{Inf}$  disables this stopping criterion (default).

[algorithm](#) Defaults to `NLOPT_GN_DIRECT_L`. DIRECT-L is a deterministic-search algorithm based on systematic division of the search domain into smaller and smaller hyperrectangles.

[xtol\\_rel](#) Stop when an optimization step (or an estimate of the optimum) changes every parameter by less than [xtol\\_rel](#) multiplied by the absolute value of the parameter. Criterion is disabled if [xtol\\_rel](#) is non-positive. Defaults to  $1e-8$ .

[ftol\\_rel](#) Stop when an optimization step (or an estimate of the optimum) changes the objective function value by less than [ftol\\_rel](#) multiplied by the absolute value of the function value. Criterion is disabled if [ftol\\_rel](#) is non-positive. Defaults to  $1e-10$ .

[maxeval](#) Stop when the number of function evaluations exceeds [maxeval](#). Criterion is disabled if [maxeval](#) is non-positive. Defaults to  $6000$ . See 'Note' on when to change its value.

More details are available by the function `nloptr.print.options()` in package [nloptr](#).

**Value**

A list of three lists each contains the control parameters for verifying the general equivalence theorem with respect to the Bayesian optimality criteria.

**Note**

Whenever the value of ELB is computationally larger than 1, it is a sign that the obtained  $\epsilon$  is not global maximum. To overcome this issue, please increase the value of the parameter maxeval to allow the optimization algorithm to find the global maximum of the sensitivity (derivative) function over the design space.

Please note the difference between maxEval in cubature and maxeval in optslst. They are quite two types of different tuning parameters. The earlier is the (approximate) maximum number of function evaluations in the hcubature adaptive integration method and the latter is the maximum number of function evaluations in the maximization problem.

**Examples**

```
sens.bayes.control()
sens.bayes.control(cubature = list(maxEval = 50000))
sens.bayes.control(optslst = list(maxeval = 3000))
sens.bayes.control(quadrature = list(level = 4))
```

---

sens.minimax.control    *Control Parameters for Verifying General Equivalence Theorem*

---

**Description**

This function returns a list of control parameters that are used to find the “answering set” for minimax and standardized maximin generated designs. The answering set is required to obtain the sensitivity (derivative) function in order to verify the optimality of a given design. Moreover, it contains a list of `nloptr` control parameters to find maximum of the sensitivity (derivative) function over the design space, used to calculate the efficiency lower bound (ELB).

**Usage**

```
sens.minimax.control(answering.set = list(n_seg = 6, merge_tol = 0.005),
  x0 = NULL, optslst = list(stopval = -Inf, algorithm =
  "NLOPT_GN_DIRECT_L", xtol_rel = 1e-08, ftol_rel = 1e-10, maxeval = 6000),
  ...)
```

**Arguments**

answering.set	A list of control parameters to find the answering set in minimax and standardized maximin optimal design problems. See 'Details'.
x0	Vector of starting values for maximizing the sensitivity (derivative) function over the design space $x$ . It will be passed to the optimization function <code>nloptr</code> .

optslist      A list. It will be passed to the argument `opts` of the function `nloptr` to find the maximum of the sensitivity function over the design space. See 'Details'.

...            Further arguments will be passed to `nl.opts` from package `nloptr`.

## Details

Given a design, an “answering set” is a subset of all local optima of the optimality criterion over the parameter space (only applicable for minimax and standardized maximin design problems). For an optimality verification by the general equivalence theorem, finding the true answering set is important because the sensitivity (derivative) function in a minimax or standardized maximin problem is obtained based on it. An invalid answering set may result in a false sensitivity plot and ELB. Unfortunately, there is no theoretical rule on how to choose the number of elements of an answering set, and they would have to be found by trial and error. Given a design, the answering set for a minimax criterion is obtained as follows:

- Step 1: Find all the local maxima of the optimality criterion (minimax) over the parameter space. For this purpose, in our algorithm, the parameter space is divided into  $(n\_seg + 1)^p$  segments, where  $p$  is the number of unknown model parameters. Then, each boundary point of the resulted segments (intervals) is assigned to the argument `par` of the function `optim` in order to start a local search using the “L-BFGS-B” method.
- Step 2: Pick the ones nearest to the global minimum subject to a merging tolerance `merge_tol` (default `0.005`).

The resulting maxima establish the answering set. Obviously, the answering set is a subset of all the local maxima. Therefore, it is very important to be able to find all the local maxima to create the true answering set with no missing elements. Otherwise, even when the design is optimal, the sensitivity (derivative) plot may not verify the optimality of the design. Finding all the local optimal of a multimodal function like the sensitivity function is not an easy task and this is the main reason that checking the general equivalence theorem (even plotting) in minimax and standardized maximin problems is complicated. As a result, the argument `answering.set` is a list with two elements:

`n_seg` Please increase the value of `n_seg` for models with larger number of unknown parameters or large parameter space.

`merge_tol` We advise to not change the default value of the parameter `merge_tol` as it has been successfully tested for many examples.

ELB is a measure of proximity of a design to the optimal design without knowing the latter. Given a design, let  $\epsilon$  be the global maximum of the sensitivity (derivative) function with respect the vector of the model predictors  $x$  over the design space. ELB is given by

$$ELB = p/(p + \epsilon),$$

where  $p$  is the number of model parameters. Obviously, calculating ELB requires finding  $\epsilon$  and therefore, a maximization problem to be solved. The function `nloptr` is used here to solve this maximization problem. The arguments `x0` and `optslist` will be passed to this function as follows: Argument `x0` provides the user initial values for this maximization problem and will be passed to the argument with the same name in the function `nloptr`.

Argument `optslist` will be passed to the argument `opts` of the function `nloptr`. `optslist` is a list and the most important components are listed as follows:

`stopval` Stop minimization when an objective value  $\leq$  `stopval` is found. Setting `stopval` to `-Inf` disables this stopping criterion (default).

`algorithm` Defaults to `NLOPT_GN_DIRECT_L`. `DIRECT-L` is a deterministic-search algorithm based on systematic division of the search domain into smaller and smaller hyperrectangles.

`xtol_rel` Stop when an optimization step (or an estimate of the optimum) changes every parameter by less than `xtol_rel` multiplied by the absolute value of the parameter. Criterion is disabled if `xtol_rel` is non-positive. Defaults to  $1e-8$ .

`ftol_rel` Stop when an optimization step (or an estimate of the optimum) changes the objective function value by less than `ftol_rel` multiplied by the absolute value of the function value. Criterion is disabled if `ftol_rel` is non-positive. Defaults to  $1e-10$ .

`maxeval` Stop when the number of function evaluations exceeds `maxeval`. Criterion is disabled if `maxeval` is non-positive. Defaults to `6000`. See 'Note' on when to change its value.

A detailed explanation of all the options is shown by the function `nloptr.print.options()` in package `nloptr`.

### Value

A list of control parameters for verifying the general equivalence theorem for minimax, standardized maximin and locally optimal designs.

### Note

Whenever the value of ELB is computationally larger than 1, it is a sign that the obtained  $\epsilon$  is not global maximum. To overcome this issue, please increase the value of the parameter `maxeval` to allow the optimization algorithm to find the global maximum of the sensitivity (derivative) function over the design space.

### Examples

```
sens.minimax.control()
sens.minimax.control(answering.set = list(n_seg = 4))
sens.minimax.control(answering.set = list(n_seg = 4), optslist = list(maxeval = 1000))
# faster checking process
sens.minimax.control(answering.set = list(n_seg = 4), optslist = list(maxeval = 2000))
```

### Description

Plots the sensitivity (derivative) function and calculates the efficiency lower bound (ELB) for a given Bayesian design. Let  $\boldsymbol{x}$  belongs to  $\chi$  that denotes the design space. Based on the general equivalence theorem, a design  $\xi^*$  is optimal if and only if the value of the sensitivity (derivative) function is non-positive for all  $\boldsymbol{x}$  in  $\chi$  and zero when  $\boldsymbol{x}$  belongs to the support of  $\xi^*$  (be equal to the one of the design points).

For an approximate (continuous) design, when the design space is one or two-dimensional, the user can visually verify the optimality of the design by observing the sensitivity plot. Furthermore, the approximity of the design to the optimal design can be measured by the ELB without knowing the latter.

## Usage

```
sensbayes(formula, predvars, parvars, family = gaussian(), x, w, lx, ux,
  fimfunc = NULL, prior = list(), sens.bayes.control = list(),
  crt.bayes.control = list(), crt_method = c("cubature", "quadrature"),
  sens_method = c("cubature", "quadrature"), plot_3d = c("lattice",
  "rgl"), plot_sens = TRUE, npar = NULL, calculate_criterion = TRUE,
  silent = FALSE)
```

## Arguments

formula	A nonlinear model <a href="#">formula</a> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <a href="#">formula</a> if necessary.
predvars	A vector of characters. Denotes the predictors in the <a href="#">formula</a> .
parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .
family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <a href="#">gaussian()</a> is applied.
x	A vector of candidate design (support) points. When is not set to NULL (default), the algorithm only finds the optimal weights for the candidate points in x. Should be set when the user has a finite number of candidate design points and the purpose is to find the optimal weight for each of them (when zero, they will be excluded from the design). For design points with more than one dimension, see 'Details' of <a href="#">sensminimax</a> .
w	Vector of the corresponding design weights for x.
lx	Vector of lower bounds for the predictors. Should be in the same order as predvars.
ux	Vector of upper bounds for the predictors. Should be in the same order as predvars.
fimfunc	A function. Returns the FIM as a matrix. Required when formula is missing. See 'Details' of <a href="#">minimax</a> .
prior	An object of class cprior. User can also use one of the functions <a href="#">uniform</a> , <a href="#">normal</a> , <a href="#">skewnormal</a> or <a href="#">student</a> to create the prior. See 'Details' of <a href="#">bayes</a> .
sens.bayes.control	A list. Control parameters to verify the general equivalence theorem. For details, see <a href="#">sens.bayes.control</a> .

<code>crt.bayes.control</code>	A list. Control parameters to approximate the integral in the Bayesian criterion at a given design over the parameter space. For details, see <code>crt.bayes.control</code> .
<code>crt_method</code>	A character denotes which method to be used to approximate the integrals in Bayesian criteria. "cubature" corresponds to the adaptive multivariate integration method using the <code>hcubature</code> algorithm (default). "quadrature" corresponds the traditional quadrature formulas and calls the function <code>createNIGrid</code> . The tuning parameters are adjusted by <code>crt.bayes.control</code> . Default is set to "cubature".
<code>sens_method</code>	Similar to <code>crt_method</code> , but for approximating the integrals in the sensitivity (derivative) function. The tuning parameters are adjusted by <code>sens.bayes.control</code> . Default is set to "cubature".
<code>plot_3d</code>	Which package should be used to plot the sensitivity (derivative) function for two-dimensional design space. Defaults to "lattice".
<code>plot_sens</code>	Plot the sensitivity (derivative) function? Defaults to TRUE.
<code>npar</code>	Number of model parameters. Used when <code>fimfunc</code> is given instead of formula to specify the number of model parameters. If not specified correctly, the sensitivity (derivative) plot may be shifted below the y-axis. When NULL (default), it will be set to <code>length(parvars)</code> or <code>prior\$npar</code> when missing(formula).
<code>calculate_criterion</code>	Calculate the optimality criterion? See 'Details' of <code>sensminimax</code> .
<code>silent</code>	Do not print anything? Defaults to FALSE.

## Details

Let  $\Xi$  be the space of all approximate designs with  $k$  design points (support points) at  $x_1, x_2, \dots, x_k$  from design space  $\chi$  with corresponding weights  $w_1, \dots, w_k$ . Let  $M(\xi, \theta)$  be the Fisher information matrix (FIM) of a  $k$ -point design  $\xi$  and  $\pi(\theta)$  is a user-given prior distribution for the vector of unknown parameters  $\theta$ . A design  $\xi^*$  is Bayesian D-optimal among all designs on  $\chi$  if and only if the following inequality holds for all  $\mathbf{x} \in \chi$

$$c(\mathbf{x}, \xi^*) = \int_{\theta \in \Theta} \text{tr} M^{-1}(\xi^*, \theta) I(\mathbf{x}, \theta) - p\pi(\theta) d\theta \leq 0,$$

with equality at all support points of  $\xi^*$ . Here,  $p$  is the number of model parameters.  $c(\mathbf{x}, \xi^*)$  is called **sensitivity** or **derivative** function.

Depending on the complexity of the problem at hand, sometimes, the CPU time can be considerably reduced by choosing a set of less conservative values for the tuning parameters `tol` and `maxEval` in the function `sens.bayes.control` when `sens_method = "cubature"`. Similarly, this applies when `sens_method = "quadrature"`. In general, if the CPU time matters, the user should find an appropriate speed-accuracy trade-off for her/his own problem. See 'Examples' for more details.

## Note

The default values of the tuning parameters in `sens.bayes.control` are set in a way that having accurate plots for the sensitivity (derivative) function and calculating the ELB to a high precision to be the primary goals, although the process may take too long. The user should choose a set of less conservative values via the argument `sens.bayes.control` when the CPU-time is too long or matters.

## Examples

```
#####
# Checking the Bayesian D-optimality of a design for the 2Pl model
#####
skew2 <- skewnormal(xi = c(0, 1), Omega = matrix(c(1, -0.17, -0.17, .5), nrow = 2),
              alpha = c(-1, 0), lower = c(-3, .1), upper = c(3, 2))

## Not run:
sensbayes(formula = ~1/(1 + exp(-b *(x - a))),
          predvars = "x", parvars = c("a", "b"),
          family = binomial(),
          x= c(-2.50914, -1.16780, -0.36904, 1.29227),
          w =c(0.35767, 0.11032, 0.15621, 0.37580),
          lx = -3, ux = 3,
          prior = skew2)
# took 29 seconds on my system!

## End(Not run)

# It took very long.
# We re-adjust the tuning parameters in sens.bayes.control to be faster
# See how we drastically reduce the maxEval and increase the tolerance
## Not run:
sensbayes(formula = ~1/(1 + exp(-b *(x - a))),
          predvars = "x", parvars = c("a", "b"),
          family = binomial(),
          x= c(-2.50914, -1.16780, -0.36904, 1.29227),
          w =c(0.35767, 0.11032, 0.15621, 0.37580),
          lx = -3, ux = 3,prior = skew2,
          sens.bayes.control = list(cubature = list(tol = 1e-4, maxEval = 300)))
# took 5 Seconds on my system!

## End(Not run)

# Compare it with the following:
sensbayes(formula = ~1/(1 + exp(-b *(x - a))),
          predvars = "x", parvars = c("a", "b"),
          family = binomial(),
          x= c(-2.50914, -1.16780, -0.36904, 1.29227),
          w =c(0.35767, 0.11032, 0.15621, 0.37580),
          lx = -3, ux = 3,prior = skew2,
          sens.bayes.control = list(cubature = list(tol = 1e-4, maxEval = 200)))
# Look at the plot!
# took 3 seconds on my system

#####
# Checking the Bayesian D-optimality of a design for the 4-parameter sigmoid emax model
#####
lb <- c(4, 11, 100, 5)
ub <- c(9, 17, 140, 10)
```

```

## Not run:
sensbayes(formula = ~ theta1 + (theta2 - theta1)*(x^theta4)/(x^theta4 + theta3^theta4),
  predvars = c("x"), parvars = c("theta1", "theta2", "theta3", "theta4"),
  x = c(0.78990, 95.66297, 118.42964,147.55809, 500),
  w = c(0.23426, 0.17071, 0.17684, 0.1827, 0.23549),
  lx = .001, ux = 500, prior = uniform(lb, ub))
# took 200 seconds on my system

## End(Not run)

# Re-adjust the tuning parameters to have it faster
## Not run:
sensbayes(formula = ~ theta1 + (theta2 - theta1)*(x^theta4)/(x^theta4 + theta3^theta4),
  predvars = c("x"), parvars = c("theta1", "theta2", "theta3", "theta4"),
  x = c(0.78990, 95.66297, 118.42964,147.55809, 500),
  w = c(0.23426, 0.17071, 0.17684, 0.1827, 0.23549),
  lx = .001, ux = 500, prior = uniform(lb, ub),
  sens.bayes.control = list(cubature = list(tol = 1e-3, maxEval = 300)))
# took 4 seconds on my system. See how much it makes difference

## End(Not run)

## Not run:
# Now we try it with quadrature. Default is 6 nodes
sensbayes(formula = ~ theta1 + (theta2 - theta1)*(x^theta4)/(x^theta4 + theta3^theta4),
  predvars = c("x"), parvars = c("theta1", "theta2", "theta3", "theta4"),
  x = c(0.78990, 95.66297, 118.42964,147.55809, 500),
  w = c(0.23426, 0.17071, 0.17684, 0.1827, 0.23549),
  sens_method = "quadrature",
  lx = .001, ux = 500, prior = uniform(lb, ub))
# 166.519 s

# use less number of nodes to see if we can reduce the CPU time
sensbayes(formula = ~ theta1 + (theta2 - theta1)*(x^theta4)/(x^theta4 + theta3^theta4),
  predvars = c("x"), parvars = c("theta1", "theta2", "theta3", "theta4"),
  x = c(0.78990, 95.66297, 118.42964,147.55809, 500),
  w = c(0.23426, 0.17071, 0.17684, 0.1827, 0.23549),
  sens_method = "quadrature",
  sens.bayes.control = list(quadrature = list(level = 3)),
  lx = .001, ux = 500, prior = uniform(lb, ub))
# we don't have an accurate plot

# use less number of levels: use 4 nodes
sensbayes(formula = ~ theta1 + (theta2 - theta1)*(x^theta4)/(x^theta4 + theta3^theta4),
  predvars = c("x"), parvars = c("theta1", "theta2", "theta3", "theta4"),
  x = c(0.78990, 95.66297, 118.42964,147.55809, 500),
  w = c(0.23426, 0.17071, 0.17684, 0.1827, 0.23549),
  sens_method = "quadrature",
  sens.bayes.control = list(quadrature = list(level = 4)),
  lx = .001, ux = 500, prior = uniform(lb, ub))

## End(Not run)

```

## Description

This function plot the sensitivity (derivative) function given an approximate (continuous) design and calculate the efficiency lower bound (ELB) for Bayesian DP-optimal designs. Let  $x$  belongs to  $\chi$  that denotes the design space. Based on the general equivalence theorem, generally, a design  $\xi^*$  is optimal if and only if the value of its sensitivity (derivative) function be non-positive for all  $x$  in  $\chi$  and it only reaches zero when  $x$  belong to the support of  $\xi^*$  (be equal to one of the design point). Therefore, the user can look at the sensitivity plot and the ELB and decide whether the design is optimal or close enough to the true optimal design (ELB tells us that without knowing the latter).

## Usage

```
sensbayescomp(formula, predvars, parvars, family = gaussian(), x, w, lx,
  ux, fimfunc = NULL, prior = list(), prob, alpha,
  sens.bayes.control = list(), crt.bayes.control = list(),
  crt_method = c("cubature", "quadrature"), sens_method = c("cubature",
  "quadrature"), plot_3d = c("lattice", "rgl"), plot_sens = TRUE,
  npar = NULL, calculate_criterion = TRUE, silent = FALSE)
```

## Arguments

formula	A nonlinear model <a href="#">formula</a> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <a href="#">formula</a> if necessary.
predvars	A vector of characters. Denotes the predictors in the <a href="#">formula</a> .
parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .
family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <code>gaussian()</code> is applied.
x	A vector of candidate design (support) points. When is not set to NULL (default), the algorithm only finds the optimal weights for the candidate points in $x$ . Should be set when the user has a finite number of candidate design points and the purpose is to find the optimal weight for each of them (when zero, they will be excluded from the design). For design points with more than one dimension, see 'Details' of <a href="#">sensminimax</a> .
w	Vector of the corresponding design weights for $x$ .
lx	Vector of lower bounds for the predictors. Should be in the same order as <code>predvars</code> .

<code>ux</code>	Vector of upper bounds for the predictors. Should be in the same order as <code>predvars</code> .
<code>fimfunc</code>	A function. Returns the FIM as a matrix. Required when <code>formula</code> is missing. See 'Details' of <a href="#">minimax</a> .
<code>prior</code>	An object of class <code>cprior</code> . User can also use one of the functions <a href="#">uniform</a> , <a href="#">normal</a> , <a href="#">skewnormal</a> or <a href="#">student</a> to create the prior. See 'Details' of <a href="#">bayes</a> .
<code>prob</code>	Either <code>formula</code> or a function. When function, its argument are <code>x</code> and <code>param</code> , and they are the same as the arguments in <code>fimfunc</code> . <code>prob</code> as a function takes the design points and vector of parameters and returns the probability of success at each design points. See 'Examples'.
<code>alpha</code>	A value between 0 and 1. Compound or combined DP-criterion is the product of the efficiencies of a design with respect to D- and average P- optimality, weighted by <code>alpha</code> .
<code>sens.bayes.control</code>	A list. Control parameters to verify the general equivalence theorem. For details, see <a href="#">sens.bayes.control</a> .
<code>crt.bayes.control</code>	A list. Control parameters to approximate the integral in the Bayesian criterion at a given design over the parameter space. For details, see <a href="#">crt.bayes.control</a> .
<code>crt_method</code>	A character denotes which method to be used to approximate the integrals in Bayesian criteria. "cubature" corresponds to the adaptive multivariate integration method using the <a href="#">hcubature</a> algorithm (default). "quadrature" corresponds the traditional quadrature formulas and calls the function <a href="#">createNIGrid</a> . The tuning parameters are adjusted by <code>crt.bayes.control</code> . Default is set to "cubature".
<code>sens_method</code>	Similar to <code>crt_method</code> , but for approximating the integrals in the sensitivity (derivative) function. The tuning parameters are adjusted by <code>sens.bayes.control</code> . Default is set to "cubature".
<code>plot_3d</code>	Which package should be used to plot the sensitivity (derivative) function for two-dimensional design space. Defaults to "lattice".
<code>plot_sens</code>	Plot the sensitivity (derivative) function? Defaults to TRUE.
<code>npar</code>	Number of model parameters. Used when <code>fimfunc</code> is given instead of <code>formula</code> to specify the number of model parameters. If not specified correctly, the sensitivity (derivative) plot may be shifted below the y-axis. When NULL (default), it will be set to <code>length(parvars)</code> or <code>prior\$npar</code> when <code>missing(formula)</code> .
<code>calculate_criterion</code>	Calculate the optimality criterion? See 'Details' of <a href="#">sensminimax</a> .
<code>silent</code>	Do not print anything? Defaults to FALSE.

## Details

Depending on the complexity of the problem at hand, sometimes, the CPU time can be considerably reduced by choosing a set of less conservative values for the tuning parameters `tol` and `maxEval` in the function [sens.bayes.control](#) when `sens_method = "cubature"`. Similarly, this applies when `sens_method = "quadrature"`. In general, if the CPU time matters, the user should find an appropriate speed-accuracy trade-off for her/his own problem. See 'Examples' for more details.

**Note**

The default values of the tuning parameters in `sens.bayes.control` are set in a way that having accurate plots for the sensitivity (derivative) function and calculating the ELB to a high precision to be the primary goals, although the process may take too long. The user should choose a set of less conservative values via the argument `sens.bayes.control` when the CPU-time is too long or matters.

**See Also**

[bayescomp](#)

**Examples**

```
#####
# Verifying the DP-optimality of a design
# The logistic model with two predictors
#####

# The design points and corresponding weights are as follows:
# Point1    Point2    Point3    Point4    Point5    Point6    Point7
# 0.07410  -0.31953   -1.00000    1.00000   -1.00000    1.00000    0.30193
# -1.00000  1.00000   -1.00000    1.00000   -0.08251   -1.00000    1.00000
# Weight1   Weight2   Weight3   Weight4   Weight5   Weight6   Weight7
# 0.020     0.275     0.224     0.131     0.092     0.156     0.103

# It should be given to the function as two separate vectors:
x1 <- c(0.07409639, -0.3195265, -1, 1, -1, 1, 0.3019317, -1, 1, -1, 1, -0.08251169, -1, 1)
w1 <- c(0.01992863, 0.2745394, 0.2236575, 0.1312331, 0.09161503, 0.1561454, 0.1028811)

p <- c(1, -2, 1, -1)

## Not run:
sensbayescomp(formula = ~exp(b0+b1*x1+b2*x2+b3*x1*x2)/(1+exp(b0+b1*x1+b2*x2+b3*x1*x2)),
  predvars = c("x1", "x2"),
  parvars = c("b0", "b1", "b2", "b3"),
  family = binomial(),
  x = x1, w = w1,
  lx = c(-1, -1), ux = c(1, 1),
  prior = uniform(p - 1.5, p + 1.5),
  prob = ~1-1/(1+exp(b0 + b1 * x1 + b2 * x2 + b3 * x1 * x2)),
  alpha = .5, plot_3d = "rgl",
  sens.bayes.control = list(cubature = list(tol = 1e-3, maxEval = 1000)))

## End(Not run)
```

## Description

It plots the sensitivity (derivative) function of the locally D-optimal criterion at a given approximate (continuous) design and also calculates its efficiency lower bound (ELB) with respect to the optimality criterion. For an approximate (continuous) design, when the design space is one or two-dimensional, the user can visually verify the optimality of the design by observing the sensitivity plot. Furthermore, the proximity of the design to the optimal design can be measured by the ELB without knowing the latter. See, for more details, Masoudi et al. (2017).

## Usage

```
senslocally(formula, predvars, parvars, family = gaussian(), x, w, lx,
            ux, inipars, fimfunc = NULL, sens.minimax.control = list(),
            calculate_criterion = TRUE, plot_3d = c("lattice", "rgl"),
            plot_sens = TRUE, npar = length(inipars), silent = FALSE)
```

## Arguments

formula	A nonlinear model <a href="#">formula</a> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <a href="#">formula</a> if necessary.
predvars	A vector of characters. Denotes the predictors in the <a href="#">formula</a> .
parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .
family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <code>gaussian()</code> is applied.
x	Vector of the design (support) points. See 'Details' of <a href="#">sensminimax</a> for models with more than one predictors.
w	Vector of the corresponding design weights for x.
lx	Vector of lower bounds for the predictors. Should be in the same order as predvars.
ux	Vector of upper bounds for the predictors. Should be in the same order as predvars.
inipars	A vector of initial estimates for the unknown parameters. It must match parvars or the argument param of the function fimfunc, when provided.
fimfunc	A function. Returns the FIM as a matrix. Required when formula is missing. See 'Details' of <a href="#">minimax</a> .

<code>sens.minimax.control</code>	Control parameters to verify the general equivalence theorem. For details, see the function <a href="#">sens.minimax.control</a> .
<code>calculate_criterion</code>	Calculate the optimality criterion? See 'Details' of <a href="#">sensminimax</a> .
<code>plot_3d</code>	Which package should be used to plot the sensitivity (derivative) function for models with two predictors. Either "rgl" or "lattice" (default).
<code>plot_sens</code>	Plot the sensitivity (derivative) function? Defaults to TRUE.
<code>npar</code>	Number of model parameters. Used when <code>fimfunc</code> is given instead of formula to specify the number of model parameters. If not given, the sensitivity plot may be shifted below the y-axis. When NULL, it is set to <code>length(inipars)</code> .
<code>silent</code>	Do not print anything? Defaults to FALSE.

### Details

Let  $\theta_0$  denotes the vector of initial estimates for the unknown parameters. A design  $\xi^*$  is locally D-optimal among all designs on  $\chi$  if and only if the following inequality holds for all  $\mathbf{x} \in \chi$

$$c(\mathbf{x}, \xi^*, \theta_0) = \text{tr}M^{-1}(\xi^*, \theta_0)I(\mathbf{x}, \theta_0) - p \leq 0,$$

with equality at all support points of  $\xi^*$ . Here,  $p$  is the number of model parameters.  $c(\mathbf{x}, \xi^*, \theta_0)$  is called **sensitivity** or **derivative** function.

ELB is a measure of proximity of a design to the optimal design without knowing the latter. Given a design, let  $\epsilon$  be the global maximum of the sensitivity (derivative) function over  $x \in \chi$ . ELB is given by

$$ELB = p/(p + \epsilon),$$

where  $p$  is the number of model parameters. Obviously, calculating ELB requires finding  $\epsilon$  and another optimization problem to be solved. The tuning parameters of this optimization can be regulated via the argument [sens.minimax.control](#). See, for more details, Masoudi et al. (2017).

### Value

an object of class `sensminimax` that is a list with the following elements:

`type` Argument type that is required for print methods.

`optima` A matrix that stores all the local optima over the parameter space. The cost (criterion) values are stored in a column named `Criterion_Value`. The last column (`Answering_Set`) shows if the optimum belongs to the answering set (1) or not (0). See 'Details' of [sens.minimax.control](#). Only applicable for minimax or standardized maximin designs.

`mu` Probability measure on the answering set. Corresponds to the rows of `optima` for which the associated row in column `Answering_Set` is equal to 1. Only applicable for minimax or standardized maximin designs.

`max_deriv` Global maximum of the sensitivity (derivative) function ( $\epsilon$  in 'Details').

`ELB` D-efficiency lower bound. Can not be larger than 1. If negative, see 'Note' in [sensminimax](#) or [sens.minimax.control](#).

`merge_tol` Merging tolerance to create the answering set from the set of all local optima. See 'Details' in `sens.minimax.control`. Only applicable for minimax or standardized maximin designs.

`crtval` Criterion value. Compare it with the column `Crtiterion_Value` in `optima` for minimax and standardized maximin designs.

`time` Used CPU time (rough approximation).

## Note

Theoretically, ELB can not be larger than 1. But if so, it may have one of the following reasons:

- `max_deriv` is not a GLOBAL maximum. Please increase the value of the parameter `maxeval` in `sens.minimax.control` to find the global maximum.
- The sensitivity function is shifted below the y-axis because the number of model parameters has not been specified correctly (less value given). Please specify the correct number of model parameters via the argument `npar`.

## References

Masoudi E, Holling H, Wong W.K. (2017). Application of Imperialist Competitive Algorithm to Find Minimax and Standardized Maximin Optimal Designs. *Computational Statistics and Data Analysis*, 113, 330-345.

## Examples

```
#####
# Exponential growth model
#####
# Verifying optimailty of a locally D-optimal design
senslocally(formula = ~a + exp(-b*x),
             predvars = "x", parvars = c("a", "b"),
             x = c(.1, 1), w = c(.5, .5),
             lx = 0, ux = 1, inipars = c(1, 10))

#####
# A model with two predictors
#####
x0 <- c(30, 3.861406, 30, 4.600633, 0, 0, 5.111376, 4.168798)
w0 <- rep(.25, 4)
senslocally(formula = ~ V*S/(Km * (1 + I/Kic)+ S * (1 + I/Kiu)),
             predvars = c("S", "I"),
             parvars = c("V", "Km", "Kic", "Kiu"),
             x = x0, w = w0,
             lx = c(0, 0), ux = c(30, 60),
             inipars = c(1.5, 5.2, 3.4, 5.6))

## Not run:
# using package rgl for 3d plot:
res<- senslocally(formula = ~ V*S/(Km * (1 + I/Kic)+ S * (1 + I/Kiu)),
                  predvars = c("S", "I"),
```

```

parvars = c("V", "Km", "Kic", "Kiu"),
x = x0, w = w0,
lx = c(0, 0), ux = c(30, 60),
inipars = c(1.5, 5.2, 3.4, 5.6),
plot_3d = "rgl")

```

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

---

senslocallycomp

*Verifying Optimality of The Locally DP-optimal Designs*


---

## Description

This function plot the sensitivity (derivative) function given an approximate (continuous) design and calculate the efficiency lower bound (ELB) for locally DP-optimal designs. Let  $x$  belongs to  $\chi$  that denotes the design space. Based on the general equivalence theorem, generally, a design  $\xi^*$  is optimal if and only if the value of its sensitivity (derivative) function be non-positive for all  $x$  in  $\chi$  and it only reaches zero when  $x$  belong to the support of  $\xi^*$  (be equal to one of the design point). Therefore, the user can look at the sensitivity plot and the ELB to decide whether the design is optimal or close enough to the true optimal design.

## Usage

```

senslocallycomp(formula, predvars, parvars, alpha, prob,
  family = gaussian(), x, w, lx, ux, inipars, fimfunc = NULL,
  sens.minimax.control = list(), calculate_criterion = TRUE,
  plot_3d = c("lattice", "rgl"), plot_sens = TRUE,
  npar = length(inipars), silent = FALSE)

```

## Arguments

formula	A nonlinear model <a href="#">formula</a> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <a href="#">formula</a> if necessary.
predvars	A vector of characters. Denotes the predictors in the <a href="#">formula</a> .
parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .
alpha	A value between 0 and 1. Compound or combined DP-criterion is the product of the efficiencies of a design with respect to D- and average P- optimality, weighted by alpha.
prob	Either formula or a function. When function, its argument are x and param, and they are the same as the arguments in fimfunc. prob as a function takes the design points and vector of parameters and returns the probability of success at each design points. See 'Examples'.

family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <code>gaussian()</code> is applied.
x	Vector of the design (support) points. See 'Details' of <a href="#">sensminimax</a> for models with more than one predictors.
w	Vector of the corresponding design weights for x.
lx	Vector of lower bounds for the predictors. Should be in the same order as <code>predvars</code> .
ux	Vector of upper bounds for the predictors. Should be in the same order as <code>predvars</code> .
inipars	Vector of initial estimates for the unknown parameters. It must match <code>parvars</code> or argument <code>param</code> of the function provided in <code>fimfunc</code> .
fimfunc	A function. Returns the FIM as a matrix. Required when <code>formula</code> is missing. See 'Details' of <a href="#">minimax</a> .
<code>sens.minimax.control</code>	Control parameters to verify the general equivalence theorem. For details, see the function <a href="#">sens.minimax.control</a> .
<code>calculate_criterion</code>	Calculate the optimality criterion? See 'Details' of <a href="#">sensminimax</a> .
<code>plot_3d</code>	Which package should be used to plot the sensitivity (derivative) function for models with two predictors. Either <code>"rgl"</code> or <code>"lattice"</code> (default).
<code>plot_sens</code>	Plot the sensitivity (derivative) function? Defaults to TRUE.
<code>npar</code>	Number of model parameters. Used when <code>fimfunc</code> is given instead of <code>formula</code> to specify the number of model parameters. If not given, the sensitivity plot may be shifted below the y-axis. When NULL, it is set to <code>length(inipars)</code> .
<code>silent</code>	Do not print anything? Defaults to FALSE.

## Value

an object of class `sensminimax` that is a list with the following elements:

`type` Argument type that is required for print methods.

`optima` A matrix that stores all the local optima over the parameter space. The cost (criterion) values are stored in a column named `Criterion_Value`. The last column (`Answering_Set`) shows if the optimum belongs to the answering set (1) or not (0). See 'Details' of [sens.minimax.control](#). Only applicable for minimax or standardized maximin designs.

`mu` Probability measure on the answering set. Corresponds to the rows of `optima` for which the associated row in column `Answering_Set` is equal to 1. Only applicable for minimax or standardized maximin designs.

`max_deriv` Global maximum of the sensitivity (derivative) function ( $\epsilon$  in 'Details').

`ELB` D-efficiency lower bound. Can not be larger than 1. If negative, see 'Note' in [sensminimax](#) or [sens.minimax.control](#).

merge\_tol Merging tolerance to create the answering set from the set of all local optima. See 'Details' in `sens.minimax.control`. Only applicable for minimax or standardized maximin designs.

crtval Criterion value. Compare it with the column `Crtiterion_Value` in `optima` for minimax and standardized maximin designs.

time Used CPU time (rough approximation).

## References

McGree, J. M., Eccleston, J. A., and Duffull, S. B. (2008). Compound optimal design criteria for nonlinear models. *Journal of Biopharmaceutical Statistics*, 18(4), 646-661.

## Examples

```
p <- c(1, -2, 1, -1)
prior4.4 <- uniform(p -1.5, p + 1.5)
formula4.4 <- ~exp(b0+b1*x1+b2*x2+b3*x1*x2)/(1+exp(b0+b1*x1+b2*x2+b3*x1*x2))
prob4.4 <- ~1-1/(1+exp(b0 + b1 * x1 + b2 * x2 + b3 * x1 * x2))
predvars4.4 <- c("x1", "x2")
parvars4.4 <- c("b0", "b1", "b2", "b3")
lb <- c(-1, -1)
ub <- c(1, 1)

## That is the optimal design when alpha = .25, see ?locallycomp on how to find it
xopt <- c(-1, -0.389, 1, 0.802, -1, 1, -1, 1)
wopt <- c(0.198, 0.618, 0.084, 0.1)

# We want to verify the optimality of the optimal design by the general equivalence theorem.

senslocallycomp(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4,
                family = binomial(), prob = prob4.4, lx = lb, ux = ub,
                alpha = .25, inipars = p, x = xopt, w = wopt)

## Not run:
# is this design also optimal when alpha = .3

senslocallycomp(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4,
                family = binomial(), prob = prob4.4, lx = lb, ux = ub,
                alpha = .3, inipars = p, x = xopt, w = wopt)

# when alpha = .3
senslocallycomp(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4,
                family = binomial(), prob = prob4.4, lx = lb, ux = ub,
                alpha = .5, inipars = p, x = xopt, w = wopt)

# when alpha = .8
senslocallycomp(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4,
                family = binomial(), prob = prob4.4, lx = lb, ux = ub,
                alpha = .8, inipars = p, x = xopt, w = wopt)

# when alpha = .9
```

```

senslocallycomp(formula = formula4.4, predvars = predvars4.4, parvars = parvars4.4,
                family = binomial(), prob = prob4.4, lx = lb, ux = ub,
                alpha = .9, inipars = p, x = xopt, w = wopt)

## As can be seen, the design loses efficiency as alpha increases.

## End(Not run)

```

---

sensminimax	<i>Verifying Optimality of The Minimax and Standardized maximin D-optimal Designs</i>
-------------	---

---

## Description

It plots the sensitivity (derivative) function of the minimax or standardized maximin D-optimal criterion at a given approximate (continuous) design and also calculates its efficiency lower bound (ELB) with respect to the optimality criterion. For an approximate (continuous) design, when the design space is one or two-dimensional, the user can visually verify the optimality of the design by observing the sensitivity plot. Furthermore, the proximity of the design to the optimal design can be measured by the ELB without knowing the latter. See, for more details, Masoudi et al. (2017).

## Usage

```

sensminimax(formula, predvars, parvars, family = gaussian(), x, w, lx,
            ux, lp, up, fimfunc = NULL, standardized = FALSE, localdes = NULL,
            sens.minimax.control = list(), calculate_criterion = TRUE,
            crt.minimax.control = list(), plot_3d = c("lattice", "rgl"),
            plot_sens = TRUE, npar = length(lp), silent = FALSE)

```

## Arguments

formula	A nonlinear model <a href="#">formula</a> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <a href="#">formula</a> if necessary.
predvars	A vector of characters. Denotes the predictors in the <a href="#">formula</a> .
parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .
family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <code>gaussian()</code> is applied.
x	Vector of the design (support) points. See 'Details' of <a href="#">sensminimax</a> for models with more than one predictors.
w	Vector of the corresponding design weights for x.

lx	Vector of lower bounds for the predictors. Should be in the same order as predvars.
ux	Vector of upper bounds for the predictors. Should be in the same order as predvars.
lp	Vector of lower bounds for the model parameters. Should be in the same order as parvars or param in the argument fimfunc.
up	Vector of upper bounds for the model parameters. Should be in the same order as parvars or param in the argument fimfunc. When a parameter is known (has a fixed value), its associated lower and upper bound values in lp and up must be set equal.
fimfunc	A function. Returns the FIM as a matrix. Required when formula is missing. See 'Details' of <a href="#">minimax</a> .
standardized	Maximin standardized design? When standardized = TRUE, the argument localdes must be given. Defaults to FALSE. See 'Details' of <a href="#">minimax</a> .
localdes	A function that takes the parameter values as inputs and returns the design points and weights of the locally optimal design. Required when standardized = "TRUE". See 'Details' of <a href="#">minimax</a> .
sens.minimax.control	Control parameters to verify the general equivalence theorem. For details, see the function <a href="#">sens.minimax.control</a> .
calculate_criterion	Calculate the optimality criterion? See 'Details' of <a href="#">sensminimax</a> .
crt.minimax.control	Control parameters to calculate the value of the minimax or standardized maximin optimality criterion over the continuous parameter space. Only applicable when calculate_criterion = TRUE. For more details, see <a href="#">crt.minimax.control</a> .
plot_3d	Which package should be used to plot the sensitivity (derivative) function for models with two predictors. Either "rgl" or "lattice" (default).
plot_sens	Plot the sensitivity (derivative) function? Defaults to TRUE.
npar	Number of model parameters. Used when fimfunc is given instead of formula to specify the number of model parameters. If not specified truly, the sensitivity (derivative) plot may be shifted below the y-axis. When NULL (default), it is set to length(lp).
silent	Do not print anything? Defaults to FALSE.

## Details

Let the unknown parameters belong to  $\Theta$ . A design  $\xi^*$  is minimax D-optimal among all designs on  $\chi$  if and only if there exists a probability measure  $\mu^*$  on

$$A(\xi^*) = \left\{ \nu \in \Theta \mid -\log|M(\xi^*, \nu)| = \max_{\theta \in \Theta} -\log|M(\xi^*, \theta)| \right\},$$

such that the following inequality holds for all  $\mathbf{x} \in \chi$

$$c(\mathbf{x}, \mu^*, \xi^*) = \int_{A(\xi^*)} \text{tr} M^{-1}(\xi^*, \nu) I(\mathbf{x}, \nu) \mu^* d(\nu) - p \leq 0,$$

with equality at all support points of  $\xi^*$ . Here,  $p$  is the number of model parameters.  $c(\mathbf{x}, \mu^*, \xi^*)$  is called **sensitivity** or **derivative** function. The set  $A(\xi^*)$  is sometimes called **answering set** of  $\xi^*$  and the measure  $\mu^*$  is a sub-gradient of the non-differentiable criterion evaluated at  $M(\xi^*, \nu)$ . For the standardized maximin D-optimal designs, the answering set  $N(\xi^*)$  is

$$N(\xi^*) = \left\{ \nu \in \Theta \mid \text{eff}_D(\xi^*, \nu) = \min_{\theta \in \Theta} \text{eff}_D(\xi^*, \theta) \right\}.$$

where  $\text{eff}_D(\xi, \theta) = \left( \frac{|M(\xi, \theta)|}{|M(\xi_\theta, \theta)|} \right)^{\frac{1}{p}}$  and  $\xi_\theta$  is the locally D-optimal design with respect to  $\theta$ . See 'Details' of [sens.minimax.control](#) on how we find the answering set.

The argument  $\mathbf{x}$  is the vector of design points. For design points with more than one dimension (the models with more than one predictors), it is a concatenation of the design points, but **dimension-wise**. For example, let the model has three predictors  $(I, S, Z)$ . Then, a two-point optimal design has the following points:  $\{\text{point1} = (I_1, S_1, Z_1), \text{point2} = (I_2, S_2, Z_2)\}$ . Then, the argument  $\mathbf{x}$  is equal to  $\mathbf{x} = c(I1, I2, S1, S2, Z1, Z2)$ .

ELB is a measure of proximity of a design to the optimal design without knowing the latter. Given a design, let  $\epsilon$  be the global maximum of the sensitivity (derivative) function with respect  $x$  where  $x \in \chi$ . ELB is given by

$$ELB = p/(p + \epsilon),$$

where  $p$  is the number of model parameters. Obviously, calculating ELB requires finding  $\epsilon$  and another optimization problem to be solved. The tuning parameters of this optimization can be regulated via the argument [sens.minimax.control](#). See, for more details, Masoudi et al. (2017).

The criterion value for the minimax D-optimal design is (global maximum over  $\Theta$ )

$$\max_{\theta \in \Theta} -\log |M(\xi, \theta)|;$$

for the standardized maximin D-optimal design is (global minimum over  $\Theta$ )

$$\inf_{\theta \in \Theta} \left[ \left( \frac{|M(\xi, \theta)|}{|M(\xi_\theta, \theta)|} \right)^{\frac{1}{p}} \right].$$

This function confirms the optimality assuming only a continuous parameter space  $\Theta$ .

## Value

an object of class `sensminimax` that is a list with the following elements:

`type` Argument type that is required for print methods.

`optima` A matrix that stores all the local optima over the parameter space. The cost (criterion) values are stored in a column named `Criterion_Value`. The last column (`Answering_Set`) shows if the optimum belongs to the answering set (1) or not (0). See 'Details' of [sens.minimax.control](#). Only applicable for minimax or standardized maximin designs.

`mu` Probability measure on the answering set. Corresponds to the rows of `optima` for which the associated row in column `Answering_Set` is equal to 1. Only applicable for minimax or standardized maximin designs.

`max_deriv` Global maximum of the sensitivity (derivative) function ( $\epsilon$  in 'Details').

ELB D-efficiency lower bound. Can not be larger than 1. If negative, see 'Note' in [sensminimax](#) or [sens.minimax.control](#).

merge\_tol Merging tolerance to create the answering set from the set of all local optima. See 'Details' in [sens.minimax.control](#). Only applicable for minimax or standardized maximin designs.

crtval Criterion value. Compare it with the column Crtiterion\_Value in optima for minimax and standardized maximin designs.

time Used CPU time (rough approximation).

### Note

Theoretically, ELB can not be larger than 1. But if so, it may have one of the following reasons:

- `max_deriv` is not a GLOBAL maximum. Please increase the value of the parameter `maxeval` in [sens.minimax.control](#) to find the global maximum.
- The sensitivity function is shifted below the y-axis because the number of model parameters has not been specified correctly (less value given). Please specify the correct number of model parameters via argument `npar`.

Please increase the value of the parameter `n_seg` in [sens.minimax.control](#) for models with larger number of parameters or large parameter space to find the true answering set for minimax and standardized maximin designs. See [sens.minimax.control](#) for more details.

### References

Masoudi E, Holling H, Wong W.K. (2017). Application of Imperialist Competitive Algorithm to Find Minimax and Standardized Maximin Optimal Designs. *Computational Statistics and Data Analysis*, 113, 330-345.

### Examples

```
#####
# Power logistic model
#####
# verifying the minimax D-optimality of a design with points x0 and weights w0
x0 <- c(-4.5515, 0.2130, 2.8075)
w0 <- c(0.4100, 0.3723, 0.2177)
# Power logistic model when s = .2
sensminimax(formula = ~ (1/(1 + exp(-b * (x-a))))^2,
             predvars = "x",
             parvars = c("a", "b"),
             family = binomial(),
             x = x0, w = w0,
             lx = -5, ux = 5,
             lp = c(0, 1), up = c(3, 1.5))

#####
# A model with two predictors
#####
```

```

# Verifying the minimax D-optimality of a design for a model with two predictors
# The model is the mixed inhibition model.
# X0 is the vector of four design points that are:
# (3.4614, 0) (4.2801, 3.1426) (30, 0) (30, 4.0373)
x0 <- c(3.4614, 4.2801, 30, 30, 0, 3.1426, 0, 4.0373)
w0 <- rep(1/4, 4)
sensminimax(formula = ~ V*S/(Km * (1 + I/Kic)+ S * (1 + I/Kiu)),
             predvars = c("S", "I"),
             parvars = c("V", "Km", "Kic", "Kiu"),
             family = "gaussian",
             x = x0, w = w0,
             lx = c(0, 0), ux = c(30, 60),
             lp = c(1, 4, 2, 4), up = c(1, 5, 3, 5))

#####
# Standardized maximin D-optimal designs
#####
# Verifying the standardized maximin D-optimality of a design for
# the loglinear model
# First we should define the function for 'localdes' argument
# The function LDOD takes the parameters and returns the points and
# weights of the locally D-optimal design
LDOD <- function(theta0, theta1, theta2){
  ## param is the vector of theta = (theta0, theta1, theta2)
  lx <- 0 # lower bound of the design space
  ux <- 150 # upper bound of the design space
  param <- c()
  param[1] <- theta0
  param[2] <- theta1
  param[3] <- theta2
  xstar <- (ux+param[3]) * (lx + param[3]) *
    (log(ux + param[3]) - log(lx + param[3]))/(ux - lx) - param[3]
  return(list(x = c(lx, xstar, ux) , w = rep(1/3, 3)))
}
x0 <- c(0, 4.2494, 17.0324, 149.9090)
w0 <- c(0.3204, 0.1207, 0.2293, 0.3296)
## Not run:
sensminimax(formula = ~theta0 + theta1* log(x + theta2),
             predvars = c("x"),
             parvars = c("theta0", "theta1", "theta2"),
             x = x0, w = w0,
             lx = 0, ux = 150,
             lp = c(2, 2, 1), up = c(2, 2, 15),
             localdes = LDOD,
             standardized = TRUE,
             sens.minimax.control = list(answering.set = list(n_seg = 10)))

## End(Not run)
#####
# Not necessary!
# The rest of the examples here are only for professional uses.
#####
# Imagine you have written your own FIM, say in Rcpp that is faster than

```

```

# the FIM created by the formula interface here.

#####
# Power logistic model
#####
# For example, th cpp FIM function for the power logistic model is named:
FIM_power_logistic
args(FIM_power_logistic)
# The arguments do not match the standard of the argument 'fimfunc'
# in 'sensminimax'
# So we reparameterize it:
myfim1 <- function(x, w, param)
  FIM_power_logistic(x = x, w = w, param =param, s = .2)

args(myfim1)
## Not run:
# Verify minimax D-optimality of a design
sensminimax(fimfunc = myfim1,
            x = c(-4.5515, 0.2130, 2.8075),
            w = c(0.4100, 0.3723, 0.2177),
            lx = -5, ux = 5,
            lp = c(0, 1), up = c(3, 1.5))

## End(Not run)
#####
# A model with two predictors
#####
# An example of a model with two-predictors: mixed inhibition model
# Fisher information matrix:
FIM_mixed_inhibition
args(FIM_mixed_inhibition)

# We should first reparameterize the FIM to match the standard of the
# argument 'fimfunc'
myfim2 <- function(x, w, param){
  npoint <- length(x)/2
  S <- x[1:npoint]
  I <- x[(npoint+1):(npoint*2)]
  out <- FIM_mixed_inhibition(S = S, I = I, w = w, param = param)
  return(out)
}
args(myfim2)
## Not run:
# Verifyng minimax D-optimality of a design
sensminimax(fimfunc = myfim2,
            x = c(3.4614, 4.2801, 30, 30, 0, 3.1426, 0, 4.0373),
            w = rep(1/4, 4),
            lx = c(0, 0), ux = c(30, 60),
            lp = c(1, 4, 2, 4), up = c(1, 5, 3, 5))

## End(Not run)

#####

```

```

# Standardized maximin D-optimal designs
#####
# An example of a user-written FIM function:
help(FIM_loglin)
# An example of verifying standardized maximin D-optimality for a design
# Look how we re-define the function LDOD above
LDOD2 <- function(param){
  ## param is the vector of theta = (theta0, theta1, theta2)
  lx <- 0 # lower bound of the design space
  ux <- 150 # upper bound of the design space
  xstar <- (ux + param[3]) * (lx + param[3]) *
    (log(ux + param[3]) - log(lx + param[3]))/(ux - lx) - param[3]
  return(list(x = c(lx, xstar, ux) , w = rep(1/3, 3)))
}

args(LDOD2)

sensminimax(fimfunc = FIM_loglin,
            x = x0,
            w = w0,
            lx = 0, ux = 150,
            lp = c(2, 2, 1), up = c(2, 2, 15),
            localdes = LDOD2,
            standardized = TRUE)

```

---

sensmultiple

*Verifying Optimality of The Multiple Objective Designs for The 4-Parameter Hill Model*


---

## Description

This function uses general equivalence theorem to verify the optimality of a multiple objective optimal design found for the 4-Parameter Hill model and the 4-parameter logistic model. For more details, See Hyun and Wong (2015).

## Usage

```

sensmultiple(dose, w, minDose, maxDose, inipars, lambda, delta,
            Hill_par = FALSE, sens.minimax.control = list(),
            calculate_criterion = TRUE, plot_sens = TRUE,
            tol = sqrt(.Machine$double.xmin), silent = FALSE)

```

## Arguments

**dose** A vector of design points. It is either dose values or logarithm of dose values when `Hill_par = TRUE`.

w	A vector of design weights.
minDose	Minimum dose $D$ . For the 4-parameter logistic model, i.e. when Hill_par = FALSE, it is the minimum of $\log(D)$ .
maxDose	Maximum dose $D$ . For the 4-parameter logistic model, i.e. when Hill_par = FALSE, it is the maximum of $\log(D)$ .
inipars	A vector of initial estimates for the vector of parameters $(a, b, c, d)$ . For the 4-parameter logistic model, i.e. when Hill_par = FALSE, it is a vector of initial estimates for $(\theta_1, \theta_2, \theta_3, \theta_4)$ .
lambda	A vector of relative importance of each of the three criteria, i.e. $\lambda = (\lambda_1, \lambda_2, \lambda_3)$ . Here $0 < \lambda_i < 1$ and $\sum \lambda_i = 1$ .
delta	Predetermined meaningful value of the minimum effective dose MED. When $\delta < 0$ , then $\theta_2 > 0$ or when $\delta > 0$ , then $\theta_2 < 0$ .
Hill_par	Hill model parameterization? Defaults to TRUE.
sens.minimax.control	Control parameters to verify the general equivalence theorem. For details, see the function <a href="#">sens.minimax.control</a> .
calculate_criterion	Calculate the criterion? Defaults to TRUE.
plot_sens	Plot the sensitivity (derivative) function? Defaults to TRUE.
tol	Tolerance for finding the general inverse of the Fisher information matrix. Defaults to <code>.Machine\$double.xmin</code> .
silent	Do not print anything? Defaults to FALSE.

## Details

ELB is a measure of proximity of a design to the optimal design without knowing the latter. Given a design, let  $\epsilon$  be the global maximum of the sensitivity (derivative) function over  $x \in \chi$ . ELB is given by

$$ELB = p/(p + \epsilon),$$

where  $p$  is the number of model parameters. Obviously, calculating ELB requires finding  $\epsilon$  and another optimization problem to be solved. The tuning parameters of this optimization can be regulated via the argument [sens.minimax.control](#). See, for more details, Masoudi et al. (2017).

## Value

an object of class `sensminimax` that is a list with the following elements:

`type` Argument type that is required for print methods.

`optima` A matrix that stores all the local optima over the parameter space. The cost (criterion) values are stored in a column named `Criterion_Value`. The last column (`Answering_Set`) shows if the optimum belongs to the answering set (1) or not (0). See 'Details' of [sens.minimax.control](#). Only applicable for minimax or standardized maximin designs.

`mu` Probability measure on the answering set. Corresponds to the rows of `optima` for which the associated row in column `Answering_Set` is equal to 1. Only applicable for minimax or standardized maximin designs.

`max_deriv` Global maximum of the sensitivity (derivative) function ( $\epsilon$  in 'Details').

`ELB` D-efficiency lower bound. Can not be larger than 1. If negative, see 'Note' in [sensminimax](#) or [sens.minimax.control](#).

`merge_tol` Merging tolerance to create the answering set from the set of all local optima. See 'Details' in [sens.minimax.control](#). Only applicable for minimax or standardized maximin designs.

`crtval` Criterion value. Compare it with the column `Crtiterion_Value` in `optima` for minimax and standardized maximin designs.

`time` Used CPU time (rough approximation).

### Note

DO NOT use this function to verify c-optimal designs for estimating 'MED' or 'ED50' (verifying single objective optimal designs) because the results may be unstable. The reason is that for the c-optimal criterion the generalized inverse of the Fisher information matrix is not stable and depends on the tolerance value (`tol`).

Theoretically, ELB can not be larger than 1. But if so, it may have one of the following reasons:

- `max_deriv` is not a GLOBAL maximum. Please increase the value of the parameter `maxeval` in [sens.minimax.control](#) to find the global maximum.
- The sensitivity function is shifted below the y-axis because the number of model parameters has not been specified correctly (less value given). Please specify the correct number of model parameters via argument `npar`.

### References

Hyun, S. W., and Wong, W. K. (2015). Multiple-Objective Optimal Designs for Studying the Dose Response Function and Interesting Dose Levels. *The international journal of biostatistics*, 11(2), 253-271.

### See Also

[multiple](#)

### Examples

```
#####
# Verifying optimality of a design for the 4-parameter Hill model
#####

## initial estiamtes for the parameters of the Hill model
a <- 0.008949 # ED50
b <- -1.79 # Hill constant
c <- 0.137 # lower limit
d <- 1.7 # upper limit
# D belongs to c(.001, 1000) ## dose in mg
## Hill parameters are c(a, b, c, d)
# dose, minDose and maxDose vector in mg scale
```

```

sensmultiple (dose = c(0.001, 0.009426562, 0.01973041, 999.9974),
              w = c(0.4806477, 0.40815, 0.06114173, 0.05006055),
              minDose = .001, maxDose = 1000,
              Hill_par = TRUE,
              inipars = c(a, b, c, d),
              lambda = c(0.05, 0.05, .90),
              delta = -1)

```

---

sensrobust

*Verifying Optimality of The Robust Designs*


---

### Description

It plots the sensitivity (derivative) function of the robust criterion at a given approximate (continuous) design and also calculates its efficiency lower bound (ELB) with respect to the optimality criterion. For an approximate (continuous) design, when the design space is one or two-dimensional, the user can visually verify the optimality of the design by observing the sensitivity plot. Furthermore, the proximity of the design to the optimal design can be measured by the ELB without knowing the latter. See, for more details, Masoudi et al. (2017).

### Usage

```

sensrobust(formula, predvars, parvars, family = gaussian(), x, w, lx, ux,
           prob, parset, fimfunc = NULL, sens.minimax.control = list(),
           calculate_criterion = TRUE, plot_3d = c("lattice", "rgl"),
           plot_sens = TRUE, npar = dim(parset)[2], silent = FALSE)

```

### Arguments

formula	A nonlinear model <a href="#">formula</a> . A symbolic description of the model consists of predictors and the unknown model parameters. Will be coerced to a <a href="#">formula</a> if necessary.
predvars	A vector of characters. Denotes the predictors in the <a href="#">formula</a> .
parvars	A vector of characters. Denotes the unknown parameters in the <a href="#">formula</a> .
family	A description of the response distribution and the link function to be used in the model. This can be a family function, a call to a family function or a character string naming the family. Every family function has a link argument allowing to specify the link function to be applied on the response variable. If not specified, default links are used. For details see <a href="#">family</a> . By default, a linear gaussian model <code>gaussian()</code> is applied.

<code>x</code>	Vector of the design (support) points. See 'Details' of <a href="#">sensminimax</a> for models with more than one predictors.
<code>w</code>	Vector of the corresponding design weights for <code>x</code> .
<code>lx</code>	Vector of lower bounds for the predictors. Should be in the same order as <code>predvars</code> .
<code>ux</code>	Vector of upper bounds for the predictors. Should be in the same order as <code>predvars</code> .
<code>prob</code>	A vector of the probability measure $\pi$ associated with each row of <code>parset</code> .
<code>parset</code>	A matrix that provides the vector of initial estimates for the model parameters, i.e. support of $\pi$ . Every row is one vector ( <code>nrow(parset) == length(prob)</code> ). See 'Details'.
<code>fimfunc</code>	A function. Returns the FIM as a matrix. Required when <code>formula</code> is missing. See 'Details' of <a href="#">minimax</a> .
<code>sens.minimax.control</code>	Control parameters to verify the general equivalence theorem. For details, see the function <a href="#">sens.minimax.control</a> .
<code>calculate_criterion</code>	Calculate the optimality criterion? See 'Details' of <a href="#">sensminimax</a> .
<code>plot_3d</code>	Which package should be used to plot the sensitivity (derivative) function for models with two predictors. Either "rgl" or "lattice" (default).
<code>plot_sens</code>	Plot the sensitivity (derivative) function? Defaults to TRUE.
<code>npar</code>	Number of model parameters. Used when <code>fimfunc</code> is given instead of <code>formula</code> to specify the number of model parameters. If not given, the sensitivity plot may be shifted below the y-axis. When NULL, it is set to <code>length(inipars)</code> .
<code>silent</code>	Do not print anything? Defaults to FALSE.

## Details

Let  $\Theta$  be the set initial estimates for the model parameters and  $\pi$  be a probability measure having support in  $\Theta$ . A design  $\xi^*$  is robust with respect to  $\pi$  if the following inequality holds for all  $\mathbf{x} \in \chi$ :

$$c(\mathbf{x}, \pi, \xi^*) = \int_{\pi} \text{tr} M^{-1}(\xi^*, \theta) I(\mathbf{x}, \theta) \pi(\theta) d(\theta) - p \leq 0,$$

with equality at all support points of  $\xi^*$ . Here,  $p$  is the number of model parameters.

ELB is a measure of proximity of a design to the optimal design without knowing the latter. Given a design, let  $\epsilon$  be the global maximum of the sensitivity (derivative) function over  $\mathbf{x} \in \chi$ . ELB is given by

$$ELB = p/(p + \epsilon),$$

where  $p$  is the number of model parameters. Obviously, calculating ELB requires finding  $\epsilon$  and another optimization problem to be solved. The tuning parameters of this optimization can be regulated via the argument [sens.minimax.control](#).

**Value**

an object of class `sensminimax` that is a list with the following elements:

`type` Argument type that is required for print methods.

`optima` A matrix that stores all the local optima over the parameter space. The cost (criterion) values are stored in a column named `Criterion_Value`. The last column (`Answering_Set`) shows if the optimum belongs to the answering set (1) or not (0). See 'Details' of [sens.minimax.control](#). Only applicable for minimax or standardized maximin designs.

`mu` Probability measure on the answering set. Corresponds to the rows of `optima` for which the associated row in column `Answering_Set` is equal to 1. Only applicable for minimax or standardized maximin designs.

`max_deriv` Global maximum of the sensitivity (derivative) function ( $\epsilon$  in 'Details').

`ELB` D-efficiency lower bound. Can not be larger than 1. If negative, see 'Note' in [sensminimax](#) or [sens.minimax.control](#).

`merge_tol` Merging tolerance to create the answering set from the set of all local optima. See 'Details' in [sens.minimax.control](#). Only applicable for minimax or standardized maximin designs.

`crtval` Criterion value. Compare it with the column `Criterion_Value` in `optima` for minimax and standardized maximin designs.

`time` Used CPU time (rough approximation).

**Note**

Theoretically, ELB can not be larger than 1. But if so, it may have one of the following reasons:

- `max_deriv` is not a GLOBAL maximum. Please increase the value of the parameter `maxeval` in [sens.minimax.control](#) to find the global maximum.
- The sensitivity function is shifted below the y-axis because the number of model parameters has not been specified correctly (less value given). Please specify the correct number of model parameters via the argument `npar`.

**See Also**

[bayes](#) [sensbayes](#) [robust](#)

**Examples**

```
# Verifying a robust design for the two-parameter logistic model
sensrobust(formula = ~1/(1 + exp(-b *(x - a))),
  predvars = c("x"),
  parvars = c("a", "b"),
  family = binomial(),
  prob = rep(1/4, 4),
  parset = matrix(c(0.5, 1.5, 0.5, 1.5, 4.0, 4.0, 5.0, 5.0), 4, 2),
  x = c(0.260, 1, 1.739), w = c(0.275, 0.449, 0.275),
  lx = -5, ux = 5)
```

---

skewnormal	<i>Assumes A Multivariate Skewed Normal Prior Distribution for The Model Parameters</i>
------------	---

---

### Description

Creates a multivariate skewed normal prior distribution for the unknown parameters as an object of class `cprior`.

### Usage

```
skewnormal(xi, Omega, alpha, lower, upper)
```

### Arguments

<code>xi</code>	A numeric vector of length $d = \text{length}(\alpha)$ representing the location parameter of the distribution. For more details, see 'Background' in <a href="#">dmsn</a> .
<code>Omega</code>	A symmetric positive-definite matrix of dimension $(d, d)$ . For more details, see 'Background' in <a href="#">dmsn</a> .
<code>alpha</code>	A numeric vector which regulates the slant of the density. For more details, see 'Background' in <a href="#">dmsn</a> .
<code>lower</code>	A vector of lower bounds for the model parameters.
<code>upper</code>	A vector of upper bounds for the model parameters.

### Value

An object of class `cprior` that is a list with the following components:

- `fn`: prior distribution as an R function with argument `param` that is the vector of the unknown parameters. See below.
- `npar`: Number of unknown parameters and is equal to the length of `param`.
- `lower`: Argument `lower`. It has the same length as `param`.
- `upper`: Argument `lower`. It has the same length as `param`.

The list will be passed to the argument `prior` of the function [bayes](#). The order of the argument `param` in `fn` has the same order as the argument `parvars` when the model is specified by a formula. Otherwise, it is equal to the argument `param` in the function `fimfunc`.

### See Also

[bayes](#) [sensbayes](#)

### Examples

```
skewnormal(xi = c(0, 1),
  Omega = matrix(c(1, -0.17, -0.17, .5), nrow = 2),
  alpha = c(1, 0), lower = c(-3, .1), upper = c(3, 2))
```

---

student	<i>Multivariate Student's t Prior Distribution for Model Parameters</i>
---------	---

---

**Description**

Creates the prior distribution for the parameters as an object of class `cprior`.

**Usage**

```
student(mean, S, df, lower, upper)
```

**Arguments**

mean	A vector of length $d = \text{ncol}(S)$ , representing the location parameter (equal to the mean vector when $df > 1$ ). For more details, see 'Arguments' in <a href="#">dmt</a> .
S	A symmetric positive-definite matrix representing the scale matrix of the distribution, such that $S * df / (df - 2)$ is the variance-covariance matrix when $df > 2$ . For more details, see 'Arguments' in <a href="#">dmt</a> .
df	Degrees of freedom; it must be a positive integer. For more details, see 'Arguments' in <a href="#">dmt</a> .
lower	A vector of lower bounds for the model parameters.
upper	A vector of upper bounds for the model parameters.

**Value**

An object of class `cprior` that is a list with the following components:

- `fn`: prior distribution as an R function with argument `param` that is the vector of the unknown parameters. See below.
- `npar`: Number of unknown parameters and is equal to the length of `param`.
- `lower`: Argument `lower`. It has the same length as `param`.
- `upper`: Argument `upper`. It has the same length as `param`.

The list will be passed to the argument `prior` of the function [bayes](#). The order of the argument `param` in `fn` has the same order as the argument `parvars` when the model is specified by a formula. Otherwise, it is equal to the argument `param` in the function `fimfunc`.

**See Also**

[bayes](#) [sensbayes](#)

**Examples**

```
skewnormal(xi = c(0, 1),
  Omega = matrix(c(1, -0.17, -0.17, .5), nrow = 2),
  alpha = c(1, 0), lower = c(-3, .1), upper = c(3, 2))
```

---

uniform	<i>Assume A Multivariate Uniform Prior Distribution for The Model Parameters</i>
---------	--

---

**Description**

Creates independent uniform prior distributions for the unknown model parameters as an object of class `cprior`.

**Usage**

```
uniform(lower, upper)
```

**Arguments**

lower	A vector of lower bounds for the model parameters.
upper	A vector of upper bounds for the model parameters.

**Value**

An object of class `cprior` that is a list with the following components:

- `fn`: prior distribution as an R function with argument `param` that is the vector of the unknown parameters. See below.
- `npar`: Number of unknown parameters and is equal to the length of `param`.
- `lower`: Argument `lower`. It has the same length as `param`.
- `upper`: Argument `lower`. It has the same length as `param`.

The list will be passed to the argument `prior` of the function `bayes`. The order of the argument `param` in `fn` has the same order as the argument `parvars` when the model is specified by a formula. Otherwise, it is equal to the argument `param` in the function `fimfunc`.

**Note**

The order of the argument `param` in `fn` has the same order as the argument `parvars` when the model is specified by a formula. Otherwise, it is the same as the argument `param` in the function `fimfunc`.

**See Also**

[bayes](#) [sensbayes](#)

**Examples**

```
uniform(lower = c(-3, .1), upper = c(3, 2))
```

# Index

- bayes, [3](#), [3](#), [5](#), [13](#), [17](#), [22](#), [23](#), [35–37](#), [62](#), [64](#),  
[65](#), [67](#), [69](#), [70](#), [75](#), [80](#), [99–102](#)
- bayescomp, [12](#), [35](#), [64](#), [81](#)
- beff, [16](#)
  
- createNIGrid, [4](#), [13](#), [22](#), [23](#), [63](#), [71](#), [76](#), [80](#)
- crt.bayes.control, [4](#), [13](#), [17](#), [22](#), [63](#), [71](#), [76](#),  
[80](#)
- crt.minimax.control, [23](#), [50](#), [51](#), [64](#), [65](#), [89](#)
  
- dmsn, [100](#)
- dmt, [101](#)
  
- family, [3](#), [13](#), [17](#), [39](#), [41](#), [45](#), [50](#), [68](#), [75](#), [79](#),  
[82](#), [86](#), [88](#), [97](#)
- FIM\_2par\_exp\_censor1, [24](#)
- FIM\_2par\_exp\_censor2, [25](#)
- FIM\_3par\_exp\_censor1, [26](#)
- FIM\_3par\_exp\_censor2, [26](#)
- FIM\_exp\_2par, [27](#)
- FIM\_kinetics\_alcohol, [28](#)
- FIM\_logistic, [28](#)
- FIM\_logistic\_2pred, [29](#)
- FIM\_logistic\_4par, [30](#)
- FIM\_loglin, [31](#)
- FIM\_mixed\_inhibition, [31](#)
- FIM\_power\_logistic, [32](#)
- FIM\_sig\_emax, [33](#)
- formula, [3](#), [13](#), [17](#), [38](#), [41](#), [45](#), [49](#), [50](#), [67](#), [75](#),  
[79](#), [82](#), [85](#), [88](#), [97](#)
  
- hcubature, [4](#), [13](#), [22](#), [23](#), [36](#), [63](#), [71](#), [76](#), [80](#)
  
- ICA.control, [3](#), [5](#), [13](#), [14](#), [29](#), [33](#), [35](#), [41](#), [42](#),  
[46](#), [50](#), [53](#), [59](#), [68](#), [69](#)
- ICAOD, [35](#)
- ICAOD-package (ICAOD), [35](#)
- iterate, [37](#)
- iterate.bayes, [37](#), [37](#)
- iterate.minimax, [37](#), [38](#)
  
- leff, [17](#), [38](#), [39](#)
- locally, [35](#), [36](#), [40](#), [46](#), [65](#), [66](#), [69](#)
- locallycomp, [45](#)
  
- minimax, [3](#), [4](#), [13](#), [14](#), [17](#), [35](#), [36](#), [38](#), [39](#), [41](#),  
[46](#), [49](#), [50](#), [51](#), [59](#), [65](#), [66](#), [68](#), [75](#), [80](#),  
[82](#), [86](#), [89](#), [98](#)
- multiple, [30](#), [35](#), [57](#), [96](#)
  
- nl.opts, [24](#), [71](#), [73](#)
- nloptr, [23](#), [24](#), [36](#), [51](#), [70–74](#)
- normal, [3](#), [5](#), [13](#), [17](#), [61](#), [75](#), [80](#)
  
- plot, [5](#), [14](#), [46](#)
- plot.bayes, [62](#)
- plot.minimax, [64](#)
- print.bayes, [65](#)
- print.minimax, [66](#)
- print.sensbayes, [66](#)
- print.sensminimax, [67](#)
  
- robust, [3](#), [35](#), [36](#), [65](#), [66](#), [67](#), [99](#)
  
- sens.bayes.control, [4](#), [13](#), [63](#), [70](#), [75](#), [76](#), [80](#)
- sens.minimax.control, [41](#), [46](#), [50](#), [59](#), [64](#),  
[68](#), [72](#), [83](#), [84](#), [86](#), [87](#), [89–91](#), [95](#), [96](#),  
[98](#), [99](#)
- sensbayes, [6](#), [36](#), [62](#), [66](#), [74](#), [99–102](#)
- sensbayescomp, [15](#), [36](#), [66](#), [79](#)
- senslocally, [36](#), [43](#), [67](#), [82](#)
- senslocallycomp, [85](#)
- sensminimax, [4](#), [36](#), [42](#), [51](#), [53](#), [54](#), [67](#), [68](#), [75](#),  
[76](#), [79](#), [80](#), [82](#), [83](#), [86](#), [88](#), [88](#), [89](#), [91](#),  
[96](#), [98](#), [99](#)
- sensmultiple, [36](#), [60](#), [94](#)
- sensrobust, [36](#), [67](#), [69](#), [70](#), [97](#)
- skewnormal, [3](#), [5](#), [13](#), [17](#), [75](#), [80](#), [100](#)
- student, [3](#), [5](#), [13](#), [17](#), [75](#), [80](#), [101](#)
  
- uniform, [3](#), [5](#), [13](#), [17](#), [75](#), [80](#), [102](#)