

Package ‘rstpm2’

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Description R implementation of generalized survival models (GSMs), smooth accelerated failure time (AFT) models and Markov multi-state models. For the GSMs, $g(S(t|x))=\eta(t,x)$ for a link function g , survival S at time t with covariates x and a linear predictor $\eta(t,x)$. The main assumption is that the time effect(s) are smooth <doi:10.1177/0962280216664760>. For fully parametric models with natural splines, this re-implements Stata's 'stpm2' function, which are flexible parametric survival models developed by Royston and colleagues. We have extended the parametric models to include any smooth parametric smoothers for time. We have also extended the model to include any smooth penalized smoothers from the 'mgcv' package, using penalized likelihood. These models include left truncation, right censoring, interval censoring, gamma frailties and normal random effects <doi:10.1002/sim.7451>, and copulas. For the smooth AFTs, $S(t|x) = S_0(t*\eta(t,x))$, where the baseline survival function $S_0(t)=\exp(-\exp(\eta_0(t)))$ is modelled for natural splines for η_0 , and the time-dependent cumulative acceleration factor $\eta(t,x)=\int_0^t \exp(\eta_1(u,x)) du$ for log acceleration factor $\eta_1(u,x)$. The Markov multi-state models allow for a range of models with smooth transitions to predict transition probabilities, length of stay, utilities and costs, with differences, ratios and standardisation.

URL <https://github.com/mclements/rstpm2>

BugReports <https://github.com/mclements/rstpm2/issues>

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aft *Parametric accelerated failure time model with smooth time functions*

Description

This implements the accelerated failure time models $S_0(t \exp(\beta x))$ and $S_0(\int_0^t \exp(\beta x(u)) du)$. The baseline function $S_0(t^*)$ is modelled as $\exp(-\exp(\eta_0(\log(t^*))))$, where $\eta_0(\log(t^*))$ is a linear predictor using natural splines.

Usage

```
aft(formula, data, smooth.formula = NULL, df = 3,
    tvc = NULL, cure.formula = ~1, control = list(),
    init = NULL, weights = NULL, tvc.intercept = TRUE,
    tvc.integrated = FALSE,
    timeVar = "", time0Var = "",
    cure = FALSE, mixture = FALSE, contrasts = NULL, subset = NULL, ...)
```

Arguments

formula	a formula object, with the response on the left of a \sim operator, and the regression terms (excluding time) on the right. The response should be a survival object as returned by the Surv function. The terms can include linear effects for any time-varying coefficients. [required]
data	a data-frame in which to interpret the variables named in the formula argument. [at present: required]
smooth.formula	a formula for describing the time effects for the linear predictor, excluding the baseline $S_0(t^*)$, but including time-dependent acceleration factors. The time-dependent acceleration factors can be modelled with any smooth functions.
df	an integer that describes the degrees of freedom for the ns function for modelling the baseline log-cumulative hazards function (default=3).
tvc	a list with the names of the time-varying coefficients. This uses natural splines (e.g. <code>tvc=list(hormon=3)</code> is equivalent to <code>smooth.formula=~...+hormon:nsx(log(time),df=3)</code>), which by default does <i>not</i> include an intercept (or main effect) term.
cure.formula	a formula for describing the cure fraction.
control	control argument passed to <code>optim</code> .

<code>init</code>	<code>init</code> should either be <code>FALSE</code> , such that initial values will be determined using Cox regression, or a numeric vector of initial values.
<code>weights</code>	an optional vector of 'prior weights' to be used in the fitting process. Should be <code>NULL</code> or a numeric vector.
<code>tvc.intercept</code>	logical for whether to include an intercept in the time-varying acceleration factor (defaults to <code>TRUE</code>)
<code>tvc.integrated</code>	logical for whether the time-varying acceleration factor should be based on a integration, rather than a cumulative effect (defaults to <code>FALSE</code>)
<code>timeVar</code>	string variable defining the time variable. By default, this is determined from the survival object, however this may be ambiguous if two variables define the time.
<code>time0Var</code>	string variable to determine the entry variable; useful for when more than one data variable is used in the entry time.
<code>cure</code>	logical for whether to model for cure using a non-mixture model (default= <code>FALSE</code>)
<code>mixture</code>	logical for whether to model for cure using a mixture model (default= <code>FALSE</code>)
<code>contrasts</code>	an optional list. See the <code>contrasts.arg</code> of <code>model.matrix.default</code> .
<code>subset</code>	an optional vector specifying a subset of observations to be used in the fitting process.
<code>...</code>	additional arguments to be passed to the <code>mle2</code> .

Details

The implementation extends the `mle2` object from the `bbmle` package. The model inherits all of the methods from the `mle2` class.

Value

An `aft`-class object that inherits from `mle2`-class.

Author(s)

Mark Clements.

See Also

[survreg](#), [coxph](#)

Examples

```
summary(aft(Surv(rectime,censrec==1)~hormon,data=brcancer,df=4))
```

aft-class	<i>Class "stm2" ~~~</i>
-----------	-------------------------

Description

Regression object for aft.

Objects from the Class

Objects can be created by calls of the form `new("aft", ...)` and `aft(...)`.

Slots

`args`: Object of class "list" ~~

Extends

Class for `mle2`, directly.

Methods

plot signature(x = "aft", y = "missing"): ...

lines signature(x = "aft"): ...

predict signature(object = "aft"): ...

predictnl signature(object = "aft", ...): ...

Examples

```
showClass("aft")
```

bhazard	<i>Placemaker function for a baseline hazard function.</i>
---------	--

Description

Defined as the identity function.

Usage

```
bhazard(x)
```

Arguments

`x` Input (and output) value

Value

Returns the input value

brcancer

German breast cancer data from Stata.

Description

See <https://www.stata-press.com/data/r11/brcancer.dta>.

Usage

```
data(brcancer)
```

Format

A data frame with 686 observations on the following 15 variables.

id a numeric vector

hormon hormonal therapy

x1 age, years

x2 menopausal status

x3 tumour size, mm

x4 tumour grade

x5 number of positive nodes

x6 progesterone receptor, fmol

x7 estrogen receptor, fmol

rectime recurrence free survival time, days

censrec censoring indicator

x4a tumour grade \geq 2

x4b tumour grade $=$ 3

x5e $\exp(-0.12 \cdot x5)$

Examples

```
data(brcancer)
## maybe str(brcancer) ; plot(brcancer) ...
```

coef<-	<i>Generic method to update the coef in an object.</i>
--------	--

Description

Generic method to update the coef in an object.

Usage

```
coef(x) <- value
```

Arguments

x	object to be updated
value	value of the coefficient to be updated.

Details

This simple generic method is used for the numerical delta method.

Value

The updated object is returned.

Examples

```
##---- Should be DIRECTLY executable !! ----  
##-- ==> Define data, use random,  
##--or do help(data=index) for the standard data sets.  
  
## The function is currently defined as  
function (x, value)  
  UseMethod("coef<-")
```

colon	<i>Colon cancer.</i>
-------	----------------------

Description

Diagnoses of colon cancer.

Usage

```
data(colon)
```

Format

A data frame with 15564 observations on the following 13 variables.

sex Sex (1=male, 2=female))
 age Age at diagnosis
 stage Clinical stage at diagnosis (1=Unknown, 2=Localised, 3=Regional, 4=Distant)
 mmdx Month of diagnosis
 yydx Year of diagnosis
 surv_mm Survival time in months
 surv_yy Survival time in years
 status Vital status at last contact (1=Alive, 2=Dead: cancer, 3=Dead; other, 4=Lost to follow-up)
 subsite Anatomical subsite of tumour (1=Coecum and ascending, 2=Transverse, 3=Descending and sigmoid, 4=Other and NOS)
 year8594 Year of diagnosis (1=Diagnosed 75-84, 2=Diagnosed 85-94)
 agegrp Age in 4 categories (1=0-44, 2=45-59, 3=60-74, 4=75+)
 dx Date of diagnosis
 exit Date of exit

Details

Caution: there is a colon dataset in the survival package. We recommend using `data(colon, package="rstpm2")` to ensure the correct dataset is used.

Examples

```
data(colon, package="rstpm2") # avoids name conflict with survival::colon
## maybe str(colon) ; ...
```

 cox.tvc

Test for a time-varying effect in the coxph model

Description

Test for a time-varying effect in the coxph model by re-fitting the partial likelihood including a time-varying effect, plot the effect size, and return the re-fitted model. The main advantage of this function over the `tt()` special is that it scales well for moderate sized datasets (cf. `tt` which expands the dataset and scales very poorly).

Usage

```
cox.tvc(obj, var=NULL, method="logt")
```

Arguments

obj	A coxph object. Currently restricted to right censoring with Breslow ties and without stratification, etc.
var	String for the effect name. Currently assumes simple continuous effects.
method	A string representing the possible time transformations. Currently only "logt".

Value

Returns a tvcCoxph object (which inherits from the m1e2 class) of the re-fitted model.

See Also

[coxph](#), [cox.zph](#)

Examples

```
## As per the example for cox.zph:
fit <- coxph(Surv(futime, fustat) ~ age + ecog.ps,
             data=ovarian)
temp <- rstpm2::cox.tvc(fit, "age")
print(temp)           # display the results
plot(temp)           # plot curves
```

eform.stpm2	<i>S3 method for to provide exponentiated coefficients with confidence intervals.</i>
-------------	---

Description

S3 method for to provide exponentiated coefficients with confidence intervals.

Usage

```
eform(object, ...)
## S3 method for class 'stpm2'
eform(object, parm, level = 0.95, method = c("Profile", "Delta"),
       name = "exp(beta)", ...)
## Default S3 method:
eform(object, parm, level = 0.95, method =
c("Delta", "Profile"), name = "exp(beta)", ...)
```

Arguments

object	regression object
parm	not currently used
level	significance level for the confidence interval

method	method for confidence interval estimation
name	name for the fitted value
...	other arguments

grad	<i>gradient function (internal function)</i>
------	--

Description

Numerical gradient for a function at a given value (internal).

Usage

```
grad(func, x, ...)
```

Arguments

func	Function taking a vector argument x (returns a vector of length ≥ 1)
x	vector of arguments for where the gradient is wanted.
...	other arguments to the function

Details

$(\text{func}(x+\text{delta}, \dots) - \text{func}(x-\text{delta}, \dots)) / (2 \text{ delta})$ where delta is the third root of the machine precision times $\text{pmax}(1, \text{abs}(x))$.

Value

A vector if $\text{func}(x)$ has length 1, otherwise a matrix with rows for x and columns for $\text{func}(x)$.

Author(s)

Mark Clements.

See Also

numDelta()

Description

This implements the generalised survival model $g(S(t|x)) = \eta$, where g is a link function, S is survival, t is time, x are covariates and η is a linear predictor. The linear predictor can include either parametric or penalised smoothers for the time effects, for time:covariate interactions and for covariate effects. The main model assumption is that the time effects in the linear predictor are smooth. This extends the class of flexible parametric survival models developed by Royston and colleagues. The model has been extended to include relative survival (excess hazards), Gamma frailties and normal random effects.

Usage

```
gsm(formula, data, smooth.formula = NULL, smooth.args = NULL,
    df = 3, cure = FALSE,
    tvc = NULL, tvc.formula = NULL,
    control = list(), init = NULL,
    weights = NULL, robust = FALSE, baseoff = FALSE,
    timeVar = "", time0Var = "", use.gr = NULL,
    optimiser=NULL, log.time.transform=TRUE,
    reltol=NULL, trace = NULL,
    link.type=c("PH","PO","probit","AH","AO"), theta.A0=0,
    contrasts = NULL, subset = NULL,
    robust_initial=NULL,
    coxph.strata = NULL, coxph.formula = NULL,
    logH.formula = NULL, logH.args = NULL,
    bhazard = NULL, bhazinit=NULL, copula=FALSE,
    frailty = !is.null(cluster) & !robust & !copula,
    cluster = NULL, logtheta=NULL,
    nodes=NULL, RandDist=c("Gamma","LogN"), recurrent = FALSE,
    adaptive = NULL, maxkappa = NULL,
    sp=NULL, criterion=NULL, penalty=NULL,
    smoother.parameters=NULL, Z=~1, outer_optim=NULL,
    alpha=1, sp.init=1,
    penalised=FALSE,
    ...)
stpm2(formula, data, weights=NULL, subset=NULL, coxph.strata=NULL, ...)
pstpm2(formula, data, weights=NULL, subset=NULL, coxph.strata=NULL, ...)
```

Arguments

formula	a formula object, with the response on the left of a \sim operator, and the parametric terms on the right. The response must be a survival object as returned by the Surv function. Specials include <code>cluster</code> and <code>bhazard</code> . [required]
data	a data.frame in which to interpret the variables named in the formula argument.

<code>smooth.formula</code>	either a parametric formula or a penalised mgcv: : gam formula for describing the time effects and time-dependent effects and smoothed covariate effects on the linear predictor scale (default=NULL). The default model is equal to $\sim s(\log(\text{time}), k=-1)$ where <code>time</code> is the time variable.
<code>df</code>	an integer that describes the degrees of freedom for the <code>ns</code> function for modelling the baseline log-cumulative hazard (default=3). Parametric model only.
<code>smooth.args</code>	a list describing the arguments for the <code>s</code> function for modelling the baseline time effect on the linear predictor scale (default=NULL).
<code>tvc</code>	a list with the names of the time-varying coefficients. For a parametric model, this uses natural splines (e.g. <code>tvc=list(hormon=3)</code> is equivalent to <code>smooth.formula=~...+as.numeric</code> which by default does <i>not</i> include an intercept term, hence you should include a main effect. Note that this will convert a logical or factor variable to a numeric value, so the user should use indicators for factor terms. For a penalised model, this uses cubic splines (e.g. <code>tvc=list(hormon=-1)</code> is equivalent to <code>smooth.formula=~...+s(log(time), by=hormon, k=-1)</code>), which by default <i>does</i> include an intercept (or main effect) term (and this code will remove any main effect from formula).
<code>tvc.formula</code>	separate formula for the time-varying effects. This is combined with <code>smooth.formula</code> or the default <code>smooth.formula</code> .
<code>baseoff</code>	Boolean used to determine whether fully define the model using <code>tvc.formula</code> rather than combining <code>logH.formula</code> and <code>tvc.formula</code>
<code>logH.args</code>	as per <code>smooth.args</code> . Deprecated.
<code>logH.formula</code>	as per <code>smooth.formula</code> . Deprecated.
<code>cure</code>	logical for whether to estimate a cure model (parametric model only).
<code>control</code>	list of arguments passed to gsm.control .
<code>init</code>	<code>init</code> should either be NULL, such that initial values will be determined using Cox regression, or a numeric vector of initial values.
<code>coxph.strata</code>	variable in the data argument for stratification of the coxph model fit for estimating initial values.
<code>weights</code>	an optional vector of 'prior weights' to be used in the fitting process. Should be NULL or a numeric vector.
<code>robust</code>	Boolean used to determine whether to use a robust variance estimator.
<code>bhazard</code>	variable for the baseline hazard for relative survival
<code>bhazinit</code>	scalar used to adjust the background cumulative hazards for calculating initial values. Default=0.1. Deprecated argument: use of the <code>control</code> argument is preferred.
<code>copula</code>	logical to indicate whether to use a copula model (experimental)
<code>timeVar</code>	variable defining the time variable. By default, this is determined from the survival object, however this may be ambiguous if two variables define the time
<code>sp</code>	fix the value of the smoothing parameters.
<code>use.gr</code>	in R, a Boolean to determine whether to use the gradient in the optimisation. Default=TRUE, Deprecated argument: use of the <code>control</code> argument is preferred.

criterion	in Rcpp, determine whether to use "GCV" or "BIC" for for the smoothing parameter selection.
penalty	use either the "logH" penalty, which is the default penalty from mgcv, or the "h" hazard penalty. Default="logH". Deprecated argument: use of the control argument is preferred.
smoother.parameters	for the hazard penalty, a list with components which are lists with components var, transform and inverse.
alpha	an ad hoc tuning parameter for the smoothing parameter.
sp.init	initial values for the smoothing parameters.
trace	integer for trace reporting; 0 represents no additional reporting. Default=0. Deprecated argument: use of the control argument is preferred.
contrasts	an optional list. See the contrasts.arg of <code>model.matrix.default</code> .
subset	an optional vector specifying a subset of observations to be used in the fitting process.
coxph.formula	additional formula used to improve the fitting of initial values [optional and rarely used].
time0Var	string variable to determine the entry variable; useful for when more than one data variable is used in the entry time.
link.type	type of link function. For "PH" (generalised proportional hazards), $g(S)=\log(-\log(S))$; for "PO" (generalised proportional odds), $g(S)=-\text{logit}(S)$; for "probit" (generalised probit), $g(S)=-\text{probit}(S)$; for "AH" (generalised additive hazards), $g(S)=-\log(S)$; for "AO" (generalised Aranda-Ordaz), $g(S)=\log((S^{-(\text{theta.AO})}-1)/\text{theta.AO})$.
theta.AO	theta parameter for the Aranda-Ordaz link type.
optimiser	select which optimiser is used. Default="BFGS". Deprecated argument: use of the control argument is preferred.
log.time.transform	should a log-transformation be used for calculating the derivative of the design matrix with respect to time? (default=TRUE)
recurrent	logical for whether clustered, left truncated data are recurrent or for first event (where the latter requires an adjustment for the frailties or random effects)
frailty	logical for whether to fit a shared frailty model
cluster	variable that determines the cluster for the frailty. This can be a vector, a string for the column, or a name. This can also be specified using a special.
logtheta	initial value for log-theta used in the gamma shared frailty model (defaults to value from a coxph model fit)
nodes	number of integration points for Gaussian quadrature. Default=9. Deprecated argument: use of the control argument is preferred.
RandDist	type of distribution for the random effect or frailty
adaptive	logical for whether to use adaptive or non-adaptive quadrature, Default=TRUE. Deprecated argument: use of the control argument is preferred.

<code>maxkappa</code>	double float value for the maximum value of the weight used in the constraint. Default=1000. Deprecated argument: use of the <code>control</code> argument is preferred.
<code>Z</code>	formula for the design matrix for the random effects
<code>reltol</code>	list with components for search and final relative tolerances. Default=list(search=1e-10, final=1e-10, outer=1e-5). Deprecated argument: use of the <code>control</code> argument with arguments <code>reltol.search</code> , <code>reltol.final</code> and <code>reltol.outer</code> is preferred.
<code>outer_optim</code>	Integer to indicate the algorithm for outer optimisation. If <code>outer_optim=1</code> (default), then use Nelder-Mead, otherwise use Nlm.
<code>robust_initial</code>	logical for whether to use Nelder-Mead to find initial values (max 50 iterations). This is useful for ill-posed initial values. Default= FALSE. Deprecated argument: use of the <code>control</code> argument is preferred.
<code>penalised</code>	logical to show whether to use penalised models with <code>pstpm</code> (penalised=TRUE) or parametrics models with <code>stpm2</code> (penalised=FALSE).
<code>...</code>	additional arguments to be passed to the <code>mle2</code> .

Details

The implementation extends the `mle2` object from the `bbmle` package.

The default smoothers for time on the linear predictor scale are `nsxs(log(time),df=3)` for the parametric model and `s(log(time))` for the penalised model.

A frequently asked question is: why does `rstpm2` give different spline estimates to `flexsurv` and Stata's `stpm2`? The short answer is that `rstpm2` uses a different natural spline basis compared with `flexsurv` and Stata's `stpm2` and slightly different knot placement than Stata's `stpm2`. If the knot placement is the same, then the predictions and other coefficients are expected to be very similar. As a longer answer, the default smoother in `rstpm2` is to use an extension of the `splines::ns` function (`rstpm2::nsx`), which uses a QR projection of B-splines for natural splines. In contrast, `flexsurv` and Stata's `stpm2` use truncated power splines for the natural spline basis (also termed 'restricted cubic splines'). The B-splines are known to have good numerical properties, while Stata's `stpm2` implementation defaults to using matrix orthogonalisation to account for any numerical instability in the truncated power basis. Furthermore, `rstpm2` allows for any smooth parametric function to be used as a smoother in `stpm2/gsm`, which is an extension over `flexsurv` and Stata's `stpm2`. Finally, it may be difficult to get `rstpm2` and Stata's `stpm2` to return the same estimates: although `nsx` includes an argument `stata.stpm2.compatible = FALSE` (change to TRUE for compatibility), the design matrix for `rstpm2` is based on individuals with events, while Stata's `stpm2` determines the spline knots from the individuals with events and the design matrix is otherwise based on all individuals.

Value

Either a `stpm2-class` or `pstpm2-class` object.

Author(s)

Mark Clements, Xing-Rong Liu, Benjamin Christoffersen.

Examples

```

## Not run:
data(brcancer)
summary(fit <- stpm2(Surv(rectime,censrec==1)~hormon,data=brcancer,df=3))

## some predictions
head(predict(fit,se.fit=TRUE,type="surv"))
head(predict(fit,se.fit=TRUE,type="hazard"))

## some plots
plot(fit,newdata=data.frame(hormon=0),type="hazard")
plot(fit,newdata=data.frame(hormon=0),type="surv")

## time-varying coefficient
summary(fit.tvc <- stpm2(Surv(rectime,censrec==1)~hormon,data=brcancer,df=3,
  tvc=list(hormon=3)))
anova(fit,fit.tvc) # compare with and without tv

## some more plots
plot(fit.tvc,newdata=data.frame(hormon=0),type="hr",var="hormon",ylim=c(0,2))
lines(fit.tvc,newdata=data.frame(hormon=1),type="hr",var="hormon",
  col=2)

plot(fit.tvc,newdata=data.frame(hormon=0),type="sdiff",var="hormon")

plot(fit.tvc,newdata=data.frame(hormon=0),type="hdiff",var="hormon")

library(scales)
cols <- c(alpha("red",alpha=0.2), alpha("blue",alpha=0.2))
plot(fit.tvc,newdata=data.frame(hormon=0),type="hazard",ci.col=cols[1])
lines(fit.tvc,newdata=data.frame(hormon=1),type="hazard",lty=2,ci.col=cols[2],
  ci=TRUE)
legend("topright",legend=c("No hormonal treatment", "(95
  lty=c(1,1,2,1), lwd=c(1,10,1,10), col=c("black",cols[1],"black",cols[2]), bty="n")

## compare number of knots
hormon0 <- data.frame(hormon=0)
plot(fit,type="hazard",newdata=hormon0)
AIC(fit)
for (df in 4:6) {
  fit.new <- stpm2(Surv(rectime,censrec==1)~hormon,data=brcancer,df=df)
  plot(fit.new,type="hazard",newdata=hormon0,add=TRUE,ci=FALSE,line.col=df)
  print(AIC(fit.new))
}

## compatibility with Stata's stpm2 using the smooth.formula argument (see Details)
summary(stpm2(Surv(rectime,censrec==1)~hormon,data=brcancer,
  smooth.formula=~nsx(log(rectime),df=3,stata.stpm2.compatible=TRUE)))
summary(stpm2(Surv(rectime,censrec==1)~hormon,data=brcancer,
  smooth.formula=~nsx(log(rectime),df=3,stata=TRUE)+
  hormon:nsx(log(rectime),df=3,stata=TRUE)))

```

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

```
gsm.control           Defaults for the gsm call
```

Description

Set useful default and allow changes for the gsm call. This is meant to make the gsm call simpler.

Usage

```
gsm.control(parscale = 1, maxit = 300, optimiser = c("BFGS", "NelderMead"), trace = 0,
            nodes = 9, adaptive = TRUE, kappa.init = 1, maxkappa = 1000,
            suppressWarnings.coxph.frailty = TRUE, robust_initial = FALSE, bhazinit = 0.1,
            eps.init = 1e-5, use.gr = TRUE, penalty = c("logH", "h"), outer_optim = 1,
            reltol.search = 1e-10, reltol.final = 1e-10, reltol.outer = 1e-05,
            criterion = c("GCV", "BIC"))
```

Arguments

<code>parscale</code>	numeric vector or scalar for the scaling of the parameter values; default 1
<code>maxit</code>	integer for the maximum number of iterations for the optimisation process
<code>optimiser</code>	which optimiser to use for the outer optimisation
<code>trace</code>	integer indicating the trace level for each optimiser
<code>nodes</code>	number of quadrature nodes
<code>adaptive</code>	logical for whether to use adaptive or non-adaptive quadrature, Default=TRUE.
<code>kappa.init</code>	initial value for the quadratic penalty for inequality constraints
<code>eps.init</code>	initial value for epsilon
<code>maxkappa</code>	double float value for the maximum value of the weight used in the constraint.
<code>suppressWarnings.coxph.frailty</code>	logical
<code>robust_initial</code>	Not currently documented.
<code>bhazinit</code>	Not currently documented.
<code>use.gr</code>	Logical for whether to use gradients.
<code>penalty</code>	Not currently documented.
<code>outer_optim</code>	Not currently documented.
<code>reltol.search</code>	Relative tolerance. Not currently documented.
<code>reltol.final</code>	Relative tolerance. Not currently documented.
<code>reltol.outer</code>	Relative tolerance. Not currently documented.
<code>criterion</code>	Not currently documented.

gsm_design	<i>Extract design information from an stpm2/gsm object and newdata for use in C++</i>
------------	---

Description

Extract design information from an stpm2/gsm object and newdata for use in C++

Usage

```
gsm_design(object, newdata, newdata0 = NULL, t0 = NULL, inflate = 100)
```

Arguments

object	stpm2/gsm object
newdata	list or data-frame used for evaluation
newdata0	list or data-frame used for evaluation at the entry time
t0	possible delayed entry time (numeric scalar)
inflate	double value to inflate minimum and maximum times for root finding

Value

list that can be read by 'gsm ssim::read_gsm(SEX args)' in C++

incrVar	<i>Utility that returns a function to increment a variable in a data-frame.</i>
---------	---

Description

A functional approach to defining an increment in one or more variables in a data-frame. Given a variable name and an increment value, return a function that takes any data-frame to return a data-frame with incremented values.

Usage

```
incrVar(var, increment = 1)
```

Arguments

var	String for the name(s) of the variable(s) to be incremented
increment	Value that the variable should be incremented.

Details

Useful for defining transformations for calculating rate ratios.

Value

A function with a single data argument that increments the variables in the data list/data-frame.

Examples

```
##---- Should be DIRECTLY executable !! ----
##-- ==> Define data, use random,
##--or do help(data=index) for the standard data sets.

## The function is currently defined as
function (var, increment = 1)
{
  n <- length(var)
  if (n > 1 && length(increment)==1)
    increment <- rep(increment, n)
  function(data) {
    for (i in 1:n) {
      data[[var[i]]] <- data[[var[i]]] + increment[i]
    }
    data
  }
}
```

legendre.quadrature.rule.200

Legendre quadrature rule for n=200.

Description

Legendre quadrature rule for n=200.

Usage

```
data(legendre.quadrature.rule.200)
```

Format

A data frame with 200 observations on the following 2 variables.

x x values between -1 and 1

w weights

Examples

```
data(legendre.quadrature.rule.200)
## maybe str(legendre.quadrature.rule.200) ; ...
```

lines.stpm2

S3 methods for lines

Description

S3 methods for lines

Usage

```
## S3 method for class 'stpm2'
lines(x, newdata = NULL, type = "surv", col = 1, ci.col= "grey",
      lty = par("lty"), ci = FALSE, rug = FALSE, var = NULL,
      exposed = NULL, times = NULL,
      type.relsurv = c("excess", "total", "other"),
      ratetable = survival::survexp.us, rmap, scale = 365.24, ...)
## S3 method for class 'pstpm2'
lines(x, newdata = NULL, type = "surv", col = 1,
      ci.col= "grey",
      lty = par("lty"), ci = FALSE, rug = FALSE, var = NULL,
      exposed = NULL, times = NULL, ...)
```

Arguments

x	an stpm2 object
newdata	required list of new data. This defines the unexposed newdata (<i>excluding</i> the event times).
type	specify the type of prediction
col	line colour
lty	line type
ci.col	confidence interval colour
ci	whether to plot the confidence interval band (default=TRUE)
rug	whether to add a rug plot of the event times to the current plot (default=TRUE)
var	specify the variable name or names for the exposed/unexposed (names are given as characters)
exposed	function that takes newdata and returns the exposed dataset. By default, this increments var (except for cure models, where it defaults to the last event time).
times	specifies the times. By default, this uses a span of the observed times.
type.relsurv	type of predictions for relative survival models: either "excess", "total" or "other"
scale	scale to go from the days in the ratetable object to the analysis time used in the analysis
rmap	an optional list that maps data set names to the ratetable names. See survexp
ratetable	a table of event rates used in relative survival when type.relsurv is "total" or "other"
...	additional arguments (add to the plot command)

 markov_msm

Predictions for continuous time, nonhomogeneous Markov multi-state models using parametric and penalised survival models.

Description

A numerically efficient algorithm to calculate predictions from a continuous time, nonhomogeneous Markov multi-state model. The main inputs are the models for the transition intensities, the initial values, the transition matrix and the covariate patterns. The predictions include state occupancy probabilities (possibly with discounting and utilities), length of stay and costs. Standard errors are calculated using the delta method. Includes, differences, ratios and standardisation.

Usage

```
markov_msm(x, trans, t = c(0,1), newdata = NULL, init=NULL,
           tmvar = NULL,
           sing.inf = 1e+10, method="adams", rtol=1e-10, atol=1e-10, slow=FALSE,
           min.tm=1e-8,
           utility=function(t) rep(1, nrow(trans)),
           utility.sd=rep(0,nrow(trans)),
           use.costs=FALSE,
           transition.costs=function(t) rep(0, sum(!is.na(trans))), # per transition
           transition.costs.sd=rep(0,sum(!is.na(trans))),
           state.costs=function(t) rep(0,nrow(trans)), # per unit time
           state.costs.sd=rep(0,nrow(trans)),
           discount.rate = 0,
           block.size=500,
           spline.interpolation=FALSE,
           debug=FALSE,
           ...)
## S3 method for class 'markov_msm'
vcov(object, ...)
## S3 method for class 'markov_msm'
as.data.frame(x, row.names=NULL, optional=FALSE,
              ci=TRUE,
              P.conf.type="logit", L.conf.type="log",
              C.conf.type="log",
              P.range=c(0,1), L.range=c(0,Inf),
              C.range=c(0,Inf),
              state.weights=NULL, obs.weights=NULL,
              ...)
## S3 method for class 'markov_msm_diff'
as.data.frame(x, row.names=NULL, optional=FALSE,
              P.conf.type="plain", L.conf.type="plain",
              C.conf.type="plain",
              P.range=c(-Inf,Inf), L.range=c(-Inf,Inf),
              C.range=c(-Inf,Inf),
```

```

        ...)
## S3 method for class 'markov_msm_ratio'
as.data.frame(x, row.names=NULL, optional=FALSE, ...)
standardise(x, ...)
## S3 method for class 'markov_msm'
standardise(x,
            weights = rep(1,nrow(x$newdata)),
            normalise = TRUE, ...)
## S3 method for class 'markov_msm'
plot(x, y, stacked=TRUE, which=c('P','L'),
     xlab="Time", ylab=NULL, col=2:6, border=col,
     ggplot2=FALSE, lattice=FALSE, alpha=0.2,
     strata=NULL,
     ...)
## S3 method for class 'markov_msm'
subset(x, subset, ...)
## S3 method for class 'markov_msm'
diff(x, y, ...)
ratio_markov_msm(x, y, ...)
## S3 method for class 'markov_msm'
rbind(..., deparse.level=1)
## S3 method for class 'markov_msm'
transform(`_data`, ...)
collapse_markov_msm(object, which=NULL, sep="; ")
zeroModel(object)
hrModel(object,hr=1,ci=NULL,seloghr=NULL)
aftModel(object,af=1,ci=NULL,selogaf=NULL)
addModel(...)
hazFun(f, tmvar="t", ...)
splineFun(time,rate,method="natural",scale=1,...)

```

Arguments

For markov_msm:

- | | |
|-------|---|
| x | list of functions or parametric or penalised survival models. Currently the models include combinations of stpm2 , pstpm2 , glm , gam , survPen or an object of class "zeroModel" from zeroModel based on one of the other classes. The order in the list matches the indexing in the trans argument. The functions can optionally use a t argument for time and/or a newdata argument. Uncertainty in the models are incorporated into the gradients, while uncertainty in the functions are currently not modelled. |
| trans | Transition matrix describing the states and transitions in the multi-state model. If S is the number of states in the multi-state model, trans should be an S x S matrix, with (i,j)-element a positive integer if a transition from i to j is possible in the multi-state model, NA otherwise. In particular, all diagonal elements should be NA. The integers indicating the possible transitions in the multi-state model should be sequentially numbered, 1, ..., K, with K the number of transitions. See msprep |

<code>t</code>	numerical vector for the times to evaluation the predictions. Includes the start time
<code>newdata</code>	data.frame of the covariates to use in the predictions
<code>init</code>	vector of the initial values with the same length as the number of states. Defaults to the first state having an initial value of 1 (i.e. " <code>[<-</code> "(<code>rep(0, nrow(trans))</code>), 1, 1)).
<code>tmvar</code>	specifies the name of the time variable. This should be set for regression models that do not specify this (e.g. glm) or where the time variable is ambiguous
<code>sing.inf</code>	If there is a singularity in the observed hazard, for example a Weibull distribution with shape < 1 has infinite hazard at $t=0$, then as a workaround, the hazard is assumed to be a large finite number, <code>sing.inf</code> , at this time. The results should not be sensitive to the exact value assumed, but users should make sure by adjusting this parameter in these cases.
<code>method</code>	For <code>markov_msm</code> , the method used by the ordinary differential equation solver. Defaults to Adams method (" <code>adams</code> ") for non-stiff differential equations. For <code>splineFun</code> , the method used for spline interpolation; see <code>splinefun</code> .
<code>rtol</code>	relative error tolerance, either a scalar or an array as long as the number of states. Passed to lsode
<code>atol</code>	absolute error tolerance, either a scalar or an array as long as the number of states. Passed to lsode
<code>slow</code>	logical to show whether to use the slow R-only implementation. Useful for debugging. Currently needed for costs.
<code>min.tm</code>	Minimum time used for evaluations. Avoids <code>log(0)</code> for some models.
<code>utility</code>	a function of the form <code>function(t)</code> that returns a utility for each state at time <code>t</code> for the length of stay values
<code>utility.sd</code>	a function of the form <code>function(t)</code> that returns the standard deviation for the utility for each state at time <code>t</code> for the length of stay values
<code>use.costs</code>	logical for whether to use costs. Default: FALSE
<code>transition.costs</code>	a function of the form <code>function(t)</code> that returns the cost for each transition
<code>transition.costs.sd</code>	a function of the form <code>function(t)</code> that returns the standard deviation for the cost for each transition
<code>state.costs</code>	a function of the form <code>function(t)</code> that returns the cost per unit time for each state
<code>state.costs.sd</code>	a function of the form <code>function(t)</code> that returns the standard deviation for the cost per unit time for each state
<code>discount.rate</code>	numerical value for the proportional reduction (per unit time) in the length of stay and costs
<code>block.size</code>	divide <code>newdata</code> into blocks. Uses less memory but is slower. Reduce this number if the function call runs out of memory.
<code>spline.interpolation</code>	logical for whether to use spline interpolation for the transition hazards rather than the model predictions directly (default=TRUE).

debug logical flag for whether to keep the full output from the ordinary differential equation in the res component (default=FALSE).

... other arguments. For markov_msm, these are passed to the `ode` solver from the `deSolve` package. For `plot.markov_msm`, these arguments are passed to `plot.default`

For `as.data.frame.markov_msm`:

`row.names` add in row names to the output data-frame

`optional` (not currently used)

`ci` logical for whether to include confidence intervals. Default: TRUE

`P.conf.type` type of transformation for the confidence interval calculation for the state occupancy probabilities. Default: log-log transformation. This is changed for `diff` and `ratio_markov_msm` objects

`L.conf.type` type of transformation for the confidence interval calculation for the length of stay calculation. Default: log transformation. This is changed for `diff` and `ratio_markov_msm` objects

`C.conf.type` type of transformation for the confidence interval calculation for the length of stay calculation. Default: log transformation. This is changed for `diff` and `ratio_markov_msm` objects

`P.range` valid values for the state occupancy probabilities. Default: (0,1). This is changed for `diff` and `ratio_markov_msm` objects

`L.range` valid values for the state occupancy probabilities. Default: (0,Inf). This is changed for `diff` and `ratio_markov_msm` objects

`C.range` valid values for the state occupancy probabilities. Default: (0,Inf). This is changed for `diff` and `ratio_markov_msm` objects

`state.weights` Not currently documented

`obs.weights` Not currently documented

For `standardise.markov_msm`:

`weights` numerical vector to use in standardising the state occupancy probabilities, length of stay and costs. Default: 1 for each observation.

`normalise` logical for whether to normalise the weights to 1. Default: TRUE

For `plot.markov_msm`:

`y` (currently ignored)

`stacked` logical for whether to stack the plots. Default: TRUE

`xlab` x-axis label

`ylab` x-axis label

`col` colours (ignored if `ggplot2=TRUE`)

`border` border colours for the `polygon` (ignored if `ggplot2=TRUE`)

`ggplot2` use `ggplot2`

alpha alpha value for confidence bands (ggplot)
 lattice use lattice
 strata formula for the stratification factors for the plot

For subset.markov_msm:

subset expression that is evaluated on the newdata component of the object to filter (or restrict) for the covariates used for predictions

For transform.markov_msm:

_data an object of class "markov_msm"

For rbind.markov_msm:

deparse.level not currently used

For collapse.states:

which either an index of the states to collapse or a character vector of the state names to collapse

sep separator to use for the collapsed state names

For zeroModel to predict zero rates:

object survival regression object to be wrapped

For hrModel to predict rates times a hazard ratio:

hr hazard ratio

seloghr alternative specification for the se of the log(hazard ratio); see also ci argument

For aftModel to predict accelerated rates:

af acceleration factor

selogaf alternative specification for the se of the log(acceleration factor); see also ci argument

addModel predict rates based on adding rates from different models

hazFun provides a rate function without uncertainty:

f rate function, possibly with tmvar and/or newdata as arguments

splineFun predicts rates using spline interpolation:

time exact times

rate rates as per time

scale rate multiplier (e.g. scale=365.25 for converting from daily rates to yearly rates)

Details

The predictions are calculated using an ordinary differential equation solver. The algorithm uses a single run of the solver to calculate the state occupancy probabilities, length of stay, costs and their partial derivatives with respect to the model parameters. The predictions can also be combined to calculate differences, ratios and standardised.

The current implementation supports a list of models for each transition.

The current implementation also only allows for a vector of initial values rather than a matrix. The predictions will need to be re-run for different vectors of initial values.

For `as.data.frame.markov_msm_ratio`, the data are provided in log form, hence the default transformations and bounds are as per `as.data.frame.markov_msm_diff`, with untransformed data on the real line.

TODO: allow for one model to predict for the different transitions.

Value

`markov_msm` returns an object of class `"markov_msm"`.

The function `summary` is used to obtain and print a summary and analysis of variance table of the results. The generic accessor functions `coef` and `vcov` extract various useful features of the value returned by `markov_msm`.

An object of class `"markov_msm"` is a list containing at least the following components:

<code>time</code>	a numeric vector with the times for the predictions
<code>P</code>	an <code>array</code> for the predicted state occupancy probabilities. The array has three dimensions: time, state, and observations.
<code>L</code>	an <code>array</code> for the predicted sojourn times (or length of stay). The array has three dimensions: time, state, and observations.
<code>Pu</code>	an <code>array</code> for the partial derivatives of the predicted state occupancy probabilities with respect to the model coefficients. The array has four dimensions: time, state, coefficients, and observations.
<code>Lu</code>	an <code>array</code> for the partial derivatives of the predicted sojourn times (or length of stay) with respect to the model coefficients. The array has four dimensions: time, state, coefficients, and observations.
<code>newdata</code>	a <code>data.frame</code> with the covariates used for the predictions
<code>vcov</code>	the variance-covariance matrix for the models of the transition intensities
<code>trans</code>	copy of the <code>trans</code> input argument
<code>call</code>	the call to the function

For debugging:

<code>res</code>	data returned from the ordinary differential equation solver. This may include more information on the predictions
------------------	--

Author(s)

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See Also

[pmatrix.fs](#), [probtrans](#)

Examples

```
## Not run:
if (requireNamespace("deSolve")) {
  library(readstata13)
  library(mstate)
  library(ggplot2)
  library(survival)
  ## Two states: Initial -> Final
  ## Note: this shows how to use markov_msm to estimate survival and risk probabilities based on
  ## smooth hazard models.
  two_states <- function(model, ...) {
    transmat = matrix(c(NA,1,NA,NA),2,2,byrow=TRUE)
    rownames(transmat) <- colnames(transmat) <- c("Initial","Final")
    rstpm2::markov_msm(list(model), ..., trans = transmat)
  }
  ## Note: the first argument is the hazard model. The other arguments are arguments to the
  ## markov_msm function, except for the transition matrix, which is defined by the new function.
  death = gsm(Surv(time,status)~factor(rx), data=survival::colon, subset=(etype==2), df=3)
  cr = two_states(death, newdata=data.frame(rx="Obs"), t = seq(0,2500, length=301))
  plot(cr,ggplot=TRUE)

  ## Competing risks
  ## Note: this shows how to adapt the markov_msm model for competing risks.
  competing_risks <- function(listOfModels, ...) {
    nRisks = length(listOfModels)
    transmat = matrix(NA,nRisks+1,nRisks+1)
    transmat[1,1+(1:nRisks)] = 1:nRisks
    rownames(transmat) <- colnames(transmat) <- c("Initial",names(listOfModels))
    rstpm2::markov_msm(listOfModels, ..., trans = transmat)
  }
  ## Note: The first argument for competing_risks is a list of models. Names from that list are
  ## used for labelling the states. The other arguments are as per the markov_msm function,
  ## except for the transition matrix, which is defined by the competing_risks function.
  recurrence = gsm(Surv(time,status)~factor(rx), data=survival::colon, subset=(etype==1), df=3)
  death = gsm(Surv(time,status)~factor(rx), data=survival::colon, subset=(etype==2), df=3)
  cr = competing_risks(list(Recurrence=recurrence,Death=death),
    newdata=data.frame(rx=levels(survival::colon$rx)),
    t = seq(0,2500, length=301))
  ## Plot the probabilities for each state for three different treatment arms
  plot(cr, ggplot=TRUE) + facet_grid(~ rx)
  ## And: differences in probabilities
  cr_diff = diff(subset(cr,rx=="Lev+5FU"),subset(cr,rx=="Obs"))
  plot(cr_diff, ggplot=TRUE, stacked=FALSE)

  ## Extended example: Crowther and Lambert (2017)
  ## library(rstpm2); library(readstata13); library(ggplot2)
  mex.1 <- read.dta13("http://fmwww.bc.edu/repec/bocode/m/multistate_example.dta")
  transmat <- rbind("Post-surgery"=c(NA,1,2),
```

```

      "Relapsed"=c(NA,NA,3),
      "Died"=c(NA,NA,NA))
colnames(transmat) <- rownames(transmat)
mex.2 <- transform(mex.1,osi=(osi=="deceased")+0)
levels(mex.2$size)[2] <- ">20-50 mm" # fix typo
mex <- mstate::msprep(time=c(NA,"rf","os"),status=c(NA,"rfi","osi"),
  data=mex.2,trans=transmat,id="pid",
  keep=c("age","size","nodes","pr_1","hormon"))
mex <- transform(mex,
  size2=(unclass(size)==2)+0, # avoids issues with TRUE/FALSE
  size3=(unclass(size)==3)+0,
  hormon=(hormon=="yes")+0,
  Tstart=Tstart/12,
  Tstop=Tstop/12)
##
c.ar <- stpm2(Surv(Tstart,Tstop,status) ~ age + size2 + size3 + nodes + pr_1 + hormon,
  data = mex, subset=trans==1, df=3, tvc=list(size2=1,size3=1,pr_1=1))
c.ad <- stpm2(Surv(Tstart, Tstop, status) ~ age + size + nodes + pr_1 + hormon,
  data = mex, subset=trans==2, df=1)
c.rd <- stpm2( Surv(Tstart,Tstop,status) ~ age + size + nodes + pr_1 + hormon,
  data=mex, subset=trans==3, df=3, tvc=list(pr_1=1))
##
nd <- expand.grid(nodes=seq(0,20,10), size=levels(mex$size))
nd <- transform(nd, age=54, pr_1=3, hormon=0,
  size2=(unclass(size)==2)+0,
  size3=(unclass(size)==3)+0)
## Predictions
system.time(pred1 <- rstpm2::markov_msm(list(c.ar,c.ad,c.rd), t = seq(0,15,length=301),
  newdata=nd, trans = transmat)) # ~2 seconds
pred1 <- transform(pred1, Nodes=paste("Nodes =",nodes), Size=paste("Size",size))
## Figure 3
plot(pred1, ggplot=TRUE) + facet_grid(Nodes ~ Size) + xlab("Years since surgery")
plot(pred1, ggplot=TRUE, flipped=TRUE) +
  facet_grid(Nodes ~ Size) + xlab("Years since surgery")
plot(pred1, strata=~nodes+size, xlab="Years since surgery", lattice=TRUE)
## Figure 4
plot(subset(pred1, nodes==0 & size=="<=20 mm"), stacked=FALSE, ggplot=TRUE) +
  facet_grid(. ~ state) +
  xlab("Years since surgery")
## Figure 5
a <- diff(subset(pred1,nodes==0 & size=="<=20 mm"),
  subset(pred1,nodes==0 & size==">20-50 mm"))
a <- transform(a, label = "Prob(Size<=20 mm)-Prob(20mm<Size<50mm)")
b <- ratio_markov_msm(subset(pred1,nodes==0 & size=="<=20 mm"),
  subset(pred1,nodes==0 & size==">20-50 mm"))
b <- transform(b,label="Prob(Size<=20 mm)-Prob(20mm<Size<50mm)")
##
c <- diff(subset(pred1,nodes==0 & size=="<=20 mm"),
  subset(pred1,nodes==0 & size==">50 mm"))
c <- transform(c, label = "Prob(Size<=20 mm)-Prob(Size>=50mm)")
d <- ratio_markov_msm(subset(pred1,nodes==0 & size=="<=20 mm"),
  subset(pred1,nodes==0 & size==">50 mm"))
d <- transform(d,label= "Prob(Size<=20 mm)-Prob(Size>=50mm)")

```

```

##
e <- diff(subset(pred1,nodes==0 & size==">20-50 mm"),
          subset(pred1,nodes==0 & size==">50 mm"))
e <- transform(e,label="Prob(20mm<Size<50 mm)-Prob(Size>=50mm)")
f <- ratio_markov_msm(subset(pred1,nodes==0 & size==">20-50 mm"),
                     subset(pred1,nodes==0 & size==">50 mm"))
f <- transform(f, label = "Prob(20mm<Size<50 mm)-Prob(Size>=50mm)")
## combine
diffs <- rbind(a,c,e)
ratios <- rbind(b,d,f)
## Figure 5
plot(diffs, stacked=FALSE, ggplot2=TRUE) + xlab("Years since surgery") +
  ylim(c(-0.4, 0.4)) + facet_grid(label ~ state)
##
plot(ratios, stacked=FALSE, ggplot2=TRUE) + xlab("Years since surgery") +
  ylim(c(0, 3)) + facet_grid(label ~ state)
## Figure 6
plot(subset(pred1, nodes==0 & size=="<=20 mm"), stacked=FALSE, which="L", ggplot2=TRUE) +
  facet_grid(. ~ state) + xlab("Years since surgery")
## Figure 7
plot(diffs, stacked=FALSE, which="L", ggplot2=TRUE) + xlab("Years since surgery") +
  ylim(c(-4, 4)) + facet_grid(label ~ state)
plot(ratios, stacked=FALSE, which="L", ggplot2=TRUE) + xlab("Years since surgery") +
  ylim(c(0.1, 10)) + coord_trans(y="log10") + facet_grid(label ~ state)
}

## End(Not run)

```

markov_sde

Predictions for continuous time, nonhomogeneous Markov multi-state models using Aalen's additive hazards models.

Description

A numerically efficient algorithm to calculate predictions from a continuous time, nonhomogeneous Markov multi-state model. The main inputs are a list of Aalen's additive hazards models, the initial values, the transition matrix and the covariate patterns. The predictions include state occupancy probabilities and length of stay. Standard errors are calculated using the delta method. Includes differences and standardisation.

Usage

```

markov_sde(models, trans, newdata, init = NULL, nLebesgue = 10000 + 1, los = FALSE,
           nOut = 300, weights = 1)
## S3 method for class 'markov_sde'
standardise(x, ...)
## S3 method for class 'markov_sde'
plot(x, y, stacked=TRUE, which=c("P","L"), index=NULL,
     xlab="Time", ylab=NULL, col=2:6, border=col,

```

```

    ggplot2=FALSE, lattice=FALSE, alpha=0.2,
    strata=NULL,
    ...)
## S3 method for class 'markov_sde'
as.data.frame(x, row.names=NULL, optional=NULL, ci=TRUE,
  P.conf.type="logit", L.conf.type="log",
  P.range=c(0,1), L.range=c(0,Inf),
  ...)

```

Arguments

models	list of models. Currently allows only for aalen regression models.
trans	Transition matrix describing the states and transitions in the multi-state model. If S is the number of states in the multi-state model, trans should be an $S \times S$ matrix, with (i,j)-element a positive integer if a transition from i to j is possible in the multi-state model, NA otherwise. In particular, all diagonal elements should be NA. The integers indicating the possible transitions in the multi-state model should be sequentially numbered, $1, \dots, K$, with K the number of transitions. See msprep
newdata	data.frame of the covariates to use in the predictions
init	vector of the initial values with the same length as the number of states. Defaults to the first state having an initial value of 1 (i.e. " <code>[<-</code> "(rep(0, nrow(trans)), 1, 1)).
nLebesgue	Number of steps for the continuous integration
los	logical variable for whether to estimate the length of stay
nOut	number of rows to represent the continuous changes
weights	numeric vector to represent differences or standardisation

For `plot.markov_sde`:

y	(currently ignored)
stacked	logical for whether to stack the plots. Default: TRUE
index	indicator of which row of newdata to plot
which	character to indicate either transition probabilities ("P") or length of stay ("L"). Default: "P".
xlab	x-axis label
ylab	x-axis label
col	colours (ignored if ggplot2=TRUE)
border	border colours for the polygon (ignored if ggplot=TRUE)
ggplot2	use ggplot2
alpha	alpha value for confidence bands (ggplot)
lattice	use lattice
strata	formula for the stratification factors for the plot

For `as.data.frame.markov_sde`:

row.names	add in row names to the output data-frame
optional	(not currently used)
ci	logical for whether to include confidence intervals. Default: TRUE
P.conf.type	type of transformation for the confidence interval calculation for the state occupancy probabilities. Default: logit transformation. This is changed to "identity" if any of the weights are negative
L.conf.type	type of transformation for the confidence interval calculation for the length of stay calculation. Default: log transformation. "identity" if any of the weights are negative
P.range	valid values for the state occupancy probabilities. Default: (0,1).
L.range	valid values for the state occupancy probabilities. Default: (0,Inf).

For `standardise.markov_sde`:

x	object to extract standardised values
...	other arguments. For <code>plot.markov_sde</code> , these arguments are passed to <code>plot.default</code> . For <code>standardise.markov_sde</code> , these arguments are not used, as the standardisation must be done earlier in <code>markov_sde</code> .

Details

Uses an approach developed by Ryalen and colleagues. This is a re-implementation in C++.

The current implementation only allows for a vector of initial values rather than a matrix. The predictions will need to be re-run for different vectors of initial values.

Value

`markov_sde` returns an object of class "markov_sde".

Author(s)

Mark Clements

See Also

[markov_msm](#)

nsx *Generate a Basis Matrix for Natural Cubic Splines (with eXtensions)*

Description

Generate the B-spline basis matrix for a natural cubic spline (with eXtensions).

Usage

```
nsx(x, df = NULL, knots = NULL, intercept = FALSE,
    Boundary.knots = range(x), derivs = if (cure) c(2, 1) else c(2, 2),
    log = FALSE, centre = FALSE,
    cure = FALSE, stata.stpm2.compatible = FALSE)
```

Arguments

x	the predictor variable. Missing values are allowed.
df	degrees of freedom. One can supply df rather than knots; ns() then chooses $df - 1 - intercept + 4 - \text{sum}(\text{derivs})$ knots at suitably chosen quantiles of x (which will ignore missing values).
knots	breakpoints that define the spline. The default is no knots; together with the natural boundary conditions this results in a basis for linear regression on x. Typical values are the mean or median for one knot, quantiles for more knots. See also Boundary.knots.
intercept	if TRUE, an intercept is included in the basis; default is FALSE.
Boundary.knots	boundary points at which to impose the natural boundary conditions and anchor the B-spline basis (default the range of the data). If both knots and Boundary.knots are supplied, the basis parameters do not depend on x. Data can extend beyond Boundary.knots
derivs	an integer vector of length 2 with values between 0 and 2 giving the derivative constraint order at the left and right boundary knots; an order of 2 constrains the second derivative to zero ($f''(x)=0$); an order of 1 constrains the first and second derivatives to zero ($f'(x)=f''(x)=0$); an order of 0 constrains the zero, first and second derivatives to zero ($f(x)=f'(x)=f''(x)=0$)
log	a Boolean indicating whether the underlying values have been log transformed; (deprecated: only used to calculate derivatives in rstpm2:::stpm2Old)
centre	if specified, then centre the splines at this value (i.e. $f(\text{centre})=0$) (default=FALSE)
cure	a Boolean indicated whether to estimate cure; changes the default derivs argument, such that the right boundary has the first and second derivatives constrained to zero; defaults to FALSE
stata.stpm2.compatible	a Boolean to determine whether to use Stata stpm's default knot placement; defaults to FALSE

Value

A matrix of dimension $\text{length}(x) * \text{df}$ where either `df` was supplied or if `knots` were supplied, $\text{df} = \text{length}(\text{knots}) + 1 + \text{intercept}$. Attributes are returned that correspond to the arguments to `ns`, and explicitly give the knots, `Boundary.knots` etc for use by `predict.nsx()`.

`nsx()` is based on the functions `ns` and `spline.des`. It generates a basis matrix for representing the family of piecewise-cubic splines with the specified sequence of interior knots, and the natural boundary conditions. These enforce the constraint that the function is linear beyond the boundary knots, which can either be supplied, else default to the extremes of the data. A primary use is in modeling formula to directly specify a natural spline term in a model.

The extensions from `ns` are: specification of the derivative constraints at the boundary knots; whether to centre the knots; incorporation of cure using derivatives; compatible knots with Stata's `stpm2`; and an indicator for a log-transformation of `x` for calculating derivatives.

References

Hastie, T. J. (1992) Generalized additive models. Chapter 7 of *Statistical Models in S* eds J. M. Chambers and T. J. Hastie, Wadsworth & Brooks/Cole.

See Also

`ns`, `bs`, `predict.nsx`, `SafePrediction`

Examples

```
require(stats); require(graphics); require(splines)
nsx(women$height, df = 5)
summary(fm1 <- lm(weight ~ ns(height, df = 5), data = women))

## example of safe prediction
plot(women, xlab = "Height (in)", ylab = "Weight (lb)")
ht <- seq(57, 73, length.out = 200)
lines(ht, predict(fm1, data.frame(height=ht)))
```

nsxD

Generate a Basis Matrix for the first derivative of Natural Cubic Splines (with eXtensions)

Description

Generate the B-spline basis matrix for the first derivative of a natural cubic spline (with eXtensions).

Usage

```
nsxD(x, df = NULL, knots = NULL, intercept = FALSE,
     Boundary.knots = range(x), derivs = if (cure) c(2, 1) else c(2, 2),
     log = FALSE, centre = FALSE,
     cure = FALSE, stata.stpm2.compatible = FALSE)
```

Arguments

<code>x</code>	the predictor variable. Missing values are allowed.
<code>df</code>	degrees of freedom. One can supply <code>df</code> rather than <code>knots</code> ; <code>ns()</code> then chooses $df - 1 - \text{intercept} + 4 - \text{sum}(\text{derivs})$ knots at suitably chosen quantiles of <code>x</code> (which will ignore missing values).
<code>knots</code>	breakpoints that define the spline. The default is no knots; together with the natural boundary conditions this results in a basis for linear regression on <code>x</code> . Typical values are the mean or median for one knot, quantiles for more knots. See also <code>Boundary.knots</code> .
<code>intercept</code>	if TRUE, an intercept is included in the basis; default is FALSE.
<code>Boundary.knots</code>	boundary points at which to impose the natural boundary conditions and anchor the B-spline basis (default the range of the data). If both <code>knots</code> and <code>Boundary.knots</code> are supplied, the basis parameters do not depend on <code>x</code> . Data can extend beyond <code>Boundary.knots</code>
<code>derivs</code>	an integer vector of length 2 with values between 0 and 2 giving the derivative constraint order at the left and right boundary knots; an order of 2 constrains the second derivative to zero ($f''(x)=0$); an order of 1 constrains the first and second derivatives to zero ($f'(x)=f''(x)=0$); an order of 0 constrains the zero, first and second derivatives to zero ($f(x)=f'(x)=f''(x)=0$)
<code>log</code>	a Boolean indicating whether the underlying values have been log transformed; (deprecated: only used to calculate derivatives in <code>rstpm2:::stpm2Old</code>)
<code>centre</code>	if specified, then centre the splines at this value (i.e. $f(\text{centre})=0$) (default=FALSE)
<code>cure</code>	a Boolean indicated whether to estimate cure; changes the default <code>derivs</code> argument, such that the right boundary has the first and second derivatives constrained to zero; defaults to FALSE
<code>stata.stpm2.compatible</code>	a Boolean to determine whether to use Stata <code>stpm</code> 's default knot placement; defaults to FALSE

Value

A matrix of dimension $\text{length}(x) * df$ where either `df` was supplied or if `knots` were supplied, $df = \text{length}(\text{knots}) + 1 + \text{intercept}$. Attributes are returned that correspond to the arguments to `ns`, and explicitly give the `knots`, `Boundary.knots` etc for use by `predict.nsxD()`.

`nsxD()` is based on the functions `ns` and `spline.des`. It generates a basis matrix for representing the family of piecewise-cubic splines with the specified sequence of interior knots, and the natural boundary conditions. These enforce the constraint that the function is linear beyond the boundary knots, which can either be supplied, else default to the extremes of the data. A primary use is in modeling formula to directly specify a natural spline term in a model.

The extensions from `ns` are: specification of the derivative constraints at the boundary knots; whether to centre the knots; incorporation of cure using derivatives; compatible knots with Stata's `stpm2`; and an indicator for a log-transformation of `x` for calculating derivatives.

References

Hastie, T. J. (1992) Generalized additive models. Chapter 7 of *Statistical Models in S* eds J. M. Chambers and T. J. Hastie, Wadsworth & Brooks/Cole.

See Also

[ns](#), [bs](#), [predict.nsx](#), [SafePrediction](#)

Examples

```
require(stats); require(graphics); require(splines)
nsx(women$height, df = 5)
summary(fm1 <- lm(weight ~ ns(height, df = 5), data = women))

## example of safe prediction
plot(women, xlab = "Height (in)", ylab = "Weight (lb)")
ht <- seq(57, 73, length.out = 200)
lines(ht, predict(fm1, data.frame(height=ht)))
```

numDeltaMethod

Calculate numerical delta method for non-linear predictions.

Description

Given a regression object and an independent prediction function (as a function of the coefficients), calculate the point estimate and standard errors

Usage

```
numDeltaMethod(object, fun, gd=NULL, conf.int=FALSE, level=0.95, ...)
```

Arguments

object	A regression object with methods <code>coef</code> and <code>vcov</code> .
fun	An independent prediction function with signature <code>function(coef, ...)</code> .
gd	Specified gradients
conf.int	Logical for whether to also calculate the confidence interval
level	Numeric for the level of the confidence interval
...	Other arguments passed to <code>fun</code> .

Details

A more user-friendly interface is provided by `predictnl`.

Value

fit	Point estimates
se.fit	Standard errors
Estimate	Point estimates
SE	Standard errors
conf.low	Lower confidence interval (if <code>conf.int=TRUE</code>)
conf.high	Upper confidence interval (if <code>conf.int=TRUE</code>)

See Also

See Also [predictnl](#).

Examples

```
##---- Should be DIRECTLY executable !! ----
##-- ==> Define data, use random,
##--or do help(data=index) for the standard data sets.

## The function is currently defined as
function (object, fun, ...)
{
  coef <- coef(object)
  est <- fun(coef, ...)
  Sigma <- vcov(object)
  gd <- grad(fun, coef, ...)
  se.est <- as.vector(sqrt(colSums(gd * (Sigma %*% gd))))
  data.frame(Estimate = est, SE = se.est)
}
```

plot-methods

plots for an stpm2 fit

Description

Given an stpm2 fit, return a plot

Usage

```
## S4 method for signature 'stpm2'
plot(x,y,newdata,type="surv",
      xlab="Time",line.col=1,ci.col="grey",
      add=FALSE,ci=TRUE,rug=TRUE,
      var=NULL,exposed=NULL,times=NULL,...)
## S4 method for signature 'pstpm2'
plot(x,y,newdata,type="surv",
      xlab="Time",line.col=1,ci.col="grey",
      add=FALSE,ci=TRUE,rug=TRUE,
      var=NULL,exposed=NULL,times=NULL,...)
```

Arguments

x	an stpm2 object
y	not used (for generic compatibility)
newdata	required list of new data. This defines the unexposed newdata (<i>excluding</i> the event times).
type	specify the type of prediction

xlab	x-axis label
line.col	line colour
ci.col	confidence interval colour
ci	whether to plot the confidence interval band (default=TRUE)
add	whether to add to the current plot (add=TRUE) or make a new plot (add=FALSE) (default=FALSE)
rug	whether to add a rug plot of the event times to the current plot (default=TRUE)
var	specify the variable name or names for the exposed/unexposed (names are given as characters)
exposed	function that takes newdata and returns the exposed dataset. By default, this increments var (except for cure models, where it defaults to the last event time).
times	specifies the times. By default, this uses a span of the observed times.
...	additional arguments (add to the plot command)

Methods

`x = "stpm2", y = "missing"` an stpm2 fit

See Also

[stpm2](#)

popmort

Background mortality rates for the colon dataset.

Description

Background mortality rates for the colon dataset.

Usage

```
data(popmort)
```

Format

A data frame with 10600 observations on the following 5 variables.

sex Sex (1=male, 2=female)

prob One year probability of survival

rate All cause mortality rate

age Age by single year of age through to age 105 years

year Calendar period

Examples

```
data(popmort)
## maybe str(popmort) ; ...
```

predict-methods	<i>Predicted values for an stpm2 or pstpm2 fit</i>
-----------------	--

Description

Given an stpm2 fit and an optional list of new data, return predictions

Usage

```
## S4 method for signature 'stpm2'
predict(object, newdata=NULL,
        type=c("surv", "cumhaz", "hazard", "density", "hr", "sdiff",
              "hdiff", "loghazard", "link", "meansurv", "meansurvdiff", "meanhr",
              "odds", "or", "margsurv", "marghaz", "marghr", "meanhaz", "af",
              "fail", "margfail", "meanmargsurv", "uncured", "rmst", "probcure",
              "lpmatrix", "gradh", "gradH", "rmstdiff", "lpmatrixD"),
        grid=FALSE, seqLength=300,
        type.relsurv=c("excess", "total", "other"), scale=365.24,
        rmap, ratetable=survival::survexp.us,
        se.fit=FALSE, link=NULL, exposed=NULL, var=NULL,
        keep.attributes=FALSE, use.gr=TRUE, level=0.95,
        n.gauss.quad=100, full=FALSE, ...)

## S4 method for signature 'pstpm2'
predict(object, newdata=NULL,
        type=c("surv", "cumhaz", "hazard", "density", "hr", "sdiff",
              "hdiff", "loghazard", "link", "meansurv", "meansurvdiff", "meanhr",
              "odds", "or", "margsurv", "marghaz", "marghr", "meanhaz", "af",
              "fail", "margfail", "meanmargsurv", "rmst", "lpmatrix",
              "gradh", "gradH", "rmstdiff", "lpmatrixD"),
        grid=FALSE, seqLength=300,
        se.fit=FALSE, link=NULL, exposed=NULL, var=NULL,
        keep.attributes=FALSE, use.gr=TRUE, level=0.95,
        n.gauss.quad=100, full=FALSE, ...)
```

Arguments

object	an stpm2 or pstpm2 object
newdata	optional list of new data (required if type in ("hr", "sdiff", "hdiff", "meansurvdiff", "or", "uncured")). For type in ("hr", "sdiff", "hdiff", "meansurvdiff", "or", "af", "uncured"), this defines the unexposed newdata. This can be combined with grid to get a regular set of event times (i.e. newdata would <i>not</i> include the event times).
type	specify the type of prediction: "surv" survival probabilities "cumhaz" cumulative hazard "hazard" hazard "density" density

	"hr" hazard ratio
	"sdiff" survival difference
	"hdiff" hazard difference
	"loghazard" log hazards
	"meansurv" mean survival
	"meansurvdiff" mean survival difference
	"odds" odds
	"or" odds ratio
	"margsurv" marginal (population) survival
	"marghaz" marginal (population) hazard
	"marghr" marginal (population) hazard ratio
	"meanhaz" mean hazard
	"meanhr" mean hazard ratio
	"af" attributable fraction
	"fail" failure (=1-survival)
	"margfail" marginal failure (=1-marginal survival)
	"meanmargsurv" mean marginal survival, averaged over the frailty distribution
	"uncured" distribution for the uncured
	"rmst" restricted mean survival time
	"rmstdiff" restricted mean survival time difference
	"probcure" probability of cure
	"lpmatrix" design matrix
	"lpmatrixD" design matrix for the derivative with respect to time
grid	whether to merge newdata with a regular sequence of event times (default=FALSE)
seqLength	length of the sequence used when grid=TRUE
type.relsurv	type of predictions for relative survival models: either "excess", "total" or "other"
scale	scale to go from the days in the ratetable object to the analysis time used in the analysis
rmap	an optional list that maps data set names to the ratetable names. See survexp
ratetable	a table of event rates used in relative survival when type.relsurv is "total" or "other"
se.fit	whether to calculate confidence intervals (default=FALSE)
link	allows a different link for the confidence interval calculation (default=NULL, such that switch(type,surv="cloglog",cumhaz="log",hazard="log",hr="log",sdiff="I",hdiff="I",loghazard="I",link="I",odds="log",or="log",margsurv="cloglog",marghaz="log",marghr="log"))
exposed	a function that takes newdata and returns a transformed data-frame for those exposed or the counterfactual. By default, this increments var (except for cure models, where it defaults to the last event time).
var	specify the variable name or names for the exposed/unexposed (names are given as characters)

keep.attributes	Boolean to determine whether the output should include the newdata as an attribute (default=TRUE)
use.gr	Boolean to determine whether to use gradients in the variance calculations when they are available (default=TRUE)
level	confidence level for the confidence intervals (default=0.95)
n.gauss.quad	number of Gauassian quadrature points used for integrations (default=100)
full	logical for whether to return a full data-frame with predictions and newdata combined. Useful for lattice and ggplot2 plots. (default=FALSE)
...	additional arguments (for generic compatibility)

Details

The confidence interval estimation is based on the delta method using numerical differentiation.

Value

A data-frame with components Estimate, lower and upper, with an attribute "newdata" for the newdata data-frame.

Methods

object= "stpm2" an stpm2 fit

See Also

[stpm2](#)

predict.nsx

Evaluate a Spline Basis

Description

Evaluate a predefined spline basis at given values.

Usage

```
## S3 method for class 'nsx'
predict(object, newx, ...)
```

Arguments

object	the result of a call to <code>nsx</code> having attributes describing knots, degree, etc.
newx	the x values at which evaluations are required.
...	Optional additional arguments. At present no additional arguments are used.

Value

An object just like `object`, except evaluated at the new values of `x`.

These are methods for the generic function `predict` for objects inheriting from classes "nsx". See `predict` for the general behavior of this function.

See Also

`nsx`.

Examples

```
basis <- nsx(women$height, df = 5)
newX <- seq(58, 72, length.out = 51)
# evaluate the basis at the new data
predict(basis, newX)
```

predictnl

Estimation of standard errors using the numerical delta method.

Description

A simple, yet exceedingly useful, approach to estimate the variance of a function using the numerical delta method. A number of packages provide functions that analytically calculate the gradients; we use numerical derivatives, which generalises to models that do not offer analytical derivatives (e.g. ordinary differential equations, integration), or to examples that are tedious or error-prone to calculate (e.g. sums of predictions from GLMs).

Usage

```
## Default S3 method:
predictnl(object, fun, newdata=NULL, gd=NULL, ...)
## S3 method for class 'lm'
predictnl(object, fun, newdata=NULL, ...)
## S3 method for class 'formula'
predict(object, data, newdata, na.action, type="model.matrix", ...)
## S3 method for class 'predictnl'
confint(object, parm, level=0.95, ...)
```

Arguments

<code>object</code>	An object with <code>coef</code> , <code>vcov</code> and <code>`coef<-`</code> methods (required).
<code>fun</code>	A function that takes <code>object</code> as the first argument, possibly with <code>newdata</code> and other arguments (required). See notes for why it is often useful to include <code>newdata</code> as an argument to the function.
<code>newdata</code>	An optional argument that defines <code>newdata</code> to be passed to <code>fun</code> .

gd	An optional matrix of gradients. If this is not specified, then the gradients are calculated using finite differences.
parm	currently ignored
level	significance level for 2-sided confidence intervals
data	object used to define the model frame
na.action	passed to model.frame
type	currently restricted to "model.matrix"
...	Other arguments that are passed to fun.

Details

The signature for fun is either `fun(object, ...)` or `fun(object, newdata=NULL, ...)`.

The different `predictnl` methods call the utility function `numDeltaMethod`, which in turn calls the `grad` function for numerical differentiation. The `numDeltaMethod` function calls the standard `coef` and `vcov` methods, and the non-standard ``coef<-`` method for changing the coefficients in a regression object. This non-standard method has been provided for several regression objects and essentially mirrors the `coef` method.

One potential issue is that some `predict` methods do not re-calculate their predictions for the fitted dataset (i.e. when `newdata=NULL`). As the `predictnl` function changes the fitted coefficients, it is required that the predictions are re-calculated. One solution is to pass `newdata` as an argument to both `predictnl` and `fun`; alternatively, `newdata` can be specified in `fun`. These approaches are described in the examples below. The `numDeltaMethod` method called by `predictnl` provides a warning when the variance estimates are zero, which may be due to this cause.

For completeness, it is worth discussing why the example `predictnl(fit,predict)` does not work for when `fit` is a `glm` object. First, `predict.glm` does not update the predictions for the fitted data. Second, the default `predict` method has a signature `predict(object, ...)`, which does not include a `newdata` argument. We could then either (i) require that a `newdata` argument be passed to the `fun` function for all examples, which would make this corner case work, or (ii) only pass the `newdata` argument if it is non-null or in the formals for the `fun` function, which would fail for this corner case. The current API defaults to the latter case (ii). To support this approach, the `predictnl.lm` method replaces a null `newdata` with `object$data`. We also provide a revised `numdelta:::predict.lm` method that performs the same operation, although its use is not encouraged due to its clumsiness.

Value

Returns an object of class `an object with class c("predictnl", "data.frame")` elements `c("fit", "se.fit", "Estimate"`, and with methods `print` and `confint`. Note that the `Estimate` and `SE` fields are deprecated and their use is discouraged, as we would like to remove them from future releases.

Author(s)

Mark Clements

Examples

```
df <- data.frame(x=0:1, y=c(10, 20))
fit <- glm(y ~ x, df, family=poisson)

predictnl(fit,
          function(obj,newdata)
            diff(predict(obj,newdata,type="response")))
```

predictnl-methods *~~ Methods for Function predictnl ~~*

Description

~~ Methods for function predictnl ~~

Methods

predictnl signature(object = "mle2", ...): Similar to predictnl.default, using S4 methods.

pstpm2-class *Class "pstpm2"*

Description

Regression object for pstpm2.

Objects from the Class

Objects can be created by calls of the form new("pstpm2", ...) and pstpm2(...).

Slots

xlevels: Object of class "list" ~~
 contrasts: Object of class "listOrNULL" ~~
 terms: Object of class "terms" ~~
 gam: Object of class "gam" ~~
 logli: Object of class "function" ~~
 timeVar: Object of class "character" ~~
 time0Var: Object of class "character" ~~
 time0Expr: Object of class "nameOrcall" ~~
 timeExpr: Object of class "nameOrcall" ~~
 like: Object of class "function" ~~

```

model.frame: Object of class "list" ~~
delayed: Object of class "logical" ~~
frailty: Object of class "logical" ~~
x: Object of class "matrix" ~~
xd: Object of class "matrix" ~~
termsd: Object of class "terms" ~~
Call: Object of class "character" ~~
y: Object of class "Surv" ~~
sp: Object of class "numeric" ~~
nevent: Object of class "numeric" ~~
link: Object of class "list" ~~
edf: Object of class "numeric" ~~
edf_var: Object of class "numeric" ~~
df: Object of class "numeric" ~~
call: Object of class "language" ~~
call.orig: Object of class "language" ~~
coef: Object of class "numeric" ~~
fullcoef: Object of class "numeric" ~~
vcov: Object of class "matrix" ~~
min: Object of class "numeric" ~~
details: Object of class "list" ~~
minuslogl: Object of class "function" ~~
method: Object of class "character" ~~
data: Object of class "list" ~~
formula: Object of class "character" ~~
optimizer: Object of class "character" ~~
args: Object of class "list" ~~

```

Extends

Class for [mle2](#), directly.

Methods

```

plot signature(x = "pstpm2", y = "missing"): ...
lines signature(x = "pstpm2", ...): ...
anova signature(object = "pstpm2", ...): ...
AIC signature(object = "pstpm2", ..., k=2): ...
AICc signature(object = "pstpm2", ..., nobs=NULL, k=2): ...

```

BIC signature(object = "pstpm2", ..., nobs = NULL): ...
qAICc signature(object = "pstpm2", ..., nobs = NULL, dispersion = 1, k = 2): ...
qAIC signature(object = "pstpm2", ..., dispersion = 1, k = 2): ...
summary signature(object = "pstpm2", ...): ...
eform signature(object = "pstpm2", ...): ...
predictnl signature(object = "pstpm2", ...): ...

Examples

```
showClass("pstpm2")
```

residuals-methods *Residual values for an stpm2 or pstpm2 fit*

Description

Given an stpm2 or pstpm2 fit, return residuals

Usage

```
## S4 method for signature 'stpm2'
residuals(object, type=c("li", "gradli"))
## S4 method for signature 'pstpm2'
residuals(object, type=c("li", "gradli"))
```

Arguments

object an stpm2 or pstpm2 object
type specify the type of residuals:
 "li" log-likelihood components (not strictly residuals)
 "gradli" gradient of the log-likelihood components (not strictly residuals)

Details

The gradients are analytical.

Value

A vector or matrix.

Methods

object= "stpm2" an stpm2 fit

See Also

[stpm2](#)

rstpm2-internal	<i>Internal functions for the rstpm2 package.</i>
-----------------	---

Description

Various utility functions used internally to the rstpm2 package.

Usage

```
lhs(formula)
rhs(formula)
lhs(formula) <- value
rhs(formula) <- value
```

Arguments

formula	A formula
value	A symbolic value to replace the current value.

simulate-methods	<i>Simulate values from an stpm2 or pstpm2 fit</i>
------------------	--

Description

Given an stpm2 fit and a data-frame of new data, return simulated values

Usage

```
## S4 method for signature 'stpm2'
simulate(object, nsim=1,
         seed=NULL, newdata=NULL,
         lower=1e-06, upper=1e+05, start=NULL, ...)
## S4 method for signature 'pstpm2'
simulate(object, nsim=1,
         seed=NULL, newdata=NULL,
         lower=1e-06, upper=1e+05, start=NULL, ...)
```

Arguments

object	an stpm2 or pstpm2 object
nsim	number of simulations per row in newdata
seed	optional random number seed
newdata	list of new data. If not specified, then defaults to object@data
lower	smallest possible time

upper	largest possible time
start	left truncated entry time (assumed to be zero if NULL)
...	additional arguments (for generic compatibility)

Methods

object = "stpm2" an stpm2 fit

Examples

```
set.seed(1002)
fit1 <- gsm(Surv(rectime, censrec==1)~hormon, data=brcancer, df=3)
simulate(fit1, nsim=10, newdata=data.frame(hormon=1))
simulate(fit1, newdata=data.frame(hormon=0:1))
```

smoothpwc	<i>Utility to use a smooth function in markov_msm based on piece-wise constant values</i>
-----------	---

Description

Utility to use a smooth function in markov_msm based on piece-wise constant values

Usage

```
smoothpwc(midts, rates, tmvar = "t", offsetvar = "", ...)
```

Arguments

midts	mid-point values for time in each segment
rates	rates at those mid-points (or for the interval)
tmvar	string for the time variable
offsetvar	string for a time offset variable
...	other arguments

Details

Uses splines to smooth the log-rates. This assumes that the rates are strictly greater than zero.

Value

a function that is used in markov_msm

Examples

```
##---- Should be DIRECTLY executable !! ----
##-- ==> Define data, use random,
##--or do help(data=index) for the standard data sets.

## The function is currently defined as
function (midts, rates, tmvar = "t", offsetvar = "", ...)
{
  log.smoother <- splinefunx(midts, log(rates), constant.right = TRUE)
  haz <- function(newdata) {
    t <- newdata[[tmvar]] + (if (offsetvar != "")
      newdata[[offsetvar]]
    else 0)
    exp(log.smoother(t))
  }
  structure(list(haz = haz), class = "smoothpwc")
}
```

stpm2-class

Class "stpm2" ~~~

Description

Regression object for stpm2.

Objects from the Class

Objects can be created by calls of the form `new("stpm2", ...)` and `stpm2(...)`.

Slots

xlevels: Object of class "list" ~~
contrasts: Object of class "listOrNULL" ~~
terms: Object of class "terms" ~~
logli: Object of class "function" ~~
lm: Object of class "lm" ~~
timeVar: Object of class "character" ~~
time0Var: Object of class "character" ~~
timeExpr: Object of class "nameOrcall" ~~
time0Expr: Object of class "nameOrcall" ~~
delayed: Object of class "logical" ~~
frailty: Object of class "logical" ~~
interval: Object of class "logical" ~~
model.frame: Object of class "list" ~~

```
call.formula: Object of class "formula" ~~  
x: Object of class "matrix" ~~  
xd: Object of class "matrix" ~~  
termsd: Object of class "terms" ~~  
Call: Object of class "character" ~~  
y: Object of class "Surv" ~~  
link: Object of class "list" ~~  
call: Object of class "language" ~~  
call.orig: Object of class "language" ~~  
coef: Object of class "numeric" ~~  
fullcoef: Object of class "numeric" ~~  
vcov: Object of class "matrix" ~~  
min: Object of class "numeric" ~~  
details: Object of class "list" ~~  
minuslogl: Object of class "function" ~~  
method: Object of class "character" ~~  
data: Object of class "list" ~~  
formula: Object of class "character" ~~  
optimizer: Object of class "character" ~~  
args: Object of class "list" ~~
```

Extends

Class [mle2](#), directly.

Methods

```
plot signature(x = "stpm2", y = "missing"): ...  
lines signature(x = "stpm2", ...): ...  
predictnl signature(object = "stpm2", ...): ...  
summary signature(object = "stpm2", ...): ...  
eform signature(object = "stpm2", ...): ...
```

Examples

```
showClass("stpm2")
```

tvcCoxph-class	Class "tvcCoxph"
----------------	------------------

Description

Experimental approach to modelling time-dependent effects in Cox regression.

Objects from the Class

Objects can be created by calls of the form `new("tvcCoxph", ...)` or `cox.tvc(...)`. See the [mle2](#) documentation.

Slots

call: Object of class "language" ~~
call.orig: Object of class "language" ~~
coef: Object of class "numeric" ~~
fullcoef: Object of class "numeric" ~~
vcov: Object of class "matrix" ~~
min: Object of class "numeric" ~~
details: Object of class "list" ~~
minuslogl: Object of class "function" ~~
method: Object of class "character" ~~
data: Object of class "list" ~~
formula: Object of class "character" ~~
optimizer: Object of class "character" ~~

Extends

Class [mle2](#), directly.

Methods

plot signature(x = "tvcCoxph", y = "missing"): ...

Examples

```
showClass("tvcCoxph")
```

update-methods	<i>Methods for Function update</i>
----------------	------------------------------------

Description

Methods for function update

Methods

update signature(object = "stpm2", ...): Similar to update.default, using S4 methods.

voptimize	<i>Vectorised One Dimensional Optimization</i>
-----------	--

Description

The function `voptimize` searches the interval from `lower` to `upper` for a minimum or maximum of the vectorised function `f` with respect to its first argument.

`optimise` is an alias for `optimize`.

Usage

```
voptimize(f, interval, ...,
         lower=pmin(interval[,1], interval[,2]),
         upper=pmax(interval[,1], interval[,2]),
         maximum = FALSE,
         tol = .Machine$double.eps^0.25)
optimise(f, interval, ...,
        lower=pmin(interval[,1], interval[,2]),
        upper=pmax(interval[,1], interval[,2]),
        maximum = FALSE,
        tol = .Machine$double.eps^0.25)
```

Arguments

<code>f</code>	the function to be optimized. The function is either minimized or maximized over its first argument depending on the value of <code>maximum</code> .
<code>interval</code>	a matrix with two columns containing the end-points of the interval to be searched for the minimum.
<code>...</code>	additional named or unnamed arguments to be passed to <code>f</code>
<code>lower, upper</code>	the lower and upper end points of the interval to be searched.
<code>maximum</code>	logical. Should we maximize or minimize (the default)?
<code>tol</code>	the desired accuracy.

Details

Note that arguments after `...` must be matched exactly.

The method used is a combination of golden section search and successive parabolic interpolation, and was designed for use with continuous functions. Convergence is never much slower than that for a Fibonacci search. If `f` has a continuous second derivative which is positive at the minimum (which is not at lower or upper), then convergence is superlinear, and usually of the order of about 1.324.

The function `f` is never evaluated at two points closer together than $\epsilon|x_0| + (tol/3)$, where ϵ is approximately `sqrt(.Machine$double.eps)` and x_0 is the final abscissa `optimize()`\$minimum.

If `f` is a unimodal function and the computed values of `f` are always unimodal when separated by at least $\epsilon|x| + (tol/3)$, then x_0 approximates the abscissa of the global minimum of `f` on the interval lower, upper with an error less than $\epsilon|x_0| + tol$.

If `f` is not unimodal, then `optimize()` may approximate a local, but perhaps non-global, minimum to the same accuracy.

The first evaluation of `f` is always at $x_1 = a + (1 - \phi)(b - a)$ where $(a, b) = (\text{lower}, \text{upper})$ and $\phi = (\sqrt{5} - 1)/2 = 0.61803..$ is the golden section ratio. Almost always, the second evaluation is at $x_2 = a + \phi(b - a)$. Note that a local minimum inside $[x_1, x_2]$ will be found as solution, even when `f` is constant in there, see the last example.

`f` will be called as `f(x, ...)` for a numeric value of `x`.

The argument passed to `f` has special semantics and used to be shared between calls. The function should not copy it.

The implementation is a vectorised version of the `optimize` function.

Value

A list with components `minimum` (or `maximum`) and `objective` which give the location of the minimum (or maximum) and the value of the function at that point.

Source

Based on R's C translation of Fortran code <https://netlib.org/fmm/fmin.f> (author(s) unstated) based on the Algol 60 procedure `localmin` given in the reference.

References

Brent, R. (1973) *Algorithms for Minimization without Derivatives*. Englewood Cliffs, NJ: Prentice-Hall.

See Also

[optimize](#) for the standard single optimiser solver, [nlm](#), [uniroot](#).

Examples

```
library(graphics)

f <- function(x, a) (x - a)^2
```

```
xmin <- voptimize(f, lower=c(0, 0), upper=c(1,1), tol = 0.0001, a = c(1/3,2/3))
xmin

## See where the function is evaluated:
voptimize(function(x) x^2*(print(x)-1), lower = c(0,0), upper = c(10,10))

## "wrong" solution with unlucky interval and piecewise constant f():
f <- function(x) ifelse(x > -1, ifelse(x < 4, exp(-1/abs(x - 1)), 10), 10)
fp <- function(x) { print(x); f(x) }

plot(f, -2,5, ylim = 0:1, col = 2)
voptimize(fp, cbind(-4, 20)) # doesn't see the minimum
voptimize(fp, cbind(-7, 20)) # ok
```

vuniroot

Vectorised One Dimensional Root (Zero) Finding

Description

The function `vuniroot` searches the interval from `lower` to `upper` for a root (i.e., zero) of the vectorised function `f` with respect to its first argument.

Setting `extendInt` to a non-"no" string, means searching for the correct interval = `c(lower, upper)` if `sign(f(x))` does not satisfy the requirements at the interval end points; see the 'Details' section.

Usage

```
vuniroot(f, interval, ...,
        lower, upper,
        f.lower = f(lower, ...), f.upper = f(upper, ...),
        extendInt = c("no", "yes", "downX", "upX"), check.conv = FALSE,
        tol = .Machine$double.eps^0.25, maxiter = 1000, trace = 0,
        n = NULL)
```

Arguments

<code>f</code>	the function for which the root is sought.
<code>interval</code>	a matrix with two columns containing the end-points of the interval to be searched for the root.
<code>...</code>	additional named or unnamed arguments to be passed to <code>f</code>
<code>lower, upper</code>	the lower and upper end points of the interval to be searched.
<code>f.lower, f.upper</code>	the same as <code>f(upper)</code> and <code>f(lower)</code> , respectively. Passing these values from the caller where they are often known is more economical as soon as <code>f()</code> contains non-trivial computations.

<code>extendInt</code>	character string specifying if the interval <code>c(lower, upper)</code> should be extended or directly produce an error when <code>f()</code> does not have differing signs at the endpoints. The default, "no", keeps the search interval and hence produces an error. Can be abbreviated.
<code>check.conv</code>	logical indicating whether a convergence warning of the underlying <code>vuniroot</code> should be caught as an error and if non-convergence in <code>maxiter</code> iterations should be an error instead of a warning.
<code>tol</code>	the desired accuracy (convergence tolerance).
<code>maxiter</code>	the maximum number of iterations.
<code>trace</code>	integer number; if positive, tracing information is produced. Higher values giving more details.
<code>n</code>	integer number; size of input vector to <code>f</code> (only used if <code>lower</code> and <code>upper</code> are of length 1)

Details

Note that arguments after `...` must be matched exactly.

Either `interval` or both `lower` and `upper` must be specified: the upper endpoint must be strictly larger than the lower endpoint.

The function values at the endpoints must be of opposite signs (or zero), for `extendInt="no"`, the default. Otherwise, if `extendInt="yes"`, the interval is extended on both sides, in search of a sign change, i.e., until the search interval $[l, u]$ satisfies $f(l) \cdot f(u) \leq 0$.

If it is *known how* f changes sign at the root x_0 , that is, if the function is increasing or decreasing there, `extendInt` can (and typically should) be specified as "upX" (for "upward crossing") or "downX", respectively. Equivalently, define $S := \pm 1$, to require $S = \text{sign}(f(x_0 + \epsilon))$ at the solution. In that case, the search interval $[l, u]$ possibly is extended to be such that $S \cdot f(l) \leq 0$ and $S \cdot f(u) \geq 0$.

`vuniroot()` uses a C++ subroutine based on "zeroin" (from Netlib) and algorithms given in the reference below. They assume a continuous function (which then is known to have at least one root in the interval).

Convergence is declared either if $f(x) == 0$ or the change in x for one step of the algorithm is less than `tol` (plus an allowance for representation error in x).

If the algorithm does not converge in `maxiter` steps, a warning is printed and the current approximation is returned.

`f` will be called as `f(x, ...)` for a numeric value of x .

The argument passed to `f` has special semantics and used to be shared between calls. The function should not copy it.

Value

A list with at least three components: `root` and `f.root` give the location of the root and the value of the function evaluated at that point. `iter` gives the number of iterations used.

Further components may be added in future: component `init.it` was added in R 3.1.0.

Source

Based on 'zeroin.c' in <https://netlib.org/c/brent.shar>.

References

Brent, R. (1973) *Algorithms for Minimization without Derivatives*. Englewood Cliffs, NJ: Prentice-Hall.

See Also

[vuniroot](#) for the standard single root solver [polyroot](#) for all complex roots of a polynomial; [optimize](#), [nlm](#).

Examples

```
require(utils) # for str

## some platforms hit zero exactly on the first step:
## if so the estimated precision is 2/3.
f <- function(x, a) x - a
str(xmin <- vuniroot(f, lower=c(0, 0), upper=c(1,1), tol = 0.0001, a = c(1/3,2/3)))
## same example with scalars for lower and upper -- using the n argument
str(xmin <- vuniroot(f, lower=0, upper=1, tol = 0.0001, n=2, a = c(1/3,2/3)))

## handheld calculator example: fixed point of cos(.):
vuniroot(function(x) cos(x) - x, lower = -pi, upper = pi, tol = 1e-9)$root

str(vuniroot(function(x) x*(x^2-1) + .5, lower = -2, upper = 2,
              tol = 0.0001))
str(vuniroot(function(x) x*(x^2-1) + .5, lower = -2, upper = 2,
              tol = 1e-10))

## Find the smallest value x for which exp(x) > 0 (numerically):
r <- vuniroot(function(x) 1e80*exp(x) - 1e-300, cbind(-1000, 0), tol = 1e-15)
str(r, digits.d = 15) # around -745, depending on the platform.

exp(r$root)      # = 0, but not for r$root * 0.999...
minexp <- r$root * (1 - 10*.Machine$double.eps)
exp(minexp)      # typically denormalized

##--- vuniroot() with new interval extension + checking features: -----

f1 <- function(x) (121 - x^2)/(x^2+1)
f2 <- function(x) exp(-x)*(x - 12)

tools::assertCondition(vuniroot(f1, cbind(0,10)),
                      "error", verbose=TRUE)
tools::assertCondition(vuniroot(f2, cbind(0, 2)),
                      "error", verbose=TRUE)
##--> error: f() .. end points not of opposite sign
```

```

## where as 'extendInt="yes"' simply first enlarges the search interval:
u1 <- vuniroot(f1, cbind(0,10),extendInt="yes", trace=1)
u2 <- vuniroot(f2, cbind(0,2), extendInt="yes", trace=2)
stopifnot(all.equal(u1$root, 11, tolerance = 1e-5),
          all.equal(u2$root, 12, tolerance = 6e-6))

## The *danger* of interval extension:
## No way to find a zero of a positive function, but
## numerically, f(-|M|) becomes zero :
tools::assertCondition(u3 <- vuniroot(exp, cbind(0,2), extendInt="yes", trace=TRUE),
                      "error", verbose=TRUE)

## Nonsense example (must give an error):
tools::assertCondition(vuniroot(function(x) 1, cbind(0,1), extendInt="yes"),
                      "error", verbose=TRUE)

## Convergence checking :
sinc_ <- function(x) ifelse(x == 0, 1, sin(x)/x)
curve(sinc_, -6,18); abline(h=0,v=0, lty=3, col=adjustcolor("gray", 0.8))

vuniroot(sinc_, cbind(0,5), extendInt="yes", maxiter=4) #-> "just" a warning

## now with check.conv=TRUE, must signal a convergence error :

vuniroot(sinc_, cbind(0,5), extendInt="yes", maxiter=4, check.conv=TRUE)

### Weibull cumulative hazard (example origin, Ravi Varadhan):
cumhaz <- function(t, a, b) b * (t/b)^a
froot <- function(x, u, a, b) cumhaz(x, a, b) - u

n <- 10
u <- -log(runif(n))
a <- 1/2
b <- 1
## Find failure times
ru <- vuniroot(froot, u=u, a=a, b=b, interval= cbind(rep(1.e-14,n), rep(1e4,n)),
              extendInt="yes")$root
ru2 <- vuniroot(froot, u=u, a=a, b=b, interval= cbind(rep(0.01,n), rep(10,n)),
              extendInt="yes")$root
stopifnot(all.equal(ru, ru2, tolerance = 6e-6))

r1 <- vuniroot(froot, u= 0.99, a=a, b=b, interval= cbind(0.01, 10),
              extendInt="up")
stopifnot(all.equal(0.99, cumhaz(r1$root, a=a, b=b)))

## An error if 'extendInt' assumes "wrong zero-crossing direction":
vuniroot(froot, u= 0.99, a=a, b=b, interval= cbind(0.1, 10), extendInt="down")

```

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