

# Package ‘QHScrnomo’

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**Title** Construct Nomogram for Competing Risk Models

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**Description** Nomogram is constructed to predict cumulative incidence rate, which is calculated by adjusting for competing causes other than that of interest. K-fold cross-validation function is implemented to validate the performance. Competing risk version of concordance index is calculated. Methods are as described in: Kattan MW, Heller G, Brennan MF (2003) <[doi:10.1002/sim.1574](https://doi.org/10.1002/sim.1574)>.

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---

addOffset4ModelFrame *Internal function to Calculate Offset for Data*

---

## Description

Internal function to Calculate Offset for Data

## Usage

```
addOffset4ModelFrame(Terms, newdata, offset = 0)
```

## Arguments

Terms	term value
newdata	data frame for prediction. Each row of the data frame contains values of covariates that are required in the crr model. If missing, the original data set that was used to develop the crr model will be used for prediction.
offset	offset value

## Details

Internally calculate offset for a data

## Value

newdata

---

anova.cmprsk	<i>anova table for competing risks regression</i>
--------------	---

---

## Description

generate anova table for crr

## Usage

```
## S3 method for class 'cmprsk'  
anova(object, ...)
```

## Arguments

object	a competing risks regression model object built from function <a href="#">crr.fit</a>
...	other arguments

## Details

generate anova table for competing risks regression model

## Value

anova table in matrix

## Examples

```
data(prostate.dat)  
dd <- datadist(prostate.dat)  
options(datadist = "dd")  
prostate.f <- cph(Surv(TIME_EVENT, EVENT_DOD == 1) ~ TX + rcs(PSA, 3) +  
  BX_GLSN_CAT + CLIN_STG + rcs(AGE, 3) +  
  RACE_AA, data = prostate.dat,  
  x = TRUE, y = TRUE, surv = TRUE, time.inc = 144)  
prostate.crr <- crr.fit(prostate.f, cencode = 0, failcode = 1)  
## anova test  
anova(prostate.crr)
```

---

cindex *Concordance index calculation*

---

### Description

Calculate concordance index

### Usage

```
cindex(
  prob,
  fstatus,
  ftime,
  type = "crr",
  failcode = 1,
  cencode = 0,
  tol = 1e-20
)
```

### Arguments

prob	predicted risk of failure event, either probability or risk score
fstatus	failure(event) variable
ftime	follow-up time variable for survival or competing risks predictions
type	type of regression models corresponding to different type of outcomes. 'logis' is the default value for binary outcome, 'surv' for ordinary survival outcome and 'crr' for competing risks outcome.
failcode	coding for failure(event). 1 is the default value.
cencode	coding for censoring. 0 is the default
tol	error tolerance. the default value is 1e-20.

### Details

to calculate the discrimination metric, concordance index for binary, time-to event and competing risks outcomes

### Value

a vector of returned values.

N	the total number of observations in the input data
n	the nonmissing number of observations that was used for calculation
usable	the total number of usable pairs.
concordant	the number of concordant pairs
cindex	the concordance index that equal to the number of concordant pairs divided by the total number of usable pairs.

**Author(s)**

Changhong Yu, Michael Kattan, Brian Wells, Amy Nowacki.

**Examples**

```
data(prostate.dat)
dd <- datadist(prostate.dat)
options(datadist = "dd")
prostate.f <- cph(Surv(TIME_EVENT,EVENT_DOD == 1) ~ TX + rcs(PSA,3) +
  BX_GLSN_CAT + CLIN_STG + rcs(AGE,3) +
  RACE_AA, data = prostate.dat,
  x = TRUE, y= TRUE, surv=TRUE,time.inc = 144)
prostate.crr <- crr.fit(prostate.f,cencode = 0,failcode = 1)

## ten fold cross validation
prostate.dat$preds.tenf.cv.prostate.crr.120 <-
  tenf.crr(prostate.crr,time = 120)

## calculate the CRR version of concordance index
with(prostate.dat, cindex(preds.tenf.cv.prostate.crr.120 ,
  ftime = TIME_EVENT,
  fstatus =EVENT_DOD, type = "crr"))["cindex"]
```

---

crr.fit

*Fit Competing Risks Regression Model Fits a competing risks regression model from an existing Cox proportional hazards object and allows a nomogram to be constructed from the competing risks regression object.*

---

**Description**

This function uses the `crr` function in the `cmprsk` package to construct a competing risk regression object.

**Usage**

```
crr.fit(fit, cencode = 0, failcode = 1)
```

**Arguments**

<code>fit</code>	a Cox proportional hazards regression model constructed from <code>cph</code> in <code>rms</code> library (by Frank Harrell)
<code>cencode</code>	the value of the status indicator that indicates a censored observation
<code>failcode</code>	the value of the status indicator that indicates an event of interest

**Value**

Returns a list of class `cmprsk`, with components:

<code>coef</code>	the estimated regression coefficients
<code>loglik</code>	log pseudo-likelihood evaluated at <code>coef</code>
<code>lscore</code>	derivatives of the log pseudo-likelihood evaluated at <code>coef</code>
<code>inf</code>	-second derivatives of the log pseudo-likelihood
<code>var</code>	estimated variance covariance matrix of <code>coef</code>
<code>res</code>	matrix of residuals giving the contribution to each score (columns) at each unique failure time (rows)
<code>uftime</code>	vector of unique failure times
<code>bfitj</code>	jumps in the Breslow-type estimate of the underlying sub-distribution cumulative hazard (used by <code>predict.crr()</code> )
<code>tfs</code>	the <code>tfs</code> matrix (output of <code>tf()</code> , if used)
<code>converged</code>	TRUE if the iterative algorithm converged.
<code>cencode</code>	the value of the status indicator that indicates a censored observation
<code>failcode</code>	the value of the status indicator that indicates an event of interest
<code>cph.f</code>	regular survival model fitted by <code>cph</code> which is saved for function <code>nomogram.crr</code> to adjust <code>lp</code> for competing risks
<code>cphdat</code>	data used for <code>cph</code> model, where all predictors are represented in numeric format, which is used by function <code>tenf.crr</code> to do ten fold cross-validation

**Note**

This function requires that the `rms` and `cmprsk` libraries are attached.

**Author(s)**

Michael W. Kattan, Ph.D. and Changhong Yu. Department of Quantitative Health Sciences, Cleveland Clinic

**References**

Michael W. Kattan, Glenn Heller and Murray F. Brennan (2003). A competing-risks nomogram for sarcoma-specific death following local recurrence. *Statistics in Medicine*. *Stat Med.* 2003;22:3515-3525.

**See Also**

[cph.crr](#) [nomogram.crr](#)

**Examples**

```

data(prostate.dat)
dd <- datadist(prostate.dat)
options(datadist = "dd")
prostate.f <- cph(Surv(TIME_EVENT,EVENT_DOD == 1) ~ TX + rcs(PSA,3) +
  BX_GLSN_CAT + CLIN_STG + rcs(AGE,3) +
  RACE_AA, data = prostate.dat,
  x = TRUE, y= TRUE, surv=TRUE,time.inc = 144)
prostate.crr <- crr.fit(prostate.f,cencode = 0,failcode = 1)
## anova test
anova(prostate.crr)
## hazards ratio
summary(prostate.crr)

## ten fold cross validation
prostate.dat$preds.tenf.cv.prostate.crr.120 <-
  tenf.crr(prostate.crr,time = 120)

## make a CRR nomogram
nomogram.crr(prostate.crr,failtime = 120,lp=FALSE,
funlabel = "Predicted 10-year cumulative incidence")

## calculate the CRR version of concordance index
with(prostate.dat, cindex(preds.tenf.cv.prostate.crr.120 ,
  ftime = TIME_EVENT,
  fstatus =EVENT_DOD, type = "crr"))["cindex"]

## generate the calibration curve for predicted 10-year cancer
## specific mortality

with(prostate.dat,
  groupci(preds.tenf.cv.prostate.crr.120 , ftime = TIME_EVENT,
  fstatus =EVENT_DOD, g = 5, u = 120,
  xlab = "Nomogram predicted 10-year cancerspecific mortality",
  ylab = "Observed predicted 10-year cancerspecific mortality")
)

```

---

groupci

*make calibration curve for competing risks endpoint*


---

**Description**

Cumulative Incidence Estimates vs. a Continuous Variable

**Usage**

```

groupci(
  x,
  ftime,
  fstatus,
  cencode = 0,
  failcode = 1,
  ci = TRUE,
  m = 50,
  g,
  cuts,
  u,
  pl = TRUE,
  conf.int = 0.95,
  xlab,
  ylab,
  xlim = c(0, 1),
  ylim = c(0, 1),
  lty = 1,
  add = FALSE,
  cex.subtitle = FALSE,
  ab = TRUE,
  ...
)

```

**Arguments**

<code>x</code>	a continuous variable
<code>ftime</code>	vector of follow-up time
<code>fstatus</code>	vector of failure status
<code>cencode</code>	value indicating centering.
<code>failcode</code>	value indicating event of interest
<code>ci</code>	logical flag to output event free probability if setting FALSE
<code>m</code>	desired minimum number of observations in a group
<code>g</code>	number of quantile groups
<code>cuts</code>	actual cuts in <code>x</code> , e.g. <code>c(0, 1, 2)</code> to use <code>[0,1)</code> , <code>[1,2]</code> .
<code>u</code>	time for which to estimate cumulative incidence
<code>pl</code>	TRUE to plot results
<code>conf.int</code>	defaults to .95 for 0.95 confidence bars. Set to FALSE to suppress bars
<code>xlab</code>	if <code>pl=TRUE</code> , is x-axis label. Default is <code>label(x)</code> or name of calling argument
<code>ylab</code>	if <code>pl=TRUE</code> , is y-axis label. Default is constructed from <code>u</code> and time units
<code>xlim</code>	range of x axis
<code>ylim</code>	range of y axis

lty	line type for primary line connecting estimates
add	set to TRUE if adding to an existing plot
cex.subtitle	character size for subtitle. Default is .7. Use FALSE to suppress subtitle.
ab	TRUE to add a 45 degree line
...	plotting parameters to pass to the plot and errbar functions

### Details

Function to divide a continuous variable  $x$  (e.g. age, or predicted cumulative incidence at time  $u$  created by `predict.cmprsk`) into  $g$  quantile groups, get cumulative incidence estimates at time  $u$  (a scalar), and to return a matrix with columns  $x$ =mean  $x$  in quantile,  $n$ =number of subjects,  $events$ =no. events, and  $ci$ = cumulative incidence at time  $u$ ,  $std.err$  = standard error. Instead of supplying  $g$ , the user can supply the minimum number of subjects to have in the quantile group ( $m$ , default=50). If  $cuts$  is given (e.g.  $cuts=c(0, .1, .2, \dots \{ \}, .9, .1)$ ), it overrides  $m$  and  $g$ .

### Value

matrix with columns named  $x$  (mean predictor value in interval),  $n$  (sample size in interval),  $events$  (number of events in interval),  $ci$  (cumulative incidence estimate),  $std.err$  (standard error of cumulative incidence)

### Note

This function is adapted from Harrell's function.

### Author(s)

Changhong Yu, Michael Kattan, Ph.D  
Department of Quantitative Health Sciences  
Cleveland Clinic

### See Also

[cuminc](#), [pred.ci](#)

### Examples

```
data(prostate.dat)
dd <- datadist(prostate.dat)
options(datadist = "dd")
prostate.f <- cph(Surv(TIME_EVENT, EVENT_DOD == 1) ~ TX + rcs(PSA, 3) +
  BX_GLSN_CAT + CLIN_STG + rcs(AGE, 3) +
  RACE_AA, data = prostate.dat,
  x = TRUE, y = TRUE, surv = TRUE, time.inc = 144)
prostate.crr <- crr.fit(prostate.f, cencode = 0, failcode = 1)

## ten fold cross validation
```

```

prostate.dat$preds.tenf.cv.prostate.crr.120 <-
  tenf.crr(prostate.crr,time = 120)

with(prostate.dat,
  groupci(preds.tenf.cv.prostate.crr.120 , ftime = TIME_EVENT,
    fstatus =EVENT_DOD, g = 5, u = 120,
    xlab = "Nomogram predicted 10-year cancerspecific mortality",
    ylab = "Observed predicted 10-year cancerspecific mortality")
)

```

---

Newlabels.cmprsk

*Change predictor labels for a model fit*


---

## Description

Change labels of predictors

## Usage

```

## S3 method for class 'cmprsk'
Newlabels(fit, labels, ...)

```

## Arguments

<code>fit</code>	a model fit
<code>labels</code>	a character vector specifying new labels for variables in a fit. To give new labels for all variables, you can specify labels of the form <code>labels=c("Age in Years","Cholesterol")</code> , where the list of new labels is assumed to be the length of all main effect-type variables in the fit and in their original order in the model formula. You may specify a named vector to give new labels in random order or for a subset of the variables, e.g., <code>labels=c(age="Age in Years",chol="Cholesterol")</code>
<code>...</code>	other arguments

## Details

This method function was written for competing risks regression model for facilitating to change the labels of predictors when construct a nomogram. It is used for the generic function [Newlabels](#)

## Value

a new model fit object with the levels adjusted.

**Examples**

```

data(prostate.dat)
dd <- datadist(prostate.dat)
options(datadist = "dd")
prostate.f <- cph(Surv(TIME_EVENT,EVENT_DOD == 1) ~ TX + rcs(PSA,3) +
  BX_GLSN_CAT + CLIN_STG + rcs(AGE,3) +
  RACE_AA, data = prostate.dat,
  x = TRUE, y= TRUE, surv=TRUE,time.inc = 144)
prostate.crr <- crr.fit(prostate.f,cencode = 0,failcode = 1)
prostate.g <- Newlabels(
  prostate.crr,
  c(
    TX = 'Treatment options',
    BX_GLSN_CAT = 'Biopsy Gleason Score Sum',
    CLIN_STG = 'Clinical stage'
  )
)

```

---

Newlevels.cmprsk

*Change levels of categorical variable for a model fit*


---

**Description**

```
##' Change levels of categorical variable
```

**Usage**

```
## S3 method for class 'cmprsk'
Newlevels(fit, levels, ...)
```

**Arguments**

fit	a model fit
levels	a list of named vectors specifying new level labels for categorical predictors. This will override parms as well as datadist information (if available) that were stored with the fit.
...	other arguments

**Details**

This method function was written for competing risks regression model for facilitating to change the levels of categorical predictors when construct a nomogram. It is used for the generic function [Newlevels](#)

**Value**

returns a new model fit object with the levels adjusted.

## Examples

```
data(prostate.dat)
dd <- datadist(prostate.dat)
options(datadist = "dd")
prostate.f <- cph(Surv(TIME_EVENT,EVENT_DOD == 1) ~ TX + rcs(PSA,3) +
  BX_GLSN_CAT + CLIN_STG + rcs(AGE,3) +
  RACE_AA, data = prostate.dat,
  x = TRUE, y= TRUE, surv=TRUE,time.inc = 144)
prostate.crr <- crr.fit(prostate.f,cencode = 0,failcode = 1)
prostate.g <- Newlevels(prostate.crr,
  list(TX=c('Treatment 1','Treatment 2', 'Treatment 3')))
```

---

nomo2.crr

*Estimate Cumulative Incidence Rate*


---

## Description

Calculate Estimated Cumulative Incidence Rate Calculate predicted cumulative incidence rate based on a competing risks regression model.

## Usage

```
nomo2.crr(x, f.crr, time)
```

## Arguments

x	a vector of sum of linear predictors for each subject.
f.crr	a saved model fitted by function <a href="#">crr.fit</a>
time	expected evaluation time

## Details

This function is usually used to transform regular failure probabilities to competing risks adjusted probabilities, when a nomogram of competing risks regression model is constructed started from a regular survival model. It is not often called externally.

## Value

a vector with each element being the predicted cumulative incidence rate at the expected time.

## Note

internal function

**Author(s)**

Michael W. Kattan, Ph.D. and Changhong Yu.  
Department of Quantitative Health Sciences, Cleveland Clinic

**See Also**

[pred2.crr](#) [crr.fit](#) [crr](#)

---

nomogram.crr

*Draw a Nomogram*

---

**Description**

Draws a partial nomogram that can be used to manually obtain predicted values from a regression model that was fitted with `rms` in effect.

**Usage**

```
nomogram.crr(  
  fit,  
  failtime = NULL,  
  ci = TRUE,  
  ...,  
  adj.to,  
  lp = TRUE,  
  lp.at,  
  lp.label = "Linear Predictor",  
  fun.at,  
  fun.lp.at,  
  fun.label = "Predicted Value",  
  fun.side,  
  interact = NULL,  
  intercept = 1,  
  conf.int = FALSE,  
  col.conf = c(1, 12),  
  conf.space = c(0.08, 0.2),  
  conf.lp = c("representative", "all", "none"),  
  est.all = TRUE,  
  abbrev = FALSE,  
  minlength = 4,  
  maxscale = 100,  
  nint = 10,  
  label.every = 1,  
  force.label = FALSE,  
  xfrac = 0.35,  
  cex.axis = 0.85,  
  cex.var = 1,
```

```

col.grid = FALSE,
vnames = c("labels", "names"),
varname.label = TRUE,
varname.label.sep = "=",
ia.space = 0.7,
tck = -0.009,
lmgp = 0.4,
omit = NULL,
naxes,
points.label = "Points",
total.points.label = "Total Points",
total.sep.page = FALSE,
total.fun,
verbose = FALSE,
total.min,
total.max,
mikeomit = NULL
)

```

### Arguments

<code>fit</code>	a competing risks regression model fit that was created with function <code>crr.fit</code> .
<code>failtime</code>	the expected failure time for calculating cumulative incidence.
<code>ci</code>	logical flag to output cumulative incidence or event free probability if setting FALSE.
<code>...</code>	settings of variables to use in constructing axes. If <code>datadist</code> was in effect, the default is to use <code>pretty(total range, nint)</code> for continuous variables, and the class levels for discrete ones. For <code>legend.nomabbrev, ...</code> specifies optional parameters to pass to <code>legend</code> . Common ones are <code>bty = "n"</code> to suppress drawing the box. You may want to specify a non-proportionally spaced font (e.g., <code>courier</code> ) number if abbreviations are more than one letter long. This will make the abbreviation definitions line up (e.g., specify <code>font = 2</code> , the default for <code>courier</code> ). Ignored for <code>print</code> .
<code>adj.to</code>	If you didn't define <code>datadist</code> for all predictors, you will have to define adjustment settings for the undefined ones, e.g. <code>adj.to=list(age=50, sex="female")</code> .
<code>lp</code>	Set to FALSE to suppress creation of an axis for scoring $X\beta$ .
<code>lp.at</code>	If <code>lp=TRUE</code> , <code>lp.at</code> may specify a vector of settings of $X\beta$ . Default is to use <code>pretty(range of linear predictors, nint)</code> .
<code>lplabel</code>	label for linear predictor axis. Default is "Linear Predictor".
<code>fun.at</code>	function values to label on axis. Default <code>fun</code> evaluated at <code>lp.at</code> . If more than one <code>fun</code> was specified, using a vector for <code>fun.at</code> will cause all functions to be evaluated at the same argument values. To use different values, specify a list of vectors for <code>fun.at</code> , with elements corresponding to the different functions (lists of vectors also applies to <code>fun.lp.at</code> and <code>fun.side</code> ).
<code>fun.lp.at</code>	If you want to evaluate one of the functions at a different set of linear predictor values than may have been used in constructing the linear predictor axis,

specify a vector or list of vectors of linear predictor values at which to evaluate the function. This is especially useful for discrete functions. The presence of this attribute also does away with the need for `nomogram` to compute numerical approximations of the inverse of the function. It also allows the user-supplied function to return factor objects, which is useful when e.g. a single tick mark position actually represents a range. If the `fun.lp.at` parameter is present, the `fun.at` vector for that function is ignored.

<code>funlabel</code>	label for fun axis. If more than one function was given but <code>funlabel</code> is of length one, it will be duplicated as needed. If <code>fun</code> is a list of functions for which you specified names (see the final example below), these names will be used as labels.
<code>fun.side</code>	a vector or list of vectors of side parameters for the axis function for labeling function values. Values may be 1 to position a tick mark label below the axis (the default), or 3 for above the axis. If for example an axis has 5 tick mark labels and the second and third will run into each other, specify <code>fun.side=c(1,1,3,1,1)</code> (assuming only one function is specified as <code>fun</code> ).
<code>interact</code>	When a continuous variable interacts with a discrete one, axes are constructed so that the continuous variable moves within the axis, and separate axes represent levels of interacting factors. For interactions between two continuous variables, all but the axis variable must have discrete levels defined in <code>interact</code> . For discrete interacting factors, you may specify levels to use in constructing the multiple axes. For continuous interacting factors, you must do this. Examples: <code>interact=list(age=seq(10,70,by=10),treat=c("A","B","D"))</code> .
<code>intercept</code>	for models such as the ordinal logistic model with multiple intercepts, specifies which one to use in evaluating the linear predictor.
<code>conf.int</code>	confidence levels to display for each scoring. Default is FALSE to display no confidence limits. Setting <code>conf.int</code> to TRUE is the same as setting it to <code>c(0.7,0.9)</code> , with the line segment between the 0.7 and 0.9 levels shaded using gray scale.
<code>col.conf</code>	colors corresponding to <code>conf.int</code> . Use fractions for gray scale (for UNIX S-PLUS).
<code>conf.space</code>	a 2-element vector with the vertical range within which to draw confidence bars, in units of <code>1=spacing</code> between main bars. Four heights are used within this range (8 for the linear predictor if more than 16 unique values were evaluated), cycling them among separate confidence intervals to reduce overlapping.
<code>conf.lp</code>	default is "representative" to group all linear predictors evaluated into deciles, and to show, for the linear predictor confidence intervals, only the mean linear predictor within the deciles along with the median standard error within the deciles. Set <code>conf.lp="none"</code> to suppress confidence limits for the linear predictors, and to "all" to show all confidence limits.
<code>est.all</code>	To plot axes for only the subset of variables named in <code>...{}</code> , set <code>est.all=FALSE</code> . Note: This option only works when zero has a special meaning for the variables that are omitted from the graph.
<code>abbrev</code>	Set to TRUE to use the <code>abbreviate</code> function to abbreviate levels of categorical factors, both for labeling tick marks and for axis titles. If you only want to abbreviate certain predictor variables, set <code>abbrev</code> to a vector of character strings containing their names.

<code>minlength</code>	applies if <code>abbrev=TRUE</code> . Is the minimum abbreviation length passed to the <code>abbreviate</code> function. If you set <code>minlength=1</code> , the letters of the alphabet are used to label tick marks for categorical predictors, and all letters are drawn no matter how close together they are. For labeling axes (interaction settings), <code>minlength=1</code> causes <code>minlength=4</code> to be used.
<code>maxscale</code>	default maximum point score is 100
<code>nint</code>	number of intervals to label for axes representing continuous variables. See <code>pretty</code> .
<code>label.every</code>	Specify <code>label.every=i</code> to label on every <i>i</i> th tick mark.
<code>force.label</code>	set to <code>TRUE</code> to force every tick mark intended to be labeled to have a label plotted (whether the labels run into each other or not)
<code>xfrac</code>	fraction of horizontal plot to set aside for axis titles
<code>cex.axis</code>	character size for tick mark labels
<code>cex.var</code>	character size for axis titles (variable names)
<code>col.grid</code>	If <code>col.grid=1</code> , no gray scale is used, but an ordinary line is drawn. If $0 < \text{col.grid} < 1$ , a <code>col</code> (gray scale) of <code>col.grid</code> is used to draw vertical reference lines for major axis divisions and <code>col.grid/2</code> for minor divisions. The default is <code>col.grid=FALSE</code> , i.e., reference lines are omitted. Specifying <code>col.grid=TRUE</code> is the same as specifying a gray scale level of <code>col.grid=.2</code> (5 for Windows S-PLUS).
<code>vnames</code>	By default, variable labels are used to label axes. Set <code>vnames="names"</code> to instead use variable names.
<code>varname.label</code>	In constructing axis titles for interactions, the default is to add <code>"(interacting.varname=level)"</code> on the right. Specify <code>varname.label=FALSE</code> to instead use <code>"(level)"</code> .
<code>varname.label.sep</code>	If <code>varname.label=TRUE</code> , you can change the separator to something other than <code>=</code> by specifying this parameter.
<code>ia.space</code>	When multiple axes are draw for levels of interacting factors, the default is to group combinations related to a main effect. This is done by spacing the axes for the second to last of these within a group only 0.7 (by default) of the way down as compared with normal space of 1 unit.
<code>tck</code>	see <code>tck</code> under <code>par</code>
<code>lmgp</code>	spacing between numeric axis labels and axis (see <code>par</code> for <code>mgp</code> )
<code>omit</code>	vector of character strings containing names of variables for which to suppress drawing axes. Default is to show all variables.
<code>naxes</code>	maximum number of axes to allow on one plot. If the nomogram requires more than one "page", the "Points" axis will be repeated at the top of each page when necessary.
<code>points.label</code>	a character string giving the axis label for the points scale
<code>total.points.label</code>	a character string giving the axis label for the total points scale
<code>total.sep.page</code>	set to <code>TRUE</code> to force the total points and later axes to be placed on a separate page

<code>total.fun</code>	a user-provided function that will be executed before the total points axis is drawn. Default is not to execute a function. This is useful e.g. when <code>total.sep.page=TRUE</code> and you wish to use <code>locator</code> to find the coordinates for positioning an abbreviation legend before it's too late and a new page is started (i.e., <code>total.fun=function()print(locator(1))</code> )
<code>verbose</code>	set to <code>TRUE</code> to get printed output detailing how tick marks are chosen and labeled for function axes. This is useful in seeing how certain linear predictor values cannot be solved for using inverse linear interpolation on the (requested linear predictor values, function values at these lp values). When this happens you will see NAs in the verbose output, and the corresponding tick marks will not appear in the nomogram.
<code>total.min</code>	Setting the minimal value in the total point axis on the nomogram.
<code>total.max</code>	Setting the maximal value in the total point axis.
<code>mikeomit</code>	The predictor variables specified by their names here will not be shown in the nomogram. The predicted outcome based on this reduced nomogram would be the same as if users were using the full version of the nomogram by entering the some values for the predictors remaining in the reduced nomogram but adjusted values for the hidden predictors so that 0 points will be achieved from these hidden predictor variables in the full nomogram.

### Details

The nomogram does not have lines representing sums, but it has a reference line for reading scoring points (default range 0–100). Once the reader manually totals the points, the predicted values can be read at the bottom. Non-monotonic transformations of continuous variables are handled (scales wrap around), as are transformations which have flat sections (tick marks are labeled with ranges).

### Value

a list of class "nomogram" that contains information used in plotting the axes. Please see [nomogram](#) for details.

### Note

internal use only

### Author(s)

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 Vanderbilt University  
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Draw a Competing Risks Nomogram

Draws a partial nomogram adjusting for competing risks for a cox ph survival model.

Changhong Yu, Michael Kattan, Ph.D  
 Department of Quantitative Health Sciences  
 Cleveland Clinic

## References

Banks J: Nomograms. Encyclopedia of Statistical Sciences, Vol 6. Editors: S Kotz and NL Johnson. New York: Wiley; 1985.

Lubsen J, Pool J, van der Does, E: A practical device for the application of a diagnostic or prognostic function. Meth. Inform. Med. 17:127–129; 1978.

Wikipedia: Nomogram, <https://en.wikipedia.org/wiki/Nomogram>.

Michael W. Kattan, Glenn Heller and Murray F. Brennan (2003). A competing-risks nomogram for sarcoma-specific death following local recurrence. Statistics in Medicine. Stat Med. 2003;22:3515-3525.

## See Also

[nomogram](#), [crr.fit](#), [pred2.crr](#), [nomo2.crr](#)

## Examples

```
data(prostate.dat)
dd <- datadist(prostate.dat)
options(datadist = "dd")
prostate.f <- cph(Surv(TIME_EVENT,EVENT_DOD == 1) ~ TX + rcs(PSA,3) +
  BX_GLSN_CAT + CLIN_STG + rcs(AGE,3) +
  RACE_AA, data = prostate.dat,
  x = TRUE, y= TRUE, surv=TRUE,time.inc = 144)
prostate.crr <- crr.fit(prostate.f,cencode = 0,failcode = 1)
## make a CRR nomogram
nomogram.crr(prostate.crr,failtime = 120,lp=FALSE,
  funlabel = "Predicted 10-year cumulative incidence")
```

---

nomogram.mk6

*Draw a Nomogram with modified function* Nomogram

---

## Description

Draw a Nomogram from a regression model

## Usage

```
nomogram.mk6(fit, ..., adj.to, lp = TRUE, lp.at,
  lplabel = "Linear Predictor",
  fun, fun.at, fun.lp.at, funlabel = "Predicted Value", fun.side,
  interact = NULL, intercept = 1, conf.int = FALSE,
  col.conf = c(1, 12), conf.space = c(0.08, 0.2),
  conf.lp = c("representative", "all", "none"), est.all = TRUE,
  abbrev = FALSE, minlength = 4, maxscale = 100, nint = 10,
  label.every = 1, force.label = FALSE, xfrac = 0.35, cex.axis = 0.85,
```

```

cex.var = 1, col.grid = NULL, vnames = c("labels", "names"),
varname.label = TRUE, varname.label.sep = "=", ia.space = 0.7,
tck = NA, tcl = -0.25, lmgp = 0.4, omit = NULL, naxes,
points.label = "Points",
total.points.label = "Total Points", total.sep.page = FALSE,
total.fun, verbose = FALSE, cap.labels = FALSE, total.min,
total.max, survtime, mikeomit = NULL)

```

## Arguments

<code>fit</code>	a regression model fit that was created with <code>library(rms)</code> in effect, and (usually) with <code>options(datadist = "object.name")</code> in effect.
<code>...</code>	settings of variables to use in constructing axes. If <code>datadist</code> was in effect, the default is to use <code>pretty(total range, nint)</code> for continuous variables, and the class levels for discrete ones. For <code>legend.nomabbrev</code> , <code>...</code> specifies optional parameters to pass to <code>legend</code> . Common ones are <code>bty = "n"</code> to suppress drawing the box. You may want to specify a non-proportionally spaced font (e.g., <code>courier</code> ) number if abbreviations are more than one letter long. This will make the abbreviation definitions line up (e.g., specify <code>font = 2</code> , the default for <code>courier</code> ). Ignored for <code>print</code> .
<code>adj.to</code>	If you didn't define <code>datadist</code> for all predictors, you will have to define adjustment settings for the undefined ones, e.g. <code>adj.to = list(age = 50, sex = "female")</code> .
<code>lp</code>	Set to <code>FALSE</code> to suppress creation of an axis for scoring $X\beta$ .
<code>lp.at</code>	If <code>lp=TRUE</code> , <code>lp.at</code> may specify a vector of settings of $X\beta$ . Default is to use <code>pretty(range of linear predictors, nint)</code> .
<code>lp.label</code>	label for linear predictor axis. Default is "Linear Predictor".
<code>fun</code>	on another axis. If more than one transformation is plotted, put them in a list, e.g. <code>list(function(x) x/2, function(x) 2*x)</code> . Any function values equal to <code>NA</code> will be ignored.
<code>fun.at</code>	function values to label on axis. Default <code>fun</code> evaluated at <code>lp.at</code> . If more than one <code>fun</code> was specified, using a vector for <code>fun.at</code> will cause all functions to be evaluated at the same argument values. To use different values, specify a list of vectors for <code>fun.at</code> , with elements corresponding to the different functions (lists of vectors also applies to <code>fun.lp.at</code> and <code>fun.side</code> ).
<code>fun.lp.at</code>	If you want to evaluate one of the functions at a different set of linear predictor values than may have been used in constructing the linear predictor axis, specify a vector or list of vectors of linear predictor values at which to evaluate the function. This is especially useful for discrete functions. The presence of this attribute also does away with the need for <code>nomogram</code> to compute numerical approximations of the inverse of the function. It also allows the user-supplied function to return factor objects, which is useful when e.g. a single tick mark position actually represents a range. If the <code>fun.lp.at</code> parameter is present, the <code>fun.at</code> vector for that function is ignored.
<code>funlabel</code>	label for <code>fun</code> axis. If more than one function was given but <code>funlabel</code> is of length one, it will be duplicated as needed. If <code>fun</code> is a list of functions for which

	you specified names (see the final example below), these names will be used as labels.
<code>fun.side</code>	a vector or list of vectors of side parameters for the <code>axis</code> function for labeling function values. Values may be 1 to position a tick mark label below the axis (the default), or 3 for above the axis. If for example an axis has 5 tick mark labels and the second and third will run into each other, specify <code>fun.side=c(1,1,3,1,1)</code> (assuming only one function is specified as <code>fun</code> ).
<code>interact</code>	When a continuous variable interacts with a discrete one, axes are constructed so that the continuous variable moves within the axis, and separate axes represent levels of interacting factors. For interactions between two continuous variables, all but the axis variable must have discrete levels defined in <code>interact</code> . For discrete interacting factors, you may specify levels to use in constructing the multiple axes. For continuous interacting factors, you must do this. Examples: <code>interact = list(age = seq(10,70,by=10), treat = c("A", "B", "D"))</code> .
<code>intercept</code>	for models such as the ordinal logistic model with multiple intercepts, specifies which one to use in evaluating the linear predictor.
<code>conf.int</code>	confidence levels to display for each scoring. Default is FALSE to display no confidence limits. Setting <code>conf.int</code> to TRUE is the same as setting it to <code>c(0.7, 0.9)</code> , with the line segment between the 0.7 and 0.9 levels shaded using gray scale.
<code>col.conf</code>	colors corresponding to <code>conf.int</code> . Use fractions for gray scale (for UNIX S-PLUS).
<code>conf.space</code>	a 2-element vector with the vertical range within which to draw confidence bars, in units of <code>l=spacing</code> between main bars. Four heights are used within this range (8 for the linear predictor if more than 16 unique values were evaluated), cycling them among separate confidence intervals to reduce overlapping
<code>conf.lp</code>	default is "representative" to group all linear predictors evaluated into deciles, and to show, for the linear predictor confidence intervals, only the mean linear predictor within the deciles along with the median standard error within the deciles. Set <code>conf.lp = "none"</code> to suppress confidence limits for the linear predictors, and to "all" to show all confidence limits.
<code>est.all</code>	To plot axes for only the subset of variables named in <code>...</code> , set <code>est.all = FALSE</code> . Note: This option only works when zero has a special meaning for the variables that are omitted from the graph
<code>abbrev</code>	Set to TRUE to use the <code>abbreviate</code> function to abbreviate levels of categorical factors, both for labeling tick marks and for axis titles. If you only want to abbreviate certain predictor variables, set <code>abbrev</code> to a vector of character strings containing their names.
<code>minlength</code>	<code>abbreviate</code> function. If you set <code>minlength = 1</code> , the letters of the alphabet are used to label tick marks for categorical predictors, and all letters are drawn no matter how close together they are. For labeling axes (interaction settings), <code>minlength = 1</code> causes <code>minlength = 4</code> to be used
<code>maxscale</code>	default maximum point score is 100
<code>nint</code>	number of intervals to label for axes representing continuous variables. See <code>pretty</code> .
<code>label.every</code>	Specify <code>label.every = i</code> to label on every <code>i</code> th tick mark

<code>force.label</code>	set to TRUE to force every tick mark intended to be labeled to have a label plotted (whether the labels run into each other or not)
<code>xfrac</code>	fraction of horizontal plot to set aside for axis titles
<code>cex.axis</code>	character size for tick mark labels
<code>cex.var</code>	character size for axis titles (variable names)
<code>col.grid</code>	If left unspecified, no vertical reference lines are drawn. Specify a vector of length one (to use the same color for both minor and major reference lines) or two (corresponding to the color for the major and minor divisions, respectively) containing colors, to cause vertical reference lines to the top points scale to be drawn. For R, a good choice is <code>col.grid = gray(c(0.8, 0.95))</code> .
<code>vnames</code>	By default, variable labels are used to label axes. Set <code>vnames = "names"</code> to instead use variable names.
<code>varname.label</code>	In constructing axis titles for interactions, the default is to add ( <code>interacting.varname = level</code> ) on the right. Specify <code>varname.label = FALSE</code> to instead use "(level)".
<code>varname.label.sep</code>	If <code>varname.label = TRUE</code> , you can change the separator to something other than = by specifying this parameter.
<code>ia.space</code>	When multiple axes are draw for levels of interacting factors, the default is to group combinations related to a main effect. This is done by spacing the axes for the second to last of these within a group only 0.7 (by default) of the way down as compared with normal space of 1 unit.
<code>tck</code>	see <code>tck</code> under <a href="#">par</a>
<code>tcl</code>	length of tick marks in nomogram
<code>lmgp</code>	spacing between numeric axis labels and axis (see <a href="#">par</a> for <code>mgp</code> )
<code>omit</code>	vector of character strings containing names of variables for which to suppress drawing axes. Default is to show all variables.
<code>naxes</code>	maximum number of axes to allow on one plot. If the nomogram requires more than one "page", the "Points" axis will be repeated at the top of each page when necessary.
<code>points.label</code>	a character string giving the axis label for the points scale
<code>total.points.label</code>	a character string giving the axis label for the total points scale
<code>total.sep.page</code>	set to TRUE to force the total points and later axes to be placed on a separate page
<code>total.fun</code>	a user-provided function that will be executed before the total points axis is drawn. Default is not to execute a function. This is useful e.g. when <code>total.sep.page = TRUE</code> and you wish to use <code>locator</code> to find the coordinates for positioning an abbreviation legend before it's too late and a new page is started (i.e., <code>total.fun = function() print(locator(1))</code> ).
<code>verbose</code>	set to TRUE to get printed output detailing how tick marks are chosen and labeled for function axes. This is useful in seeing how certain linear predictor values cannot be solved for using inverse linear interpolation on the (requested linear predictor values, function values at these lp values). When this happens you will see NAs in the verbose output, and the corresponding tick marks will not appear in the nomogram.

cap.labels	logical: should the factor labels have their first letter capitalized?
total.min	the minimum point for the total point axis
total.max	the maximum point for the total point axis
survtime	specified survival time for the predicted survival probability
mikeomit	a modified version of omit

### Details

a modified version of nomogram in rms package

### Value

a nomogram object

### Note

internal use only. please reference to [nomogram](#) details.

### References

Banks J: Nomograms. Encyclopedia of Statistical Sciences, Vol 6. Editors: S Kotz and NL Johnson. New York: Wiley; 1985.

Lubsen J, Pool J, van der Does, E: A practical device for the application of a diagnostic or prognostic function. Meth. Inform. Med. 17:127–129; 1978.

Wikipedia: Nomogram, <https://en.wikipedia.org/wiki/Nomogram>.

### See Also

[rms](#), [plot.summary.rms](#), [axis](#), [pretty](#), [approx](#)

### Examples

```
data(prostate.dat)
dd <- datadist(prostate.dat)
options(datadist = "dd")
prostate.f <- cph(Surv(TIME_EVENT, EVENT_DOD == 1) ~ TX + rcs(PSA, 3) +
  BX_GLSN_CAT + CLIN_STG + rcs(AGE, 3) +
  RACE_AA, data = prostate.dat,
  x = TRUE, y = TRUE, surv = TRUE, time.inc = 144)
## make a cph nomogram
nomogram.mk6(prostate.f, survtime=120, lp=FALSE,
funlabel = "Predicted 10-year cumulative incidence")
```

---

pred.ci                      *Calculate Cumulative Incidence*

---

**Description**

Calculate Cumulative Incidence

**Usage**

```
pred.ci(cum, tm1, failcode)
```

**Arguments**

cum	a object from function <a href="#">cuminc</a>
tm1	expected failure time
failcode	value indicating the event of interest

**Details**

Extract cumulative incidence and its variance from a object generated from function [cuminc](#).

**Value**

a data frame with 3 columns.

column 1:	Group name.
column 2:	Cumulative Incidence Probability.
column 3:	Variance

**Author(s)**

Michael W. Kattan, Ph.D. and Changhong Yu.  
Department of Quantitative Health Sciences, Cleveland Clinic

**See Also**

[cuminc](#)

**Examples**

```
data(prostate.dat) # get demo data set
cum <- cuminc(prostate.dat$TIME_EVENT,prostate.dat$EVENT_DOD,
              cencode = 0)
# calculate the expected cumulative incidence by 5 year for death from
# prostate cancer
# Here, code for cause A is 'DOA'.
pred.ci(cum,60,failcode = 1)
```

---

`pred2.crr`*Predict Cumulative Incidence Rate*

---

**Description**

Predict Cumulative Incidence Rate

**Usage**

```
pred2.crr(f.crr, lp, time)
```

**Arguments**

<code>f.crr</code>	a saved competing risks regression model created by function <a href="#">crr.fit</a>
<code>lp</code>	a scalar being the sum of linear predictors for a single subject.
<code>time</code>	expected time point, at which cumulative incidence rate will be assessed.

**Details**

Calculate the predicted cumulative incidence rate based on a saved competing risks regression model. The cumulative incidence is adjusted for other competing causes rather than the event of interest.

**Value**

Return the predicted cumulative incidence rate.

**Author(s)**

Michael Kattan, Ph.D, Changhong Yu  
Department of Quantitative Health Sciences  
Cleveland Clinic

**See Also**

[predict.crr](#) [crr.fit](#)

---

pred3.crr	<i>Predict cumulative incidence used internally</i>
-----------	---

---

**Description**

Predict cumulative incidence used internally an internal function.

**Usage**

```
pred3.crr(z, cov1, cov2, time, lps = FALSE)
```

**Arguments**

z	the fitter crr model
cov1	covariate matrix 1
cov2	covariate matrix 2
time	time point at which the prediction will make
lps	logical flag. if the liner predictor be generated.

**Details**

Internally used only.

**Value**

A list. See [crr](#) for details.

**Note**

an internal function called by [crr.fit](#).

**Author(s)**

changhong

---

predict.cmprsk      *Calculate Predicted Competing Risks Probability*

---

### Description

Calculate Predicted Competing Risks Probability.

### Usage

```
## S3 method for class 'cmprsk'  
predict(object, newdata = NULL, time,  
lps = FALSE, ...)
```

### Arguments

object	a saved crr model fit created by function <code>crr.fit</code>
newdata	data frame for prediction. Each row of the data frame contains values of covariates that are required in the crr model. If missing, the original data set that was used to develop the crr model will be used for prediction.
time	expected time point for evaluating the competing risks probability.
lps	set TRUE to return linear predictor values instead of failure probabilities.
...	other arguments

### Details

Calculate predicted probabilities for a competing risks regression model, which is fitted by function [crr.fit](#).

### Value

A vector with the length equal to the number of rows in the data frame, which was used to make prediction. Each element corresponds to a predicted failure probability at the expected time point.

### Note

This function is adapted from function [predict.crr](#) in package `cmprsk`.

### Author(s)

Michael W. Kattan, Ph.D. and Changhong Yu.  
Department of Quantitative Health Sciences, Cleveland Clinic

### References

Fine JP and Gray RJ (1999) A proportional hazards model for the subdistribution of a competing risk. *JASA* 94:496-509.

**See Also**

[crr.fit](#), [predict.crr](#)

**Examples**

```
data(prostate.dat)
dd <- datadist(prostate.dat)
options(datadist = "dd")
prostate.f <- cph(Surv(TIME_EVENT,EVENT_DOD == 1) ~ TX + rcs(PSA,3) +
  BX_GLSN_CAT + CLIN_STG + rcs(AGE,3) +
  RACE_AA, data = prostate.dat,
  x = TRUE, y= TRUE, surv=TRUE,time.inc = 144)
prostate.crr <- crr.fit(prostate.f,cencode = 0,failcode = 1)
prostate.dat$pred.60 <- predict(prostate.crr, time=60)
```

---

predictDesign

*Internal function to Calculate Predicted Competing Risks Probability.*

---

**Description**

Internal function to Calculate Predicted Competing Risks Probability.

**Usage**

```
predictDesign(
  fit,
  newdata = NULL,
  type = c("lp", "x", "data.frame", "terms", "adjto", "adjto.data.frame",
    "model.frame"),
  se.fit = FALSE,
  conf.int = FALSE,
  conf.type = c("mean", "individual"),
  incl.non.slopes = NULL,
  non.slopes = NULL,
  kint = 1,
  na.action = na.keep,
  expand.na = TRUE,
  center.terms = TRUE,
  ...
)
```

**Arguments**

<code>fit</code>	a saved crr model fit crated by function
<code>newdata</code>	data frame for prediction. Each row of the data frame contains values of covariates that are required in the crr model. If missing, the original data set that was used to develop the crr model will be used for prediction.
<code>type</code>	type of output desired
<code>se.fit</code>	return standard error
<code>conf.int</code>	specify <code>conf.int</code> as a positive fraction to obtain upper and lower confidence intervals (e.g., <code>'conf.int=0.95'</code> ).
<code>conf.type</code>	specifies the type of confidence interval.
<code>incl.non.slopes</code>	set to <code>'TRUE'</code> to include multiple intercepts, default is <code>'NULL'</code>
<code>non.slopes</code>	a vector of multiple intercepts
<code>kint</code>	a single integer specifying the number of the intercept to use in multiple-intercept models.
<code>na.action</code>	Function to handle missing values in <code>'newdata'</code> .
<code>expand.na</code>	set to <code>'FALSE'</code> to keep the naresid from having any effect
<code>center.terms</code>	set to <code>'FALSE'</code> to suppress subtracting adjust-to values from columns of the design matrix before computing terms with <code>'type="terms"'</code> .
<code>...</code>	ignored <code>crr.fit</code>

**Details**

Calculate predicted probabilities for a competing risks regression model

**Value**

A vector with the length equal to the number of rows in the data frame, which was used to make prediction. Each element corresponds to a predicted failure probability at the expected time point.

**See Also**

[predictrms](#),

---

prostate.dat

*Prostate cancer data set*

---

**Description**

This is an artificial prostate cancer dataset used for illustrating the usages of functions in R package and QHScrnomo

**Format**

A data frame with 2000 observations on the following 9 variables.

**UNIQID** patient ID

**TX** Treatment options of prostate cancer with levels EBRT, PI, RP

**PSA** Pre-treatment PSA levels

**BX\_GLSN\_CAT** Biopsy Gleason Score Sum. a factor with levels 1 for 2-6 2 for 7 and 3 for 8-10

**CLIN\_STG** Clinical stage with levels T1, T2, T3

**AGE** Age at treatment date

**RACE\_AA** patient ethnicity, a factor with levels 0 for other and 1 for African American

**TIME\_EVENT** follow up time in months

**EVENT\_DOD** followup status, 0 - censored, 1 - died of prostate cancer, 2 - died of other causes

**Details**

This is a simulated data set.

**Examples**

```
data(prostate.dat)
```

---

```
sas.cmprsk
```

```
generate prediction equation for a competing risks regression models
```

---

**Description**

Generate an equation to calculate X beta from a crr model fit. If specify a time point, the function also generates the subcumulative rate at the time point.

**Usage**

```
sas.cmprsk(f, time = NA, baseonly = FALSE, file = "", append = FALSE)
```

**Arguments**

<b>f</b>	a model fit from the competing risks regression.
<b>time</b>	time point
<b>baseonly</b>	logical variable. If true, only base survival probability will be printed.
<b>file</b>	A connection, or a character string naming the file to print to
<b>append</b>	logical. Only used if the argument file is the name of file. If TRUE output will be appended to file; otherwise, it will overwrite the content of file

**Details**

f should be fitted by the function [crr.fit](#)

**Value**

out a character vector that can be output as a formula by function `cat`  
 Rout same as out except replacing "max", "min", "=" and "\*\*" with "pmax", "pmin",  
 "==" and "^" respectively so that the formula can be pasted to R session and compute X beta directly without any further modification

**Author(s)**

changhong

**See Also**

[sascode](#), [Function](#), [crr.fit](#)

**Examples**

```
data(prostate.dat)
dd <- datadist(prostate.dat)
options(datadist = "dd")
prostate.f <- cph(Surv(TIME_EVENT, EVENT_DOD == 1) ~ TX + rcs(PSA, 3) +
  BX_GLSN_CAT + CLIN_STG + rcs(AGE, 3) +
  RACE_AA, data = prostate.dat,
  x = TRUE, y = TRUE, surv = TRUE, time.inc = 144)
prostate.crr <- crr.fit(prostate.f, cencode = 0, failcode = 1)
sas.cmprsk(prostate.crr, time = 60)
```

---

summary.cmprsk

*summary for competing risks regression*

---

**Description**

generate summary information

**Usage**

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

**Arguments**

object a crr model object from function [crr.fit](#)  
 ... other parameters

**Details**

summarize a crr regression model

**Value**

a matrix

**Author(s)**

changhong

**Examples**

```
data(prostate.dat)
dd <- datadist(prostate.dat)
options(datadist = "dd")
prostate.f <- cph(Surv(TIME_EVENT,EVENT_DOD == 1) ~ TX + rcs(PSA,3) +
  BX_GLSN_CAT + CLIN_STG + rcs(AGE,3) +
  RACE_AA, data = prostate.dat,
  x = TRUE, y= TRUE, surv=TRUE,time.inc = 144)
prostate.crr <- crr.fit(prostate.f,cencode = 0,failcode = 1)
summary(prostate.crr)
```

---

tenf.crr

*Ten fold cross validation for competing risks regression*


---

**Description**

Ten fold cross validation for crr endpoint

**Usage**

```
tenf.crr(fit, time = NA, lps = FALSE, fold = 10)
```

**Arguments**

fit	a competing risks regression model fitted by function <a href="#">crr.fit</a> .
time	the expected time point.
lps	logical flag. If true, values of predicted X beta will be output instead of cumulative incidence
fold	number of fold. the default is 10 fold cross validation.

**Details**

Do cross validation on a competing risk regression model.

**Value**

A vector of predicted values of cumulative incidence or X beta for each observation.

**Note**

Before the function is called, packages 'Hmisc', 'rms' and 'cmprsk' should be loaded as the function will call some functions in these packages.

**Author(s)**

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**See Also**

[crr.fit](#), [crr](#)

**Examples**

```
data(prostate.dat)
dd <- datadist(prostate.dat)
options(datadist = "dd")
prostate.f <- cph(Surv(TIME_EVENT,EVENT_DOD == 1) ~ TX + rcs(PSA,3) +
  BX_GLSN_CAT + CLIN_STG + rcs(AGE,3) +
  RACE_AA, data = prostate.dat,
  x = TRUE, y= TRUE, surv=TRUE,time.inc = 144)
prostate.crr <- crr.fit(prostate.f,cencode = 0,failcode = 1)

## ten fold cross validation
prostate.dat$preds.tenf<-
  tenf.crr(prostate.crr,time = 120)
```

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