

Package ‘quantregGrowth’

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Type Package

Title Non-Crossing Additive Regression Quantiles and Non-Parametric Growth Charts

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Maintainer Vito M. R. Muggeo <vito.muggeo@unipa.it>

Description Fits non-crossing regression quantiles as a function of linear covariates and multiple smooth terms, including varying coefficients, via B-splines with L1-norm difference penalties. Random intercepts and variable selection are allowed via the lasso penalties. The smoothing parameters are estimated as part of the model fitting, see Muggeo and others (2021) <[doi:10.1177/1471082X20929802](https://doi.org/10.1177/1471082X20929802)>. Monotonicity and concavity constraints on the fitted curves are allowed, see Muggeo and others (2013) <[doi:10.1007/s10651-012-0232-1](https://doi.org/10.1007/s10651-012-0232-1)>, and also <[doi:10.13140/RG.2.2.12924.85122](https://doi.org/10.13140/RG.2.2.12924.85122)> or <[doi:10.13140/RG.2.2.29306.21445](https://doi.org/10.13140/RG.2.2.29306.21445)> some code examples.

Depends R (>= 3.5.0), quantreg, splines, SparseM, methods

License GPL

Suggests knitr, rmarkdown, mgcv, markdown

VignetteBuilder knitr

NeedsCompilation no

Author Vito M. R. Muggeo [aut, cre] (ORCID: <<https://orcid.org/0000-0002-3386-4054>>)

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quantregGrowth-package

Non-Crossing Additive Regression Quantiles and Non-Parametric Growth Charts.

Description

Fits non-crossing regression quantiles as a function of linear covariates and smooth terms via P-splines with difference penalties. Random intercepts and selection of linear variables are allowed via the lasso penalties. Estimation of (possibly adaptive) smoothing/tuning parameters (for the spline terms, the random intercepts and the variables selector) are carried out efficiently as part of model fitting.

Details

Package: quantregGrowth
 Type: Package
 Version: 1.7-2
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 License: GPL

Package quantregGrowth allows estimation of growth charts via quantile regression. Given a set of percentiles (i.e. probability values), gcrq estimates non-crossing quantile curves as a flexible function of quantitative covariates (typically age in growth charts), and possibly additional linear terms. To ensure flexibility, B-splines with a difference L_1 penalty are employed to estimate non parametrically the curves wherein monotonicity and concavity constraints may be also set. Multiple smooth terms, including varying coefficients, are allowed and the amount of smoothness for each term is efficiently included in the model fitting algorithm, see Muggeo et al. (2021). plot.gcrq displays the fitted lines along with observations and pointwise confidence intervals.

Author(s)

Vito M.R. Muggeo

Maintainer: Vito M.R. Muggeo <vito.muggeo@unipa.it>

References

Muggeo VMR, Torretta F, Eilers PHC, Sciandra M, Attanasio M (2021). Multiple smoothing parameters selection in additive regression quantiles. *Statistical Modelling*, **21**, 428-448.

Muggeo VMR (2021). Additive Quantile regression with automatic smoothness selection: the R package `quantregGrowth`.

<https://www.researchgate.net/publication/350844895>

Muggeo VMR, Sciandra M, Tomasello A, Calvo S (2013). Estimating growth charts via nonparametric quantile regression: a practical framework with application in ecology. *Environ Ecol Stat*, **20**, 519-531.

Muggeo VMR (2018). Using the R package `quantregGrowth`: some examples.

<https://www.researchgate.net/publication/323573492>

Some references on growth charts (the first two papers employ the so-called LMS method)

Cole TJ, Green P (1992) Smoothing reference centile curves: the LMS method and penalized likelihood. *Statistics in Medicine* **11**, 1305-1319.

Rigby RA, Stasinopoulos DM (2004) Smooth centile curves for skew and kurtotic data modelled using the Box-Cox power exponential distribution. *Statistics in Medicine* **23**, 3053-3076.

Wei Y, Pere A, Koenker R, He X (2006) Quantile regression methods for reference growth charts. *Statistics in Medicine* **25**, 1369-1382.

Some references on regression quantiles

Koenker R (2005) *Quantile regression*. Cambridge University Press, Cambridge.

Cade BS, Noon BR (2003) A gentle introduction to quantile regression for ecologists. *Front Ecol Environ* **1**, 412-420.

See Also

[gcrq](#), [rq](#) in package `quantreg`

Examples

```
#see ?gcrq for some examples
```

charts

Easy computing growth charts

Description

Computes and returns quantiles as a function of the specified covariate values

Usage

```
charts(fit, k, file = NULL, digits=2, conf.level=0,
       dataframe=FALSE, transf=NULL, se.type=c("sandw","boot"), ...)
```

Arguments

<code>fit</code>	The object fit returned by <code>gcrq</code>
<code>k</code>	Numeric indicating the covariate values. If integer (and scalar, specified via <code>5L</code> , say), <code>k</code> equispaced values in the covariate range are taken.
<code>file</code>	If specified, the (path) file name wherein the returned matrix including the quantiles will be written via <code>write.csv()</code>
<code>digits</code>	Number of digits whereby the estimated quantiles are rounded.
<code>conf.level</code>	If larger than zero, the pointwise confidence intervals for the estimated quantiles are also returned. If <code>conf.level=0</code> the simple point estimates.
<code>dataframe</code>	Logical. If <code>TRUE</code> and <code>conf.level>0</code> a dataframe is returned having point estimate and confidence intervals collapsed. Otherwise a matrix having number of rows equal to the number of covariate values.
<code>transf</code>	An optional character string (with "y" as argument) meaning a function to apply to the predicted values. E.g. " $(\exp(y)-0.1)$ ". If <code>NULL</code> (default) it is taken as the inverse of function <code>transf</code> (*if*) supplied in <code>gcrq</code> . The standard errors (provided <code>se.fit=TRUE</code> has been set) are adjusted accordingly via the Delta method. See argument "transf" in <code>gcrq()</code> . If <code>transf</code> has been specified in <code>gcrq()</code> , use <code>transf="y"</code> to force predictions on the transformed scale, i.e. without back transforming.
<code>se.type</code>	Which covariance matrix should be used, provided that <code>conf.level>0</code> . See type in <code>predict.gcrq</code> .
<code>...</code>	Further arguments passed on to <code>write.csv()</code>

Details

This function is simply a wrapper for `predict.gcrq`

Value

A matrix having number of columns equal to the number of quantile curves and number of rows depending `k`

Note

`charts` just works with models having a single smooth term. See `predict.gcrq` when the model involves multiple covariates.

Author(s)

Vito Muggeo

See Also

`predict.gcrq`

Examples

```
## Not run:
charts(_fit_, k=1L) #prediction at the minimum of covariate
charts(_fit_, k=1) #prediction at covariate value 1.

charts(_fit_, k=10L)

## End(Not run)
```

gcrq	<i>Growth charts regression quantiles with automatic smoothness estimation</i>
------	--------------------------------------------------------------------------------

Description

Modelling unspecified nonlinear relationships between covariates and quantiles of the response conditional distribution. Typical example is estimation of nonparametric growth charts (via quantile regression). Quantile curves are estimated via B-splines with a L_1 penalty on the spline coefficient differences, while non-crossing and possible monotonicity and concavity restrictions are set to obtain estimates more biologically plausible. Linear terms can be specified in the model formula. Multiple smooth terms, including varying coefficients, with automatic selection of corresponding smoothing parameters are allowed.

Usage

```
gcrq(formula, tau=c(.1,.25,.5,.75,.9), data, subset, weights, na.action,
      transf=NULL, y=TRUE, n.boot=0, eps=0.0001, display=FALSE,
      method=c("REML","ML"), df.opt=2, df.nc=FALSE, lambda0=.1, h=0.8, lambda.max=1000,
      tol=0.01, it.max=15, single.lambda=TRUE, foldid=NULL, nfolds=10,
      lambda.ridge=0, adjX.constr=TRUE, contrasts=NULL, sparse=FALSE, nc.fit=FALSE, ...)
```

Arguments

formula	a standard R formula to specify the response in the left hand side, and the covariates in the right hand side, such as $y \sim \psi(x) + z$, see Details for further examples.
tau	a numeric vector to specify the quantile curves of interest. Default to probability values (.1, .25, .5, .75, .9).
data	the dataframe where the variables required by the formula, subset and weights arguments are stored.
subset	optional. A vector specifying a subset of observations to be used in the fitting process.
weights	optional. A numeric vector specifying weights to be assigned to the observations in the fitting process. Currently unimplemented.
na.action	a function which indicates how the possible 'NA's are handled.

<code>transf</code>	an optional character string (with "y" as argument) meaning a function to apply to the response variable before fitting. It can be useful to guarantee fitted values within a specified range; e.g. if $y \geq 0$, we could set " $\log(y+0.1)$ ". If provided, the resulting object fit refers to the model for the transformed response and it will include the corresponding inverse function (numerically computed) to be used to back transform predictions (see argument <code>transf</code> in <code>predict.gcrq</code> and <code>plot.gcrq</code>).
<code>y</code>	logical. If TRUE (default) the returned object includes also the responses vector.
<code>n.boot</code>	Number of nonparametric (cases resampling) bootstrap samples to be used. If $n.boot > 0$, the covariance matrix can be obtained as empirical covariance matrix of the bootstrap distributions, see <code>vcov.gcrq</code> . Notice that the smoothing parameter (if relevant) is assumed fixed. Namely it does change throughout the bootstrap replicates.
<code>eps</code>	A small positive constant to ensure noncrossing curves (i.e. the minimum distance between two consecutive curves). Use it at your risk! If <code>eps</code> is large, the resulting fitted quantile curves could appear unreasonable.
<code>display</code>	Logical. Should the iterative process be printed? Ignored if no smooth is specified in the formula or if all the smoothing parameters specified in <code>ps</code> terms are fixed.
<code>method</code>	character, "ML" or "REML" affecting the smoothing parameter estimation. Default is "REML" which appears to provide better performance in simulation studies. Ignored if no smoothing parameter has to be estimated.
<code>df.opt</code>	How the model and term-specific degrees of freedom are computed. <code>df.opt=1</code> means via the null penalized coefficients, and <code>df.opt=2</code> via the trace of the approximate hat matrix. Ignored if no smoothing parameter is to be estimated.
<code>df.nc</code>	logical. If TRUE and the model refers to multiple quantile curves, the degrees of freedom account for the noncrossing constraints. Ignored for single quantile fits. Default to FALSE, as it is still experimental.
<code>lambda0</code>	the starting value for the lambdas to be estimated. Ignored if all the smoothing parameters specified in <code>ps</code> terms are fixed.
<code>h</code>	The step halving factor affecting estimation of the smoothing parameters. Lower values lead to slower updates in the lambda values. Ignored if all the smoothing parameters specified in <code>ps</code> terms are fixed.
<code>lambda.max</code>	The upper bound for lambda estimation. Ignored if all the smoothing parameters specified in <code>ps</code> terms are fixed.
<code>tol</code>	The tolerance value to declare convergence. Ignored if all the smoothing parameters specified in <code>ps</code> terms are fixed.
<code>it.max</code>	The maximum number of iterations in lambdas estimation. Ignored if all the smoothing parameters specified in <code>ps</code> terms are fixed.
<code>single.lambda</code>	Logical. Should the smoothing parameter (for each smooth term) to be the same across the quantile curves being estimated? Ignored when just a single quantile curve is being estimated.
<code>foldid</code>	optional. A numeric vector identifying the group labels to perform cross validation to select the smoothing parameter. Ignored if the <code>lambda</code> argument in <code>ps()</code> is not a vector.

<code>nfolds</code>	optional. If <code>foldid</code> is not provided, it is scalar specifying the number of ‘folds’ (groups) which should be used to perform cross validation to select the smoothing parameter. Default to 10, but it is ignored if the <code>lambda</code> argument in <code>ps()</code> is not a vector.
<code>lambda.ridge</code>	Numerical value (typically very small) to stabilize model estimation.
<code>adjX.constr</code>	logical. If TRUE, each linear covariate is shifted (by subtracting its min) in order to set up effective constraints to prevent crossing. It does not need to be TRUE, if the noncrossing constraints are set on the fitted values (i.e. <code>nc.fit=TRUE</code>). Ignored if the model does not include linear terms or <code>tau</code> is a scalar.
<code>contrasts</code>	an optional list. See argument <code>contrasts.arg</code> in <code>model.matrix.default</code> .
<code>sparse</code>	logical. If TRUE, the model is fitted via sparse algebra as implemented in the SparseM package. Typically <code>sparse=TRUE</code> is used when the model involves a single smooth with a very rich basis and a large sample size, see Details.
<code>nc.fit</code>	If TRUE, the noncrossing constraints are set directly on the fitted values. Useful if the sample size is not large and there are several linear covariates.
<code>...</code>	If <code>sparse=TRUE</code> , these are further arguments to be passed to the <code>rq.fit.sfn()</code> or <code>rq.fit.sfn()</code> for sparse fits. Eg it could be <code>control=list(maxiter=300)</code>

Details

The function fits regression quantiles at probability values `tau` as a function of covariates specified in the formula argument. The formula may include linear terms and one or several `ps` terms to account for nonlinear relationships with quantitative covariates, usually age in growth charts. When the `lambda` argument in `ps()` is a negative scalar, the smoothing parameter is estimated iteratively as discussed in Muggeo et al. (2021). If a positive scalar, it represents the actual smoothing parameter value.

When the model includes a single `ps` term, setting `sparse=TRUE` (introduced since version 1.7-0) could reduce the computational time especially when the sample size is large and the basis involves several terms. However when `sparse=TRUE` is set, no linear term is allowed and for the smooth term a full and uncentred basis is used (i.e., equivalent to setting `dropc=FALSE` and `center=FALSE` along with `constr.fit=FALSE` in `ps()`). Therefore the correct call would be

```
gcrq(y ~ 0+ps(x), sparse=TRUE) #if y~ps(x), a warning is printed as the intercept is not explicitly removed
```

which fits the same model as

```
gcrq(y ~ 0+ps(x, dropc=FALSE, center=FALSE)) #sparse=FALSE is the default
```

Smoothing parameter selection via ‘K-fold’ cross validation (CV) is also allowed (but not recommended) if the model includes a single `ps` term: `lambda` should be a vector of candidate values, and the final fit is returned at the ‘optimal’ `lambda` value. To select the smoothing parameter via CV, `foldid` or `nfolds` may be supplied. If provided, `foldid` overwrites `nfolds`, otherwise `foldid` is obtained via random extraction, namely `sample(rep(seq(nfolds), length = n))`. However selection of smoothing parameter via CV is allowed only with a unique `ps` term in the formula.

Value

This function returns an object of class `gcrq`, that is a list with the following components (only the most important are listed)

<code>coefficients</code>	The matrix of estimated regression parameters; the number of columns equals the number of the fitted quantile curves.
<code>x</code>	the design matrix of the final fit (including the dummy rows used by penalty).
<code>edf.j</code>	a matrix reporting the edf values for each term at each quantile curve. See the section 'Warning' below.
<code>rho</code>	a vector including the values of the objective functions at the solution for each quantile curve.
<code>fitted.values</code>	a matrix of fitted quantiles (a column for each tau value)
<code>residuals</code>	a matrix of residuals (a column for each tau value)
<code>D.matrix</code>	the penalty matrix (multiplied by the smoothing parameter value).
<code>D.matrix.nolambda</code>	the penalty matrix.
<code>pLin</code>	number of linear covariates in the model.
<code>info.smooth</code>	some information on the smoothing term (if included in the formula via <code>ps</code>).
<code>BB</code>	further information on the smoothing term (if present in the formula via <code>ps</code>), including stuff useful for plotting via <code>plot.gcrq()</code> .
<code>Bderiv</code>	if the smooth term is included, the first derivative of the B spline basis.
<code>boot.coef</code>	The array including the estimated coefficients at different bootstrap samples (provided that <code>n.boot>0</code> has been set).
<code>y</code>	the response vector (if <code>gcrq()</code> has been called with <code>y=TRUE</code>).
<code>contrasts</code>	the contrasts used, when the model contains a factor.
<code>xlevels</code>	the levels of the factors (when included) used in fitting.
<code>taus</code>	a vector of values between 0 and 1 indicating the estimated quantile curves.
<code>call</code>	the matched call.

Warning

Variable selection is still at an experimental stage but it works :-)

When including linear terms, the covariates X_j are shifted such that $\min(X_j) = 0$, therefore the model intercept should be interpreted accordingly.

The options 'REML' or 'ML' of the argument `method`, refer to how the degrees of freedom are computed to update the lambda estimates.

Currently, standard errors are obtained via the sandwich formula or the nonparametric bootstrap (case resampling). Both methods ignore uncertainty in the smoothing parameter selection.

Since version 1.2-1, computation of the approximate edf can account for the noncrossing constraints by specifying `df.nc=TRUE`. That could affect model estimation when the smoothing parameter(s) have to be estimated, because the term specific edf are used to update the lambda value(s). When lambda is not being estimated (it is fixed or there is no `ps` term in the formula), parameter estimate

is independent of the `df.nc` value. The `summary.gcrq` method reports if the edf account for the noncrossing constraints.

Using `ps(..., center=TRUE)` in the formula leads to lower uncertainty in the fitted curve while guaranteeing noncrossing constraints.

Currently, decomposition of Bsplines (i.e. `ps(..., decom=TRUE)`) is incompatible with shape (monotonicity and concavity) restrictions and even with noncrossing constraints.

Note

This function is based upon the package `quantreg` by R. Koenker. Currently methods specific to the class "gcrq" are `print.gcrq`, `summary.gcrq`, `vcov.gcrq`, `plot.gcrq`, `predict.gcrq`, `AIC.gcrq`, and `logLik.gcrq`.

If the sample is not large, and/or the basis rank is large (i.e. a large number of columns) and/or there are relatively few distinct values in the covariate distribution, the fitting algorithm may fail returning error messages like the following

```
> Error info = 20 in stepy2: singular design
```

To remedy it, it suffices to change some arguments in `ps()`: to decrease `ndx` or `deg` (even by a small amount) or to increase (even by a small amount) the `lambda` value. Sometimes even by changing slightly the `tau` probability value (for instance from 0.80 to 0.79) can bypass the aforementioned errors.

Author(s)

Vito M. R. Muggeo, <vito.muggeo@unipa.it>

References

V.M.R. Muggeo, F. Torretta, P.H.C. Eilers, M. Sciandra, M. Attanasio (2021). Multiple smoothing parameters selection in additive regression quantiles, *Statistical Modelling*, 21: 428-448.

V. M. R. Muggeo (2021). Additive Quantile regression with automatic smoothness selection: the R package `quantregGrowth`. <https://www.researchgate.net/publication/350844895>

V. M. R. Muggeo, M. Sciandra, A. Tomasello, S. Calvo (2013). Estimating growth charts via nonparametric quantile regression: a practical framework with application in ecology, *Environ Ecol Stat*, 20, 519-531.

V. M. R. Muggeo (2018). Using the R package `quantregGrowth`: some examples. <https://www.researchgate.net/publication/323573492>

See Also

[ps](#), [plot.gcrq](#), [predict.gcrq](#)

Examples

```
## Not run:
##=== Example 1: an additive model from ?mgcv::gam

d<-mgcv::gamSim(n=200, eg=1)
```

```

o<-gcrq(y ~ ps(x0) + ps(x1)+ ps(x2) + ps(x3), data=d, tau=.5, n.boot=50)
plot(o, res=TRUE, col=2, conf.level=.9, shade=TRUE, split=TRUE)

###== Example 2: some simple examples involving just a single smooth

data(growthData) #load data
tauss<-seq(.1,.9,by=.1) #fix the percentiles of interest

m1<-gcrq(y~ps(x), tau=tauss, data=growthData) #lambda estimated..

m2<-gcrq(y~ps(x, lambda=0), tau=tauss, data=growthData) #unpenalized.. very wiggly curves
#strongly penalized models
m3<-gcrq(y~ps(x, lambda=1000, d=2), tau=tauss, data=growthData) #linear
m4<-gcrq(y~ps(x, lambda=1000, d=3), tau=tauss, data=growthData) #quadratic

#penalized model with monotonicity restrictions
m5<-gcrq(y~ps(x, monotone=1, lambda=10), tau=tauss, data=growthData)

#monotonicity constraints,lambda estimated, and varying penalty
m6<-gcrq(y~ps(x, monotone=1, lambda=10, var.pen="(1:k)"), tau=tauss, data=growthData)
m6a<-gcrq(y~ps(x, monotone=1, lambda=10, var.pen="(1:k)^2"), tau=tauss, data=growthData)

par(mfrow=c(2,3))
plot(m1, res=TRUE, col=-1)
plot(m2, pch=20, res=TRUE)
plot(m3, add=TRUE, lty=2, col=4)
plot(m4, pch=20, res=TRUE)
plot(m5, pch=4, res=TRUE, legend=TRUE, col=2)
plot(m6, lwd=2, col=3)
plot(m6a, lwd=2, col=4)

#select lambda via 'K-fold' CV (only with a single smooth term)
m7<-gcrq(y~ps(x, lambda=seq(0,10,l=20)), tau=tauss, data=growthData)
par(mfrow=c(1,2))
plot(m7, cv=TRUE) #display CV score versus lambda values
plot(m7, res=TRUE, grid=list(x=5, y=8), col=4) #fit at the best lambda (by CV)

###== Example 3: VC models

n=50
x<-1:n/n
mu0<-10+sin(2*pi*x)
mu1<- 7+4*x
y<-c(mu0,mu1)+rnorm(2*n)*.2 #small noise.. just to illustrate..
x<-c(x,x)
z<-rep(0:1, each=n)

# approach 1: a smooth in each *factor* level
g<-factor(z)
o <-gcrq(y~ g+ps(x,by=g), tau=.5)

```

```

predict(o, newdata=data.frame(x=c(.3,.7), g=factor(c(0,1))))
par(mfrow=c(1,2))
plot(x[1:50],mu0,type="l")
plot(o, term=1, add=TRUE)
points(c(.3,.7), predict(o, newdata=data.frame(x=c(.3,.7), g=factor(c(0,0)))), pch=4, lwd=3,col=2)

plot(x[1:50],mu1,type="l")
plot(o, term=2, add=T, shift=coef(o)["g1",], col=3) #note the argument 'shift'
points(c(.3,.7), predict(o, newdata=data.frame(x=c(.3,.7), g=factor(c(1,1)))), pch=4, lwd=3,col=3)

# approach 2: a general smooth plus the (smooth) 'interaction' with a continuous covariate..
b1 <-gcrq(y~ ps(x) + z+ ps(x,by=z), tau=.5)
par(mfrow=c(1,2))
plot(x[1:50],mu0,type="l")
plot(b1, add=TRUE, term=1) #plot the 1st smooth term

#plot the 2nd smooth of 'b1' (which is actually the difference mu1-mu0)
plot(x[1:50], mu1-mu0, type="l")
plot(b1, term=2, add=TRUE, interc=FALSE, shift=coef(b1)["z",])

###== Example 4: random intercepts example

n=50
x<-1:n/n

set.seed(69)
z<- sample(1:15, size=n, replace=TRUE)
#table(z)

#true model: linear effect + 3 non-null coeffs when z= 3, 7, and 13
y<-2*x+ I(z==3)- I(z==7) + 2*I(z==13) + rnorm(n)*.15
id<-factor(z)

o <-gcrq(y~x+ps(id), tau=.5)
plot(o, term=1) #plot the subject-specific intercepts

#== variable selection
n=50
p=30
p1<-10

X<-matrix(rnorm(n*p,5),n,p)
b<-rep(0,p)
id<-sample(1:p, p1)
b[id]<-round(runif(p1,.5,2),2)
b[id]<-b[id]* sign(ifelse(runif(p1)<.5,1,-1))
lp <- drop(tcrossprod(X,t(b)))
y <- 2+lp+rnorm(n)*1.5

```

```
gcrq(y~ps(X), tau=.7)
## End(Not run)
```

growthData	<i>Simulated data to illustrate capabilities of the package</i>
------------	-----------------------------------------------------------------

Description

The growthData data frame has 200 rows and 3 columns.

Usage

```
data(growthData)
```

Format

A data frame with 200 observations on the following 3 variables.

- x the supposed 'age' variable.
- y the supposed growth variable (e.g. weight).
- z an additional variable to be considered in the model.

Details

Simulated data to illustrate capabilities of the package.

Examples

```
data(growthData)
with(growthData, plot(x,y))
```

logLik.gcrq	<i>Log Likelihood, AIC and BIC for gcrq objects</i>
-------------	-----------------------------------------------------

Description

The function returns the log-likelihood value(s) evaluated at the estimated coefficients

Usage

```
## S3 method for class 'gcrq'
logLik(object, summ=TRUE, ...)
## S3 method for class 'gcrq'
AIC(object, ..., k=2, bondell=FALSE)
```

Arguments

object	A gcrq fit returned by gcrq()
summ	If TRUE, the log likelihood values (and relevant edf) are summed over the different taus to provide a unique value accounting for the different quantile curves. If FALSE, tau-specific values are returned.
k	Optional numeric specifying the penalty of the edf in the AIC formula. $k < 0$ means $k = \log(n)$.
bondell	Logical. If TRUE, the <i>SIC</i> according to formula (7) in Bondell et al. (2010) is computed.
...	optional arguments (nothing in logLik.gcrq). For AIC.gcrq, summ=TRUE or FALSE can be set.

Details

The 'logLikelihood' is computed by assuming an asymmetric Laplace distribution for the response as in [logLik.rq](#), namely $n(\log(\tau(1 - \tau)) - 1 - \log(\rho_\tau/n))$, where ρ_τ is the minimized objective function. When there are multiple quantile curves $j = 1, 2, \dots, J$ (and summ=TRUE) the formula is $n(\sum_j \log(\tau_j(1 - \tau_j)) - J - \log(\sum_j \rho_{\tau_j}/(nJ)))$

AIC.gcrq simply returns $-2 * \logLik + k * edf$ where k is 2 or $\log(n)$.

Value

The log likelihood(s) of the model fit object

Author(s)

Vito Muggeo

References

Bondell HD, Reich BJ, Wang H (2010) Non-crossing quantile regression curve estimation, *Biometrika*, 97: 825-838.

See Also

[logLik.rq](#)

Examples

```
## logLik(o) #a unique value (o is the fit object from gcrq)
## logLik(o, summ=FALSE) #vector of the log likelihood values
## AIC(o, k=-1) #BIC
```

ncross.rq.fitXB	<i>Estimation of noncrossing regression quantiles with monotonicity restrictions.</i>
-----------------	---------------------------------------------------------------------------------------

Description

These are internal functions of package `quantregGrowth` and should be not called by the user.

Usage

```
ncross.rq.fitXB(y, x, B=NULL, X=NULL, taus, monotone=FALSE, concave=FALSE,
  nomiBy=NULL, byVariabili=NULL, ndx=10, deg=3, dif=3, lambda=0, eps=.0001,
  var.pen=NULL, penMatrix=NULL, lambda.ridge=0, dropcList=FALSE,
  decomList=FALSE, vcList=FALSE, dropvcList=FALSE, centerList=FALSE,
  ridgeList=FALSE, ps.matrix.list=FALSE, colmeansB=NULL, Bconstr=NULL,
  adjX.constr=TRUE, adList=FALSE, it.j=10, myeps=NULL, nc=TRUE, lassoList=FALSE, ...)
```

```
ncross.rq.fitXBsparse(y, x, B=NULL, X=NULL, taus, monotone=FALSE, concave=FALSE,
  nomiBy=NULL, byVariabili=NULL, ndx=10, deg=3, dif=3, lambda=0, eps=.0001,
  var.pen=NULL, penMatrix=NULL, lambda.ridge=0, dropcList=FALSE, decomList=FALSE,
  vcList=FALSE, dropvcList=FALSE, centerList=FALSE, ridgeList=FALSE,
  ps.matrix.list=FALSE, colmeansB=NULL, Bconstr=NULL, adjX.constr=TRUE,
  adList=FALSE, it.j=10, myeps=NULL, nc=TRUE, ...)
```

```
ncross.rq.fitX(y, X = NULL, taus, adjX.constr=TRUE, lambda.ridge = 0,
  eps = 1e-04, sparse=FALSE, nc.fit=FALSE, ...)
```

```
gcrq.rq.cv(y, B, X, taus, monotone, concave, ndx, lambda, deg, dif, var.pen=NULL,
  penMatrix=NULL, lambda.ridge=0, dropcList=FALSE, decomList=FALSE,
  vcList=vcList, dropvcList=FALSE, nfold=10, foldid=NULL, eps=.0001,
  sparse=FALSE, ...)
```

Arguments

<code>y</code>	the responses vector. see gcrq
<code>x</code>	the covariate supposed to have a nonlinear relationship.
<code>B</code>	the B-spline basis.
<code>X</code>	the design matrix for the linear parameters.
<code>taus</code>	the percentiles of interest.
<code>monotone</code>	numerical value (-1/0/+1) to define a non-increasing, unconstrained, and non-decreasing flexible fit, respectively.
<code>concave</code>	numerical value (-1/0/+1) to possibly define concave or convex fits.
<code>nomiBy</code>	useful for VC models (when B is not provided).

byVariabili	useful for VC models (when B is not provided).
ndx	number of internal intervals within the covariate range, see ndx in ps .
deg	spline degree, see ps .
dif	difference order of the spline coefficients in the penalty term.
lambda	smoothing parameter value(s), see lambda in ps .
eps	tolerance value.
var.pen	Varying penalty, see ps .
penMatrix	Specified penalty matrix, see pen.matrix in ps .
lambda.ridge	a (typically very small) value, see lambda.ridge gcrq .
dropcList	see dropc in ps .
decomList	see decompose in ps .
vcList	to indicate if the smooth is VC or not, see by in ps .
dropvcList	see ps .
centerList	see center in ps .
ridgeList	see ridge in ps .
ps.matrix.list	nothing relevant for the user.
colmeansB	see center in ps .
Bconstr	see constr.fit in ps .
foldid	vector (optional) to perform cross validation, see the same arguments in gcrq .
nfolds	number of folds for crossvalidation, see the same arguments in gcrq .
cv	returning cv scores; see the same arguments in gcrq .
adjX.constr	logical to shift the linear covariates. Appropriate only with linear terms.
adList	see ad in ps .
it.j	Ignore.
myeps	Ignore.
sparse	logical, meaning if sparse computations have to be used.
lassoList	nothing relevant for the user.
nc	logical, nothing relevant for the user.
nc.fit	logical, nothing relevant for the user.
...	optional.

Details

These functions are called by [gcrq](#) to fit growth charts based on regression quantiles with non-crossing and monotonicity restrictions. The computational methods are based on the package [quantreg](#) by R. Koenker and details are described in the reference paper.

Value

A list of fit information.

Author(s)

Vito M. R. Muggeo

See Also[gcrq](#)**Examples**

##See ?gcrq

`plot.gcrq`*Plot method for gcrq objects*

Description

Displaying the estimated growth charts from a gcrq fit.

Usage

```
## S3 method for class 'gcrq'
plot(x, term=NULL, add = FALSE, res = FALSE, conf.level=0, axis.tau=FALSE,
     interc=TRUE, se.interc=FALSE, legend = FALSE, select.tau, deriv = FALSE,
     cv = FALSE, transf=NULL, lambda0=FALSE, shade=FALSE, overlap=NULL, rug=FALSE,
     overall.eff=TRUE, grid=NULL, smoos=NULL, split=FALSE, shift=0, type=c("sandw","boot"),
     n.points=NULL, ...)
```

Arguments

<code>x</code>	a fitted "gcrq" object.
<code>term</code>	the variable name or its index in the formula entering the model. Can be vector. Both linear and spline terms (i.e. included in the model via <code>ps</code>) can be specified and relevant fitted quantile curves (as optionally specified by <code>select.tau</code>) will be plotted. If the model includes both linear and smooth terms, the smooth terms are counted and drawn <i>*first*</i> : therefore if the model formula is $y \sim z + ps(x)$, <code>term=1</code> refers to the smooth term. If <code>NULL</code> , all smooth terms are plotted according to the <code>split</code> argument. If the model includes multiple quantile curves and <code>axis.tau=TRUE</code> , <code>term=1</code> refers to the model intercept (if in the model). The variable name should be reported within <code>'ps()'</code> , e.g. <code>'ps(age)'</code> regardless of additional arguments specified in <code>ps</code> .
<code>interc</code>	Should the smooth term be plotted along with the model intercept (provided it is included in the model)? Of course such argument is ignored if the smooth term has been called via <code>ps(, dropc=FALSE)</code> and the plot always includes implicitly the 'intercept'. Note that <code>interc=TRUE</code> is requested to display the noncrossing curves (if multiple quantile curves are being plotted).

se.interc	logical. If TRUE the standard errors of fitted quantile curves account for uncertainty of the model intercept (provided it is included in the model). If FALSE, then the uncertainty relates purely to the (usually centred) smooth itself. Ignored if <code>conf.level=0</code> .
add	logical. If TRUE the fitted quantile curves are added on the current plot.
res	logical. If TRUE 'partial residuals' are also displayed on the plot. Borrowing terminology from GLM, partial residuals for covariate X_j are defined as fitted values corresponding to X_j + residuals (from the actual fit). If there is a single covariate, the partial residuals correspond to observed data. If multiple quantile curves have been estimated, the fitted values coming from the 'middle' quantile curve are employed to compute the partial residuals. 'Middle' means 'corresponding to the τ_k closest to 0.50'. I don't know if that is the best choice.
conf.level	logical. If larger than zero, pointwise confidence intervals for the fitted quantile curve are also shown (at the confidence level specified by <code>conf.level</code>). Such confidence intervals are independent of the possible intercept accounted for via the <code>intercept</code> argument. See <code>type</code> to select different methods (bootstrap or sandwich) to compute the standard errors.
axis.tau	logical. If TRUE, the estimated coefficient term is plotted against the probability values. This graph could be useful if the model has been estimated at several tau values.
legend	logical. If TRUE a legend is drawn on on the right side of the plot.
select.tau	an optional numeric vector to draw only some of the fitted quantiles. Percentile values or integers 1 to <code>length(tau)</code> may be supplied.
deriv	logical. If TRUE the first derivative of the fitted curves are displayed.
cv	logical. If TRUE and the "gcrq" object contains a single smooth term wherein lambda has been selected via CV, then the cross-validation scores against the lambda values are plotted.
transf	An optional character string (with "y" as argument) meaning a function to apply to the predicted values (and possibly residuals) before plotting. E.g. " $\exp(y)-0.1$ ". If NULL (default) it is taken as the inverse of function <code>transf</code> (*if*) supplied in <code>gcrq</code> . See argument "transf" in <code>gcrq()</code> . If <code>transf</code> has been specified in <code>gcrq()</code> , use <code>transf="y"</code> to force plotting on the transformed scale, i.e. without back transforming.
lambda0	logical. If <code>cv=TRUE</code> , should the CV plot include also the first CV value? Usually the first CV value is at <code>lambda=0</code> , and typically it is much bigger than the other values making the plot not easy to read. Default to FALSE not to display the first CV value in the plot.
shade	logical. If TRUE and <code>conf.level>0</code> , the pointwise confidence intervals are portrayed via shaded areas.
overlap	NULL or numeric (scalar or vector). If provided and different from NULL, it represents the abscissa values (on the covariate scale) where the legends (i.e. the probability values) of each curve are set. It will be recycled, if its length differs from the number of quantile curves. If unspecified (i.e. <code>overlap=NULL</code>), the legends are placed outside the fitted lines on the right side. If specified, <code>legend=TRUE</code> is implicitly assumed.

<code>rug</code>	logical. If TRUE, the covariate distribution is displayed as a rug plot at the foot of the plot. Default to FALSE.
<code>overall.eff</code>	logical. If the smooth term has been called via <code>ps(..., decom=TRUE)</code> , by specifying <code>overall.eff=TRUE</code> the overall smooth effect is drawn, otherwise only the penalized part is portrayed (always <i>without</i> intercept).
<code>grid</code>	if provided, a grid of horizontal and vertical lines is drawn. <code>grid</code> has to be a list with the following components <code>x,y,col,lty,lwd</code> . If <code>x (y)</code> is a vector, the vertical (horizontal) lines are drawn at these locations. If <code>x (y)</code> is a scalar, the vertical (horizontal) lines are drawn at <code>x (y)</code> equispaced values. <code>col, lty,lwd</code> refer to the lines to be drawn.
<code>smoos</code>	logical, indicating if the residuals (provided that <code>res=TRUE</code>) will be drawn using a <i>smoothed</i> scatterplot. If NULL (default) the smoothed scatterplot will be employed when the number of observation is larger than 10000.
<code>split</code>	logical. If there are multiple terms (both smooth and linear) and <code>split=TRUE</code> , <code>plot.gcrq()</code> tries to split the plotting area in 2 columns and number of rows depending on the number of smooths. If <code>split=FALSE</code> , the plots are produced on the current device according to the current graphics settings. Ignored if there is single smooth term.
<code>shift</code>	Numerical value(s) to be added to the curve(s) to be plotted. If vector with length equal to the number of quantile curves to plot, the <code>shift[j]</code> is added to the <code>j</code> th quantile curve.
<code>type</code>	If <code>conf.level>0</code> , which covariance matrix should be used to compute and to portray the pointwise confidence intervals? 'boot' means case-resampling bootstrap (see <code>n.boot</code> in <code>gcrq()</code>), 'sandw' mean via the sandwich formula.
<code>n.points</code>	On how many values the plotted lines should rely on? If NULL, 100 values which suffice most of times. Increasing <code>n.points</code> can be useful when adaptive smoothing is used.
<code>...</code>	Additional graphical parameters: <code>xlab, ylab, ylim, and xlim</code> (effective when <code>add=FALSE</code>); <code>lwd, lty, and col</code> for the fitted quantile lines; <code>col<0</code> means color palette for the different curves; <code>cex</code> and <code>text.col</code> for the legend (if <code>legend=TRUE</code> or <code>overlap</code> is specified); <code>cex.p, col.p, and pch.p</code> for the points (if <code>res=TRUE</code>). When <code>axis.tau=TRUE</code> , all arguments accepted by <code>plot()</code> , <code>points()</code> , <code>matplot()</code> , and <code>matpoints()</code> but <code>pch, type, xlab, ylab, lty</code> .

Details

Takes a "gcrq" object and displays the fitted quantile curves as a function of the covariate specified in `term`. If `conf.level>0` pointwise confidence intervals are also displayed. When the object contains the component `cv`, `plot.gcrq` can display cross-validation scores against the lambda values, see argument `cv`. If a single quantile curve is being displayed, the default 'ylab' includes the relevant edf value (leaving out the basis intercept). If `axis.tau=TRUE` and the fit includes several quantile curves, `plot.gcrq()` portrays the estimated coefficients versus the probability values. If `term` refers to a categorical variable, the point estimates against the categories are plotted (`conf.level` is ignored).

Value

The function simply generates a new plot or adds fitted curves to an existing one.

Note

Plotting non-crossing curves could depend on the arguments 'interc' and 'shift', in turn depending on how the model has been specified. Take care about that!

Author(s)

Vito M. R. Muggeo

See Also

[gcrq](#), [predict.gcrq](#)

Examples

```
## Not run:
## use the fits from ?gcrq
##The additive model
plot(o, res=TRUE, col=2, conf.level=.9, shade=TRUE, split=TRUE)

par(mfrow=c(2,2))
plot(m5, select.tau=c(.1,.5,.9), overlap=0.6, legend=TRUE)
plot(m5, grid=list(x=8,y=5), lty=1) #a 8 times 5 grid..
plot(m7, cv=TRUE) #display CV score versus lambda values
plot(m7, res=TRUE, grid=list(x=5, y=8), col=4) #fitted curves at the best lambda value

## End(Not run)
```

predict.gcrq

Prediction for "gcrq" objects

Description

Takes a "gcrq" objects and computes fitted values

Usage

```
## S3 method for class 'gcrq'
predict(object, newdata, se.fit=FALSE, transf=NULL, xreg,
        type=c("sandw", "boot"), ...)
```

Arguments

object	a fitted "gcrq" object.
newdata	a dataframe including <i>all</i> the covariates of the model. The smooth term is represented by a covariate and proper basis functions will be build accordingly. If omitted, the fitted values are used. Ignored if xreg is provided.
se.fit	logical. If TRUE, standard errors of the fitted quantiles are computed using the bootstrap or the sandwich covariance matrix, according to the argument type.
transf	An optional character string (with "y" as argument) meaning a function to apply to the predicted values. E.g. "(exp(y)-0.1)". If NULL (default) it is taken as the inverse of function transf (*if*) supplied in gcrq. The standard errors (provided se.fit=TRUE has been set) are adjusted accordingly via the Delta method. See argument "transf" in gcrq(). If transf has been specified in gcrq(), use transf="y" to force predictions on the transformed scale, i.e. without back transforming.
xreg	the design matrix for which predictions are requested. If provided, xreg has to include the basis functions of the B-spline.
type	If se.fit=TRUE, which cov matrix should be used? 'boot' means case-resampling bootstrap (see n.boot in gcrq()), 'sandw' mean via the sandwich formula.
...	arguments passed to other functions

Details

predict.gcrq computes fitted quantiles as a function of observations included in newdata or xreg. Either newdata or xreg have to be supplied, but newdata is ignored when xreg is provided.

Value

If se.fit=FALSE, a matrix of fitted values with number of rows equal to number of rows of input data and number of columns depending on the number of fitted quantile curves (i.e length of taus).
If se.fit=TRUE, a list of matrices (fitted values and standard errors).

Author(s)

Vito M.R. Muggeo

See Also

[gcrq](#), [plot.gcrq](#)

Examples

```
##see ?gcrq
## predict(m1, newdata=data.frame(x=c(.3, .7)))
```

print.gcrq	<i>Print method for the gcrq class</i>
------------	----------------------------------------

Description

Printing the most important features of a gcrq model.

Usage

```
## S3 method for class 'gcrq'  
print(x, digits = max(3, getOption("digits") - 4), ...)
```

Arguments

x	object of class gcrq
digits	number of digits to be printed
...	arguments passed to other functions

Author(s)

Vito M.R. Muggeo

See Also

[summary.gcrq](#)

ps	<i>Specifying a smooth term in the gcrq formula.</i>
----	------------------------------------------------------

Description

Function used to define the smooth term (via P-splines) within the gcrq formula. The function actually does not evaluate a (spline) smooth, but simply it passes relevant information to proper fitter functions.

Usage

```
ps(..., lambda = -1, d = 3, by=NULL, ndx = NULL, deg = 3, knots=NULL,  
    monotone = 0, concave = 0, var.pen = NULL, pen.matrix=NULL, dropc=TRUE,  
    center=TRUE, K=NULL, decomp=FALSE, constr.fit=TRUE, shared.pen=FALSE,  
    st=FALSE, ad=0)
```

Arguments

...	The covariate supposed to have a nonlinear relationship with the quantile curve(s) being estimated. A B-spline is built, and a (difference) penalty is applied. In growth charts this variable is typically the age. If the covariate is a factor, category-specific coefficients are estimated subject to a lasso penalty. See the last example in <code>?gcrq</code> . A matrix of (continuous) covariates can be also supplied to perform variable selection (among its columns).
lambda	A supplied smoothing parameter for the smooth term. If it is negative scalar, the smoothing parameter is estimated iteratively as discussed in Muggeo et al. (2021). If a positive scalar, it represents the actual smoothing parameter. If it is a vector, cross validation is performed to select the ‘best’ value. See Details in gcrq .
d	The difference order of the penalty. Default to 3 Ignored if <code>pen.matrix</code> is supplied.
by	if different from NULL, a numeric or factor variable of the same dimension as the covariate in ... If numeric the elements multiply the smooth (i.e. a varying coefficient model); if factor, a smooth is fitted for each factor level. Usually the variable <code>by</code> is also included as main effect in the formula, see examples in gcrq . When <code>by</code> includes a factor, the formula should include the model intercept, i.e. $y \sim g + ps(x, by=g)$ and not $y \sim \theta + g + ps(x, by=g)$.
ndx	The number of intervals of the covariate range used to build the B-spline basis. Non-integer values are rounded by <code>round()</code> . If NULL, default, it is taken $\min(n/4, 9)$ (versions $\leq 1.1-0$ it was $\min(n/4, 40)$, the empirical rule of Ruppert). It could be reduced further (but no less than 5 or 6, say) if the sample size is not large and the default value leads to some error in the fitting procedure, see section Note in gcrq . Likewise, if the underlying relationship is strongly nonlinear, <code>ndx</code> could be increased. The returned basis will have ‘ <code>ndx+deg-1</code> ’ (if <code>dropc=TRUE</code>) basis functions.
deg	The degree of the spline polynomial. Default to 3. The B-spline basis is composed by <code>ndx+deg</code> basis functions and if <code>dropc=TRUE</code> the first column is removed for identifiability (and the model intercept is estimated without any penalty).
knots	The knots locations. If NULL, equispaced knots are set. Note if predictions outside the observed covariate range have to be computed (via <code>predict.gcrq</code>), the knots should be set enough outside the observed range.
monotone	Numeric value to set up monotonicity restrictions on the first derivative of fitted smooth function <ul style="list-style-type: none"> • ‘0’ = no constraint (default); • ‘1’ = non-decreasing smooth function; • ‘-1’ = non-increasing smooth function.
concave	Numeric value to set up monotonicity restrictions on the second derivative of fitted smooth function <ul style="list-style-type: none"> • ‘0’ = no constraint (default); • ‘1’ = concave smooth function; • ‘-1’ = convex smooth function.

<code>var.pen</code>	A character indicating the varying penalty. See Details.
<code>pen.matrix</code>	if provided, a penalty matrix A , say, such that the penalty in the objective function, apart from the smoothing parameter, is $\ Ab\ _1$ where b is the spline coefficient vector being penalized.
<code>dropc</code>	logical. Should the first column of the B-spline basis be dropped for the basis identifiability? Default to TRUE. Note, if <code>dropc=FALSE</code> is set, it is necessary to omit the model intercept AND not to center the basis, i.e. <code>center=FALSE</code> . Alternatively, both a full basis and the model intercept may be included by adding a small ridge penalty via <code>lambda.ridge>0</code> .
<code>center</code>	logical. If TRUE the smooth effects are 'centered' over the covariate values, i.e. $\sum_i \hat{f}(x_i) = 0$.
<code>K</code>	A scalar tuning the selection of wiggleness of the smoothed curve when λ has to be estimated (i.e. <code>lambda<0</code> is set). The larger K , the smoother the curve. Simulations suggest $K=2$ for the smoothing, and $K=\log(n/p^{2/3})$ for variable selection and random intercepts (p is the number of variables or number of subjects). See details.
<code>decom</code>	logical. If TRUE, the B-spline B (with a d order difference penalty) is decomposed into truncated power functions namely unpenalized polynomial terms up to degree $d-1$, and additional terms $Z = BD'(DD')^{-1}$. Only the coefficients of Z are penalized via an identity matrix, i.e. a lasso penalty. Currently <code>decom=TRUE</code> does not work with shape (monotonicity and concavity) restrictions and noncrossing constraints.
<code>constr.fit</code>	logical. If <code>monotone</code> or <code>concave</code> are different from 0, <code>constr.fit=TRUE</code> means that these constraints are set on the fitted quantiles rather than on the spline coefficients.
<code>shared.pen</code>	logical. If TRUE and the smooth is a VC term with a factor specified in <code>by</code> , the smooths in each level of the factor share the same smoothing parameter.
<code>st</code>	logical. If TRUE the variable(s) are standardized via the <code>scale()</code> function. Typically used for variable selection via lasso, i.e. when a matrix of covariates is passed in <code>ps()</code> .
<code>ad</code>	a positive number to carry out a form of <i>adaptive</i> lasso. More specifically, at each step of the iterative algorithm, the penalty is $\lambda \sum_j w_j \beta_j $ where $w_j = \tilde{\beta}_j ^{-ad}$ and $\tilde{\beta}_j$ are estimates coming from the previous iteration with a different value of λ . <code>ad=0</code> means the standard lasso and <code>ad=1</code> adaptive lasso (with weights updated during the iterative process).

Details

If a numeric variable has been supplied, `ps()` builds a B-spline basis with number of columns equal to `ndx+deg` (or `length(knots)-deg-1`). However, unless `dropc=FALSE` is specified, the first column is removed for identifiability, and the spline coefficients are penalized via differences of order d ; $d=0$ leads to a penalty on the coefficients themselves. If `pen.matrix` is supplied, d is ignored. Since versions 1.5-0 and 1.6-0, a factor or matrix can be supplied.

`lambda` is the tuning parameter, fixed or to be estimated. When `lambda=0` an unpenalized (and typically wiggly) fit is obtained, and as `lambda` increases the curve gets smoother till a $d-1$ degree

polynomial. At 'intermediate' lambda values, the fitted curve is a *piecewise* polynomial of degree $d-1$.

It is also possible to put a varying penalty via the argument `var.pen`. Namely for a constant smoothing (`var.pen=NULL`) the penalty is $\lambda \sum_k |\Delta_k^d|$ where Δ_k^d is the k -th difference (of order d) of the spline coefficients. For instance if $d = 1$, $|\Delta_k^1| = |b_k - b_{k-1}|$ where the b_k are the spline coefficients. When a varying penalty is set, the penalty becomes $\lambda \sum_k |\Delta_k^d| w_k$ where the weights w_k depend on `var.pen`; for instance `var.pen="((1:k)^2)"` results in $w_k = k^2$. See models `m6` and `m6a` in the examples of `gcrq`.

If `decom=TRUE`, the smooth can be plotted with or without the fixed part, see `overall.eff` in the function `plot.gcrq`.

Value

The function simply returns the covariate with added attributes relevant to smooth term.

Note

For shape-constrained fits, use `constr.fit=FALSE` only if you are using a single full and uncentred basis, namely something like
`gcrq(y~0+ps(x, center=FALSE, dropc=FALSE, monotone=1, constr.fit=FALSE),...)`.

Author(s)

Vito M. R. Muggeo

References

Muggeo VMR, Torretta F, Eilers PHC, Sciandra M, Attanasio M (2021). Multiple smoothing parameters selection in additive regression quantiles, *Statistical Modelling*, 21, 428-448.

For a general discussion on using B-spline and penalties in regression model see

Eilers PHC, Marx BD. (1996) Flexible smoothing with B-splines and penalties. *Statistical Sciences*, 11:89-121.

See Also

`gcrq`, `plot.gcrq`

Examples

```
##see ?gcrq
```

```
##gcrq(y ~ ps(x),...) #it works (default: center = TRUE, dropc = TRUE)
##gcrq(y ~ 0 + ps(x, center = TRUE, dropc = FALSE)) #it does NOT work
##gcrq(y ~ 0 + ps(x, center = FALSE, dropc = FALSE)) #it works
```

SiChildren

Age, height and weight in a sample of Italian children

Description

Age, height and weight in a sample of 1424 Italian children born in Sicily in the eighties

Usage

```
data("SiChildren")
```

Format

A data frame with 1424 observations on the following 3 variables.

age age in years

height child height (in centimeter)

weight child weight (in kilo)

Details

Data refer on the usual antropometric measures of Italian boys born in Sicily in the first years of 80s. Data have been kindly provided by prof M. Chiodi

Source

Gattuccio F., and Pirronello S., and Chiodi M (1988) Possibilita' di identificazione di tipologie evolutive del periodo puberale: proposta di una metodica pr finalita' predittive, *Rivista di pediatria preventiva e sociale nipiologia*, 189-199

Examples

```
data(SiChildren)
## see the package vignette for an example using such dataset
```

summary.gcrq

Summarizing model fits for growth charts regression quantiles

Description

summary and print methods for class gcrq

Usage

```
## S3 method for class 'gcrq'
summary(object, type=c("sandw", "boot"), digits = max(3, getOption("digits") - 3),
        signif.stars = getOption("show.signif.stars"), ...)
```

Arguments

object	An object of class "gcrq".
type	Which covariance matrix should be used to compute the estimate standard errors? 'boot' means case-resampling bootstrap (see <code>n.boot</code> in <code>gcrq()</code>), 'sandw' mean via the sandwich formula.
digits	controls number of digits printed in output.
signif.stars	Should significance stars be printed?
...	further arguments.

Details

`summary.gcrq` returns some information on the fitted quantile curve at different probability values, such as the estimates, standard errors, values of check (objective) function values at solution. Currently there is no `print.summary.gcrq` method, so `summary.gcrq` itself prints results.

The SIC returned by `print.gcrq` and `summary.gcrq` is computed as $\log(\rho_\tau/n) + \log(n)edf/(2n)$, where ρ_τ is the usual asymmetric sum of residuals (in absolute value). For multiple J quantiles it is $\log(\sum_\tau \rho_\tau/(nJ)) + \log(nJ)edf/(2nJ)$. Note that computation of SIC in `AIC.gcrq` relies on the Laplace assumption for the response.

Author(s)

Vito M.R. Muggeo

See Also

[gcrq](#)

Examples

```
## see ?gcrq
##summary(o)
```

`vcov.gcrq`*Variance-Covariance Matrix for a Fitted 'gcrq' Model*

Description

Returns the variance-covariance matrix of the parameter estimates of a fitted gcrq model object.

Usage

```
## S3 method for class 'gcrq'  
vcov(object, term, type=c("sandw", "boot"), ...)
```

Arguments

<code>object</code>	a fitted model object of class "gcrq" returned by <code>gcrq()</code> .
<code>term</code>	if specified, the returned covariance matrix includes entries relevant to parameter estimates for that 'term' only. If missing, the returned matrices refer to all model parameter estimates. Currently <code>term</code> is not allowed.
<code>type</code>	Which cov matrix should be returned? 'boot' means case-resampling bootstrap (see <code>n.boot</code> in <code>gcrq()</code>), 'sandw' mean via the sandwich formula.
<code>...</code>	additional arguments.

Details

Bootstrap-based covariance matrix, i.e. `type="boot"`, is computable only if the object fit has been obtained by specifying `n.boot>0` in `gcrq()`.

Value

A list (its length equal the length of `tau` specified in `gcrq`) of square matrices. Namely the list includes the covariance matrices of the parameter estimates for each regression quantile curve.

Author(s)

Vito Muggeo

See Also

[summary.gcrq](#)

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