Package ‘ordPens’

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Description Selection and/or smoothing of ordinally scaled independent variables using a group lasso or generalized ridge penalty.
Depends grplasso, mgcv, RLRsim, tcltk
Suggests IsoGene, Biobase
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ordPens-package

Selection and/or smoothing of ordinal predictors

Description

Selection and/or smoothing of ordinally scaled independent variables using a group lasso or generalized ridge penalty.

Details

Package: ordPens
Type: Package
Version: 0.3-1
Date: 2015-05-21
Depends: grplasso, mgcv, RLRsim, tcltk
Suggests: IsoGene, Biobase
License: GPL-2
LazyLoad: yes

Smoothing and selection of ordinal predictors is done by the function `ordSelect`; smoothing only, by `ordsmooth`. For ANOVA with ordinal factors, use `ordAOV`.

Author(s)

Jan Gertheiss

Maintainer: Jan Gertheiss, <jgerthe@uni-goettingen.de>

References


See Also

`ordSelect`, `ordSmooth`, `ordAOV`
Examples

# smooth modeling of a random dataset
set.seed(123)

# generate (ordinal) predictors
x1 <- sample(1:8,100,replace=TRUE)
x2 <- sample(1:6,100,replace=TRUE)
x3 <- sample(1:7,100,replace=TRUE)

# the response
y <- -1 + log(x1) + sin(3*(x2-1)/pi) + rnorm(100)

# x matrix
x <- cbind(x1,x2,x3)

# lambda values
lambda <- c(1000,500,200,100,50,30,20,10,1)

# smooth modeling
o1 <- ordSmooth(x = x, y = y, lambda = lambda)

# results
round(o1$coef,digits=3)
plot(o1)

# If for a certain plot the x-axis should be annotated in a different way,
# this can (for example) be done as follows:
plot(o1, whichx = 1, xlim = c(0,9), xaxt = "n")
axis(side = 1, at = c(1,8), labels = c("no agreement","total agreement"))

---

ICFCoreSetCWP  
ICF core set for chronic widespread pain

Description

The data set contains observed levels of ICF categories from the (comprehensive) ICF Core Set for chronic widespread pain (CWP) and a physical health component summary measure for n = 420 patients.

Usage

data(ICFCoreSetCWP)

Format

The data frame has 420 rows and 68 columns. The first 67 columns contain observed levels of ICF categories from the (comprehensive) ICF Core Set for chronic widespread pain (CWP). In the last column, the physical health component summary measure is given. Each row corresponds to one patient with CWP. ICF categories have discrete ordinal values between 0 and 4 (columns 1 - 50 and 67), or between -4 and 4 (columns 51 - 66). See the given references for details.
Details

The original data set contained some missing values, which have been imputed using R package Amelia.

The data were collected within the study Validation of ICF Core Sets for chronic conditions, which was a collaboration effort between the ICF Research Branch of the collaborating centers for the Family of International Classifications in German, the Classification, Terminology and standards Team from the World Health Organization and the International Society for Physical and Rehabilitation Medicine.

Special thanks go to the following participating study centers: Ankara University, Turkey; Azienda Ospedaliera di Sciacca, Italy; Donauspital, Vienna, Austria; Drei-Burgen-Klinik, Bad Muenster, Germany; Edertal Klinik, Bad Wildungen, Germany; Fachklinik Bad Bentheim, Germany; Hospital das Clínicas, School of Medicine, University of Sao Paulo, Brazil; Hospital San Juan Bautista, Catamarca, Argentina; Istituto Scientifico di Montescano, Italy; Istituto Scientifico di Veruno, Italy; Kaiser-Franz-Josef-Spital, Vienna, Austria; Klinik am Regenbogen, Nittenau, Germany; Klinik Bavaria Kreischa, Germany; Klinik Hoher Meissner, Bad Sooden-Allendorf, Germany; Klinikum Berchtesgadener Land, Schoenau, Germany; Kuwait Physical Medicine and Rehabilitation Society, Safat, Kuwait; National Institute for Medical Rehabilitation, Budapest, Hungary; Neuro-Orthopaedisches Krankenhaus und Zentrum fuer Rehabilitative Medizin Soltau, Germany; Praxis fuer Physikalische Medizin und Rehabilitation, Goettingen, Germany; Rehabilitationsklinik Seehof der Bundesversicherungsanstalt fuer Angestellte, Teltow, Germany; Rehaklinik Rheinfelden, Switzerland; Spanish Society of Rheumatology, Madrid, Spain; University Hospital Zurich, Switzerland; University of Santo Tomas, Quezon City, Philippines.

Most special thanks go to all the patients participating in the study.

If you use the data, please cite the following two references.

References


Examples

```r
# load the data
data(ICFCoreSetCWP)

# available variables
names(ICFCoreSetCWP)

# adequate coding of x matrix (using levels 1,2,...)
p <- ncol(ICFCoreSetCWP) - 1
n <- nrow(ICFCoreSetCWP)
add <- c(rep(1,50),rep(5,16),1)
add <- matrix(add,n,p,byrow=TRUE)
x <- ICFCoreSetCWP[,1:p] + add
```
ordAOV

# make sure that also a coefficient is fitted for levels
# that are not observed in the data
addrow <- c(rep(5,50),rep(9,16),5)
x <- rbind(x,addrow)
y <- c(ICFCoreSet$phcs,NA)

# some lambda values
lambda <- c(600,500,400,300,200,100)

# smoothing and selection
modelICF <- ordSelect(x = x, y = y, lambda = lambda)

# results
plot(modelICF)

# plot a selected ICF category (e.g. e110 'drugs')
# with adequate class labels
plot(modelICF, whichx = 51, xaxt = "n")
axis(side = 1, at = 1:9, labels = -4:4)

ordAOV

ANOVA for factors with ordered levels

Description

This function performs analysis of variance when the factor(s) of interest has/have ordinal scale level. For testing, values from the null distribution are simulated.

Usage

ordAOV(x, y, type = c("RLRT", "LRT"), nsim = 10000, null.sample = NULL, ...)

Arguments

x          a vector or matrix of integers 1,2,... giving the observed levels of the ordinal factor(s). If x is a matrix, it is assumed that each column corresponds to one ordinal factor.
y          the vector of response values.
type       the type of test to carry out: likelihood ratio ("LRT") or restricted likelihood ratio ("RLRT").
sim        number of values to simulate from the null distribution.
null.sample a vector, or a list of vectors (in case of multi-factorial ANOVA) containing values already simulated from the null distribution (overrides nsim)
...        additional arguments to LRTSim and RLRTSim, respectively.
Details

The method assumes that ordinal factor levels (contained in vector/columns of matrix `x`) take values 1,2,...,max, where max denotes the highest level of the respective factor observed in the data. Every level between 1 and max has to be observed at least once.

The method uses a mixed effects formulation of the usual one- or multi-factorial ANOVA model (with main effects only) while penalizing (squared) differences of adjacent means. Testing for equal means across factor levels is done by (restricted) likelihood ratio testing for a zero variance component in a linear mixed model. For simulating values from the finite sample null distribution of the (restricted) likelihood ratio statistic, the algorithms implemented in Package `RLRsim` are used. See `LRTSim` and `RLRTSim` for further information.

If `x` is a vector (or one-column matrix), one-factorial ANOVA is applied, and it is simulated from the exact finite sample null distribution as derived by Crainiceanu & Ruppert (2004). If `x` is a matrix, multi-factorial ANOVA (with main effects only) is done, and the approximation of the finite sample null distribution proposed by Greven et al. (2008) is used. Simulation studies by Gertheiss (2014) suggest that for ANOVA with ordinal factors RLRT should rather be used than LRT.

Value

In case of one-factorial ANOVA, a list of class `htest` containing the following components (see also `exactLRT` and `exactRLRT`):

- `statistic` the observed (restricted) likelihood ratio.
- `p` p-value for the observed test statistic.
- `method` a character string indicating what type of test was performed and how many values were simulated to determine the critical value.
- `sample` the samples from the null distribution returned by `LRTSim` and `RLRTSim`, respectively.

In case of multi-factorial ANOVA, a list (of lists) with the jth component giving the results above when testing the main effect of factor j.

Author(s)

Jan Gertheiss

References


ordGene


See Also

LRTSim, RLRTSim

Examples

# load some data
data(ICFCoresetCWP)

# the physical health component summary
y <- ICFCoresetCWP$phcs

# consider the first ordinal factor
x <- ICFCoresetCWP[,1]

# adequate coding
x <- as.integer(x - min(x) + 1)

# ANOVA
ordAOV(x, y, type = "RLRT", nsim=1000000)

ordGene

Testing for differentially expressed genes

Description

This function can be used to test for genes that are differentially expressed between levels of an ordinal factor, such as dose levels or ordinal phenotypes.

Usage

ordGene(xpr, lvs, type = c("RLRT", "LRT"), nsim = 1e6, null.sample=NULL, progressBar = TRUE, ...)

Arguments

xpr a matrix or data frame of gene expression data with Probe IDs as row names.
lvs a numeric vector containing the factor levels (e.g., dose levels) corresponding to the columns of xpr.
type the type of test to carry out: likelihood ratio ("LRT") or restricted likelihood ratio ("RLRT").
nsim number of values to simulate from the null distribution.
null.sample a vector containing values already simulated from the null distribution (overrides nsim)
ordGene

progress_bar enables or disables the progress bar; default is TRUE, set it as FALSE if problems with the tcltk package occur.

... additional arguments to `LRTSim` and `RLRTSim`, respectively.

Details

For each gene in the dataset, `ordAOV` is applied to test for differences between levels given in `lvs`. See `ordAOV` for further information on the testing procedure. Simulation studies by Gertheiss (2014) suggest that a restricted likelihood test (RLRT) should rather be used than a likelihood ratio test (LRT).

In addition to (R)LRT, results of usual one-way ANOVA (not taking the factor’s ordinal scale level into account) and a t-test assuming a linear trend across factor levels are reported. Note that the t-test does not assume linearity in the doses (such as 0, 0.5, 2.0, 5.0, ...), if given, but in the levels, i.e., 1, 2, 3, etc.

Value

A matrix containing the raw p-values for each gene (rows) when using (R)LRT, ANOVA or a t-test (columns).

Author(s)

Jan Gertheiss

References


See Also

`ordAOV`

Examples

```r
## Not run:
# use dopamine data from package IsoGene
require(IsoGene)
require(Biobase)
data(dopamine)
xpr <- data.frame(exprs(dopamine))
dose <- unlist pData(dopamine)
plot(dose, xpr[,83,], col=as.factor(dose), lwd=2, ylab="expression")
```
ordSelect

Select and smoothing of dummy coefficients of ordinal predictors

Description

Fits dummy coefficients of ordinally scaled independent variables with a group lasso penalty on differences of adjacent dummy coefficients.

Usage

ordSelect(x, y, u = NULL, z = NULL, offset = rep(0, length(y)), lambda, nu = 1, zeta = 1, model = c("linear", "logit", "poisson"), penscale = sqrt, scalex = TRUE, scalez = TRUE, scaleu = TRUE, nonpenx = NULL, nonpenz = NULL, nonpenu = NULL, intercept = TRUE, eps = 1e-3, ...)

Arguments

x the matrix of ordinal predictors, with each column corresponding to one predictor and containing numeric values from \{1,2,...\}; for each covariate, category 1 is taken as reference category with zero dummy coefficient.

y the response vector.

u a matrix (or data.frame) of additional categorical (nominal) predictors, with each column corresponding to one (additional) predictor and containing numeric values from \{1,2,...\}; corresponding dummy coefficients will be regularized using a groupwise simple Ridge penalty, and for each covariate category 1 is taken as reference category.

z a matrix (or data.frame) of additional metric predictors, with each column corresponding to one (additional) predictor; corresponding coefficients will be regularized using a simple Lasso penalty.

offset vector of offset values.

lambda vector of penalty parameters (in decreasing order). Optimization starts with the first component. See details below.
nu additional tuning parameter to control the strength of the penalty imposed on dummy coefficients corresponding to u. See details below.

zeta additional tuning parameter to control the strength of the penalty imposed on coefficients corresponding to z. See details below.

model the model which is to be fitted. Possible choices are "linear" (default), "logit" or "poisson". See details below.

penscale rescaling function to adjust the value of the penalty parameter to the degrees of freedom of the parameter group. See the references below.

scalex, scaleu, scalez logical. Should (split/dummy-coded) design matrices corresponding to x, u, resp. z be scaled to have unit variance over columns? See details below.

nonpenx, nonpenu, nonpenz vectors of indices indicating columns of x, u, resp. z whose regression coefficients are not penalized.

intercept logical. Should a (non-penalized) intercept be included in the model? Default is TRUE.

eps a (small) constant to be added to the columnwise standard deviations when scaling design matrices, to control the effect of very small stds. See details below.

... additional arguments.

Details

The method assumes that categorical covariates (contained in x and u) take values 1,2,...,max, where max denotes the (columnwise) highest level observed in the data. If any level between 1 and max is not observed, a corresponding (dummy) coefficient is fitted anyway. If any level > max is not observed but possible in principle, and a corresponding coefficient is to be fitted, the easiest way is to add a corresponding row to x (and u, z) with corresponding y value being NA.

In order to adjust the strength of penalty that is imposed on nominal or metric covariates, nu, resp. zeta can be chosen. The penalty parameter used for nominal variables (contained in u) is lambda*nu, for metric predictors (contained in z) it is lambda*zeta.

If a linear regression model is fitted, response vector y may contain any numeric values; if a logit model is fitted, y has to be 0/1 coded; if a poisson model is fitted, y has to contain count data.

If scalex, scaleu or scalez are TRUE, design matrices constructed from x, u, resp. z are scaled to have unit variance over columns. In the case of x, the design matrix is split-coded, in the case of u, it is dummy-coded, and in case of z, it is just z. If a certain x- or u- category, however, is observed only a few times, variances may become very small and scaling has enormous effects on the result and may cause numerical problems. Hence a small constant eps can be added to each standard deviation when used for scaling.

Value

An ordPen object, which is a list containing:

fitted the matrix of fitted response values of the training data. Columns correspond to different lambda values.
ordSelect

coefficients  the matrix of fitted coefficients with respect to dummy-coded (ordinal or nominal) categorical input variables (including the reference category) as well as metric predictors. Columns correspond to different lambda values.

model       the type of the fitted model: "linear", "logit", or "poisson".

lambda      the used lambda values.

xlevels     a vector giving the number of levels of the ordinal predictors.

ulevels     a vector giving the number of levels of the nominal predictors.

zcovars     the number of metric covariates.

Author(s)

Jan Gertheiss

References


See Also

plotNordPen, predictNordPen, icfcoresetCWP

Examples

# smoothing and selection of ordinal covariates using a random dataset
set.seed(123)

# generate (ordinal) predictors
x1 <- sample(1:8,100,replace=TRUE)
x2 <- sample(1:6,100,replace=TRUE)
x3 <- sample(1:7,100,replace=TRUE)

# the response
y <- -1 + log(x1) + sin(3*(x2-1)/pi) + rnorm(100)

# x matrix
x <- cbind(x1,x2,x3)

# lambda values
lambda <- c(1000,500,200,100,50,30,20,10,1)

# smoothing and selection
o2 <- ordSelect(x = x, y = y, lambda = lambda)
ordSmooth

`Smoothing dummy coefficients of ordinal predictors`

Description

Fits dummy coefficients of ordinally scaled independent variables with the sum of squared differences of adjacent dummy coefficients being penalized.

Usage

```r
ordSmooth(x, y, u = NULL, z = NULL, offset = rep(0, length(y)), lambda, nu = 1, zeta = 1, model = c("linear", "logit", "poisson"), penscale = identity, scalex = TRUE, scalez = TRUE, scaleu = TRUE, nonpenx = NULL, nonpenz = NULL, nonpenu = NULL, intercept = TRUE, eps = 1e-3, delta = 1e-6, maxit = 25, ...)
```

Arguments

- `x`  
  the matrix (or `data.frame`) of ordinal predictors, with each column corresponding to one predictor and containing numeric values from \{1,2,...\}; for each covariate, category 1 is taken as reference category with zero dummy coefficient.

- `y`  
  the response vector.

- `u`  
  a matrix (or `data.frame`) of additional categorical (nominal) predictors, with each column corresponding to one (additional) predictor and containing numeric values \{1,2,...\}; corresponding dummy coefficients will be regularized using a simple Ridge penalty, and for each covariate category 1 is taken as reference category.

- `z`  
  a matrix (or `data.frame`) of additional metric predictors, with each column corresponding to one (additional) predictor; corresponding coefficients will be regularized using a simple Ridge penalty.

- `offset`  
  vector of offset values.

- `lambda`  
  vector of penalty parameters (in decreasing order). Optimization starts with the first component. See details below.

- `nu`  
  additional tuning parameter to control the strength of the penalty imposed on dummy coefficients corresponding to `u`. See details below.
ordSmooth

zeta  additional tuning parameter to control the strength of the penalty imposed on coefficients corresponding to z. See details below.

model  the model which is to be fitted. Possible choices are "linear" (default), "logit" or "poisson". See details below.

penscale  rescaling function to adjust the value of the penalty parameter to the degrees of freedom of the parameter group.

scalex, scaleu, scalez  logical. Should (split/dummy-coded) design matrices corresponding to x, u, resp. z be scaled to have unit variance over columns? See details below.

nonpenx, nonpenu, nonpenz  vectors of indices indicating columns of x, u, resp. z whose regression coefficients are not penalized.

intercept  logical. Should a (non-penalized) intercept be included in the model? Default is TRUE.

eps  a (small) constant to be added to the columnwise standard deviations when scaling design matrices, to control the effect of very small stds. See details below.

delta  a small positive convergence tolerance which is used as stopping criterion for the penalized Fisher scoring when a logit or poisson model is fitted. See details below.

maxit  integer given the maximal number of (penalized) Fisher scoring iterations.

...  additional arguments.

Details

The method assumes that categorical covariates (contained in x and u) take values 1, 2, ..., max, where max denotes the (columnwise) highest level observed in the data. If any level between 1 and max is not observed, a corresponding (dummy) coefficient is fitted anyway. If any level > max is not observed but possible in principle, and a corresponding coefficient is to be fitted, the easiest way is to add a corresponding row to x (and u, z) with corresponding y value being NA.

In order to adjust the strength of penalty that is imposed on nominal or metric covariates, nu, resp. zeta can be chosen. The penalty parameter used for nominal variables (contained in u) is lambda*nu, for metric predictors (contained in z) it is lambda*zeta.

If a linear regression model is fitted, response vector y may contain any numeric values; if a logit model is fitted, y has to be 0/1 coded; if a poisson model is fitted, y has to contain count data.

If scalex, scaleu or scalez are TRUE, design matrices constructed from x, u, resp. z are scaled to have unit variance over columns. In the case of x, the design matrix is split-coded, in the case of u, it is dummy-coded, and in case of z, it is just z. If a certain x- or u- category, however, is observed only a few times, variances may become very small and scaling has enormous effects on the result and may cause numerical problems. Hence a small constant eps can be added to each standard deviation when used for scaling.

A logit or poisson model is fitted by penalized Fisher scoring. For stopping the iterations the criterion $\sqrt{\text{sum}(\text{b.new} - \text{b.old})^2 / \text{sum(b.old}^2)} < \text{delta}$ is used.
**Value**

An `ordPen` object, which is a list containing:

- `fitted` the matrix of fitted response values of the training data. Columns correspond to different lambda values.
- `coefficients` the matrix of fitted coefficients with respect to dummy-coded (ordinal or nominal) categorical input variables (including the reference category) as well as metric predictors. Columns correspond to different lambda values.
- `model` the type of the fitted model: "linear", "logit", or "poisson".
- `lambda` the used lambda values.
- `xlevels` a vector giving the number of levels of the ordinal predictors.
- `ulevels` a vector giving the number of levels of the nominal predictors.
- `zcovars` the number of metric covariates.

**Author(s)**

Jan Gertheiss

**References**


**See Also**

`plotNordpen`, `predictNordpen`

**Examples**

```r
# smooth modeling of a random dataset
set.seed(123)

# generate (ordinal) predictors
x1 <- sample(1:8,100,replace=TRUE)
x2 <- sample(1:6,100,replace=TRUE)
x3 <- sample(1:7,100,replace=TRUE)

# the response
y <- -1 + log(x1) + sin(3*(x2-1)/pi) + rnorm(100)

# x matrix
x <- cbind(x1,x2,x3)

# lambda values
lambda <- c(1000,500,200,100,50,30,20,10,1)

# smooth modeling
o1 <- ordSmooth(x = x, y = y, lambda = lambda)
```
plot.ordPen

# results
round(o1$coef,digits=3)
plot(o1)

# If for a certain plot the x-axis should be annotated in a different way,
# this can (for example) be done as follows:
plot(o1, whichx = 1, xlim = c(0,9), xaxt = "n")
axis(side = 1, at = c(1,8), labels = c("no agreement","total agreement"))

plot.ordPen  
Plot method for ordPen objects

Description

Takes a fitted ordPen object and plots estimated dummy coefficients of ordinal predictors for different lambda values.

Usage

## S3 method for class 'ordPen'
plot(x, whichlam = NULL, whichx = NULL,
     type = NULL, xlab = NULL, ylab = NULL, main = NULL,
     xlim = NULL, ylim = NULL, col = NULL, ...)

Arguments

x  
an ordPen object.

whichlam  
a vector of indices of lambda values corresponding to object$lambda for which plotting is done; if NULL, all values from object$lambda are considered.

whichx  
a vector of indices indicating the ordinal predictors whose dummy coefficients are plotted; e.g., set whichx=2, if you just want the plot for the second smooth term.

type  
1-character string giving the type of plot desired, see plot.default.

xlab  
a label for the x axis; if supplied then this will be used as the x label for all plots.

ylab  
a label for the y axis; if supplied then this will be used as the y label for all plots.

main  
a main title for the plot(s); if supplied then this will be used as the title for all plots.

xlim  
the x limits; if supplied then this pair of numbers are used as the x limits for each plot.

ylim  
the y limits; if supplied then this pair of numbers are used as the y limits for each plot.

col  
the plotting color; can be a vector of the same length as whichlam specifying different colors for different lambda values. Default is shades of gray: the higher lambda the darker.

...  
additional graphical parameters (see plot.default, or par).
predict.ordPen

**Value**

The function simply generates plots.

**Author(s)**

Jan Gertheiss

**See Also**

ordSelect, ordSmooth

**Examples**

```
# see for example
help(ordSelect)
```

---

### predict.ordPen

**Predict method for ordPen objects**

**Description**

Obtains predictions from an ordPen object.

**Usage**

```r
## S3 method for class 'ordPen'
predict(object, newx, newu = NULL, newz = NULL, 
offset = rep(0, nrow(as.matrix(newx))), 
type = c("link", "response"), ...)  
```

**Arguments**

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>object</td>
<td>an ordPen object.</td>
</tr>
<tr>
<td>newx</td>
<td>the matrix (or data.frame) of new observations of the considered ordinal predictors, with each column corresponding to one predictor and containing numeric values from {1,2,...}.</td>
</tr>
<tr>
<td>newu</td>
<td>a matrix (or data.frame) of new observations of the additional categorical (nominal) predictors, with each column corresponding to one (additional) predictor and containing numeric values {1,2,...}.</td>
</tr>
<tr>
<td>newz</td>
<td>a matrix (or data.frame) of new observations of the additional metric predictors, with each column corresponding to one (additional) predictor.</td>
</tr>
<tr>
<td>offset</td>
<td>potential offset values.</td>
</tr>
<tr>
<td>type</td>
<td>the type of prediction. type = &quot;link&quot; is on the scale of linear predictors, whereas type = &quot;response&quot; is on the scale of the response variable, i.e. type = &quot;response&quot; applies the inverse link function to the linear predictors.</td>
</tr>
<tr>
<td>...</td>
<td>additional arguments.</td>
</tr>
</tbody>
</table>
predict.ordPen

Value

A matrix of predictions whose columns correspond to the different values of the penalty parameter lambda of the ordPen object.

Author(s)

Jan Gertheiss

See Also

ordSelect, ordSmooth

Examples

# the training data
set.seed(123)

# generate (ordinal) predictors
x1 <- sample(1:8,100,replace=TRUE)
x2 <- sample(1:6,100,replace=TRUE)
x3 <- sample(1:7,100,replace=TRUE)

# the response
y <- -1 + log(x1) + sin(3*(x2-1)/pi) + rnorm(100)

# x matrix
x <- cbind(x1,x2,x3)

# lambda values
lambda <- c(1000,500,200,100,50,30,20,10,1)

# selecting and/or smoothing
o1 <- ordSmooth(x = x, y = y, lambda = lambda)
o2 <- ordSelect(x = x, y = y, lambda = lambda)

# new data
x1 <- sample(1:8,10,replace=TRUE)
x2 <- sample(1:6,10,replace=TRUE)
x3 <- sample(1:7,10,replace=TRUE)
newx <- cbind(x1,x2,x3)

# prediction
round(predict(o1, newx), digits=3)
round(predict(o2, newx), digits=3)
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