

Package: RcmdrPlugin.HH (via r-universe)

September 3, 2024

Type Package

Title Rcmdr Support for the HH Package

Version 1.1-51

Date 2024-03-07

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Depends R (>= 3.0.2), HH

Imports Rcmdr (>= 2.0-0), lattice, mgcv

Suggests car, leaps, latticeExtra, rgl

Description Rcmdr menu support for many of the functions in the HH package. The focus is on menu items for functions we use in our introductory courses.

License GPL (>= 2)

NeedsCompilation no

Date/Publication 2024-03-06 23:40:11 UTC

Repository <https://rmheiberger.r-universe.dev>

RemoteUrl <https://github.com/cran/RcmdrPlugin.HH>

RemoteRef HEAD

RemoteSha 75a1aecbcf03dab24d6cb471b44c7c70201fc5a0

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RcmdrPlugin.HH-package

Functions added to the Rcmdr package to support the introductory course at Temple University.

Description

Our introductory course spends time on several topics that are not yet in the R Commander. Therefore we wrote the menu items and make them available.

Details

The DESCRIPTION file:

```

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Depends:    R (>= 3.0.2), HH
Imports:    Rcmdr (>= 2.0-0), lattice, mgcv
Suggests:   car, leaps, latticeExtra, rgl
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```

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[bestSubsetsRegressionModel.HH](#) Rcmdr interface to the [regsubsets](#) function in the leaps package.

[twoWayTable.HH](#) Pearson's Chi-squared Test for Count Data (additional formats for data input)

[anovaTableI.HH](#) Sequential sums of squares on the Rcmdr menu.

[scatter3dHH](#) add the ability to plot squared residuals. The squared residuals have been adopted into Rcmdr. This interface offers a checkbox for a new 3D window and an option to draw a non-

least-squares plane for pedagogical comparison.

`ci.plot` Plot confidence and prediction intervals for simple linear regression.

`panel.bwplot.intermediate.hh` Panel function for `bwplot` that give the user control over the placement of the boxes.

`interaction2wt` Plot all main effects and twoway interactions in a multifactor design.

`scatterPlotMatrix.HH` Similar to `scatterplotMatrix` The revision uses `row1atop=FALSE` to force the main diagonal of the scatterplot matrix to go uphill from southwest to northeast.

`QQPlot.HH` Added Shapiro-Wilk test of normality.

`norm.curve` Plot a normal curve with shaded rejection regions, optionally a second curve centered at an alternative hypothesis value can be plotted. Both x and z scales are displayed.

Author(s)

Richard M. Heiberger, with contributions from Burt Holland

Maintainer: Richard M. Heiberger <rmh@temple.edu>

References

Heiberger, Richard M. and Holland, Burt (2015). *Statistical Analysis and Data Display: An Intermediate Course with Examples in R*. Second Edition. Springer-Verlag, New York. <https://link.springer.com/book/10.1007/978-1-4939-2122-5>

Heiberger, Richard M. and Holland, Burt (2004). *Statistical Analysis and Data Display: An Intermediate Course with Examples in S-Plus, R, and SAS*, First Edition. Springer Texts in Statistics. Springer. <https://link.springer.com/book/10.1007/978-1-4757-4284-8>.

See Also

`Rcmdr`

Examples

```
## Not run:
## start R
library(RcmdrPlugin.HH)      ## loads the package and opens the Rcmdr
                             ## window with the HH menu

## End(Not run)
```

anovaTableI.HH *Rcmdr interface to anova function*

Description

Rcmdr interface to anova function, specifically to get the sequential sums of squares.

Usage

```
anovaTableI.HH()  
anovaTableII.HH() ## exact copy of John Fox's anovaTable from Rcmdr/R/model-menu.R
```

Author(s)

Richard M. Heiberger <rmh@temple.edu>

See Also

[anova](#)

bestSubsetsRegressionModel.HH
Rcmdr interface to the regsubsets function in the leaps package.

Description

Menu interface to the Best Subsets Regression function. Selection boxes allow one response variables and one or more predictor variables. All subsets are calculated. Only the best k , where k is menu item, are displayed. A graph displaying one of the following statistics (R^2 , residual sum of squares, adjusted R^2 , CV_p , BIC, MAE) is displayed. The model with highest adjusted R^2 is made the active model and its summary is displayed.

Usage

```
bestSubsetsRegressionModel.HH()
```

Author(s)

Richard M. Heiberger <rmh@temple.edu>

See Also

[regsubsets](#)

BoxCox

Rcmdr BoxCox demo renamed to active function.

Description

Rcmdr menu for Box-Cox Transformations

Usage

BoxCox()

Author(s)

John Fox <jfox@mcmaster.ca>

CloseCommanderRestart *Close Rcmdr without questions and then restart.*

Description

Close Rcmdr without questions. CloseCommanderNoQuestionRestart has absolutely no questions. CloseCommanderRestart asks only about saving files. Both functions restart Rcmdr immediately and therefore have the full .GlobalEnv from the R session still available.

Usage

CloseCommanderNoQuestionRestart()
CloseCommanderRestart()

Author(s)

Richard M. Heiberger <rmh@temple.edu>

See Also

[closeCommander](#)

`confidenceIntervalsPlot`*Rcmdr interface to plot confidence and prediction intervals in simple linear regression*

Description

Rcmdr menu interface to the function `ci.plot`. Variable boxes are provided for one predictor variable, one response variable. The simple linear regression is calculated and made the active model.

Usage`confidenceIntervalsPlot()`**Author(s)**Richard M. Heiberger <rmh@temple.edu>**See Also**[ci.plot](#)

`DotplottbRcmdr`*Rcmdr menu interface to `dotplot(panel=panel.dotplot.tb)`.*

Description

Rcmdr menu interface to `dotplot(panel=panel.dotplot.tb)`.

Usage`DotplottbRcmdr()`**Author(s)**Richard M. Heiberger <rmh@temple.edu>**See Also**[panel.dotplot.tb](#)

Interaction2wtRcmdr *Rcmdr menu interface to interaction2wt*

Description

Plot all main effects and twoway interactions in a multifactor design. The main diagonal displays boxplots for the main effects of each factor. The off-diagonals show the interaction plots for each pair of factors. The *i, j* panel shows the same factors as the *j, i* but with the trace- and x-factor roles interchanged.

Usage

```
Interaction2wtRcmdr()
```

Author(s)

Richard M. Heiberger <rmh@temple.edu>

See Also

[interaction2wt](#)

latticeFunctions *Support functions for the Xyplot.HH2 function.*

Description

Support functions for the Xyplot.HH2 function.

Usage

```
latticeFunctions()
latticePanelFunctions()
splomFormula(predictor, data.frame.name)
usualFormula(functionName, response, predictor, data.frame.name)
```

Arguments

predictor, data.frame.name, functionName, response
 Arguments to functions.

Value

For `latticeFunctions` and `latticePanelFunctions`, vector of function names. For `splomFormula` and `usualFormula`, a model formula containing the specified variable names.

Author(s)

Richard M. Heiberger <rmh@temple.edu>

See Also

[Xyplot.HH](#)

MMCMenu

Menu interface to MMC plots.

Description

Menu interface to MMC (Mean–mean Multiple Comparison) plots.

Usage

```
MMCMenu()  
MMC2menu()  
AOVModelSP(n=1)
```

Arguments

n Minimum number of "aov" models.

Author(s)

Richard M. Heiberger <rmh@temple.edu>

normal.and.t.hypotheses.plot

Rcmdr normalHypothesesPlot and tHypothesesPlot menu.

Description

Rcmdr menus to draw graphs of hypotheses, critical values, and p-values.

Usage

```
normal.and.t.hypotheses.plot()
```

```
FHypothesesPlot()
```

```
ChisqHypothesesPlot()
```

Author(s)

Richard M. Heiberger <rmh@temple.edu>.

See Also

[norm.curve](#), [F.curve](#), [chisq.curve](#)

PlotLikertDialog *Rcmdr Menu function to specify a likert plot.*

Description

Please see [likert](#) for details on the `plot.likert` and related functions.

Usage

```
PlotLikertDialog()
listAllLikertCapable(envir = .GlobalEnv, ...)
LikertFormula()
LikertFormulaConstruct(functionName, response, predictor)
varPosnOriginal(variables, type = c("all", "factor", "numeric", "nonfactor",
  "twoLevelFactor"))
```

Arguments

`envir, ...` Arguments to [ls](#).
`functionName, response, predictor`
 Arguments to functions.
`variables, type` See [varPosn](#) for discussion.

Value

For `listAllLikertCapable`, a character vector of names of all objects that satisfy the search criteria in the specified environments. See [likert](#) for details on what objects are likert capable.

`LikertFormulaConstruct` constructs a model formula for use by `plot.likert.formula` from its input arguments.

`varPosnOriginal` is the same as [varPosn](#) except that it always keeps the variables in the same order as the original `data.frame`.

Author(s)

Richard M. Heiberger <rmh@temple.edu>

See Also

[likert](#)

PredictModel	<i>Rcmdr menu interface to predict</i>
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Description

Rcmdr menu interface to predict

Usage

```
PredictModel()
```

Author(s)

Richard M. Heiberger <rmh@temple.edu>

See Also

[predict](#)

Projector	<i>Set Rcmdr options for good visibility on classroom projector and on netbook screen with 600 pixel height.</i>
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Description

Set Rcmdr options for good visibility on classroom projector and on netbook screen with 600 pixel height.

Usage

```
Projector()  
H600()
```

Author(s)

Richard M. Heiberger <rmh@temple.edu>

`QQPlot.HH`*Quantile-Comparison (QQ) Plot*

Description

Rcmdr menu interface to plot the qqplot of variable against one of the following distributions: normal, t, chi-square, F, other.

Usage`QQPlot.HH()`**Details**

The `normal` gives the option to do the Shapiro-Wilk test of normality. The `other` requires you to specify the distribution. Any distribution for which quantile and density functions exist in R (with prefixes `q` and `d`, respectively) may be used.

Value

NULL. These functions are used only for their side effect (to make a graph).

Author(s)

John Fox <jfox@mcmaster.ca>. Shapiro-Wilk test added by Richard M. Heiberger <rmh@temple.edu>.

See Also

[qqPlot](#), [shapiro.test](#)

`Regr1Plot`*Rcmdr Menu function to display the squared residuals.*

Description

Rcmdr Menu function to display the squared residuals of a linear fit of one y variable on one x variable. The default model is simple linear regression $y \sim x$. Any other model of one y on one x may be used. See the last example in [regr1.plot](#) for an example of a quadratic function of x.

Usage`Regr1Plot()`**Author(s)**

Richard M. Heiberger <rmh@temple.edu>

See Also[regr1.plot](#)

ResizeEtcDialog	<i>Rcmdr Menu function to specify combining and resizing "trellis" objects.</i>
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Description

Please see [ResizeEtc](#) for details on the combination of "trellis" and related functions.

Usage

```
ResizeEtcDialog()  
listAllTrellisObjects(envir = .GlobalEnv, ...)
```

Arguments

envir, ... Arguments to [ls](#).

Details

This dialog is a template designed to help with writing commandline code.

Value

For `listAllTrellisObjects`, a character vector of names of all "trellis" objects that satisfy the search criteria in the specified environments.

Author(s)

Richard M. Heiberger <rmh@temple.edu>

See Also

[ResizeEtc](#), [resizePanels](#)

R_options

Set R options from within R commander.

Description

Set R options from within R commander.

Usage

R_options()

Author(s)

Richard M. Heiberger <rmh@temple.edu>

Scatter3DDialog.HH

Rcmdr 3D Scatterplot Dialog (HH)

Description

This dialog sets up a call to the [scatter3dHH](#) function to draw a three-dimensional scatterplot, and optionally to [Identify3d](#) to label points interactively with the mouse.

Details

The explanatory variables provide the "horizontal" and "out-of-screen" axes of the scatterplot, the response variable provides the "vertical" axis.

Data points are represented as spheres or points, depending upon the number of observations.

Several regression surfaces can be plotted: a linear least-squares surface; a full quadratic least-squares surface with squared and cross-product terms; a "smooth" regression surface — either a smoothing spline, if no degrees of freedom are specified (in which case the [gam](#) function selects the df by generalized cross validation), or a fixed-df regression spline; an additive-regression surface (also fit by [gam](#)), with either smoothing spline or regression spline components (again selected according to the specification of degrees of freedom). If only one surface is fit, then residuals are plotted as red (negative) and green (positive) lines from the surface to the points. If the squared residuals option is checked, then squared residuals are plotted. The sum of the area of these squares is the "residual sum of squares".

You can specify a factor defining groups by pressing the *Plot by groups* button. A separate surface or set of surfaces is plotted for each level of the groups factor. These surfaces can be constrained to be parallel.

The completed plot can be manipulated with the mouse: Click, hold, drag the left mouse button to rotate the display; click, hold, and drag the right button (or centre button on a three-button mouse) to zoom in and out.

If the box labelled *Identify observations with mouse* is checked, you may use the mouse to identify points interactively: Press the right mouse button (or the centre button on a three-button mouse), drag a rectangle around the points to be identified, and release the button. Repeat this procedure for each point or set of "nearby" points to be identified. To exit from point-identification mode, right-click (or centre-click) in an empty region of the plot.

Points may also be identified subsequently by selecting *Identify observations with mouse* from the R Commander *3D graph* menu: As above, click and drag the left mouse button to rotate the display, and click and drag the right (or centre) button to identify points.

Author(s)

John Fox <jfox@mcmaster.ca>. Squared residuals added by Richard M. Heiberger <rmh@temple.edu>.

See Also

[scatter3dHH](#), [Identify3d](#), [rgl.open](#), [gam](#)

scatter3dHH

Three-Dimensional Scatterplots and Point Identification

Description

The scatter3d function uses the rgl package to draw 3D scatterplots with various regression surfaces. The function Identify3d allows you to label points interactively with the mouse: Press the right mouse button (on a two-button mouse) or the centre button (on a three-button mouse), drag a rectangle around the points to be identified, and release the button. Repeat this procedure for each point or set of "nearby" points to be identified. To exit from point-identification mode, click the right (or centre) button an empty region of the plot.

This is a revision of the Rcmdr scatter3d to add the ability to plot squared residuals.

Usage

```
scatter3dHH(x, y, z,
  xlab=deparse(substitute(x)), ylab=deparse(substitute(y)),
  zlab=deparse(substitute(z)),
  revolutions=0, bg.col=c("white", "black"),
  axis.col=if (bg.col == "white") "black" else "white",
  surface.col=c("blue", "green", "orange", "magenta",
    "cyan", "red", "yellow", "gray"),
  neg.res.col="red", pos.res.col="green", point.col="yellow",
  text.col=axis.col,
  grid.col=if (bg.col == "white") "black" else "gray",
  fogtype=c("exp2", "linear", "exp", "none"),
  residuals=(length(fit) == 1), surface=TRUE, grid=TRUE,
  grid.lines=26, df.smooth=NULL, df.additive=NULL,
  sphere.size=1, threshold=0.01, speed=1, fov=60,
  fit="linear", groups=NULL, parallel=TRUE, model.summary=FALSE,
  squares = FALSE, square.color = "gray", coef.ratio = 1, ...)
```

Arguments

<code>x</code>	variable for horizontal axis.
<code>y</code>	variable for vertical axis (response).
<code>z</code>	variable for out-of-screen axis.
<code>xlab, ylab, zlab</code>	axis labels.
<code>revolutions</code>	number of full revolutions of the display.
<code>bg.col</code>	background colour; one of "white", "black".
<code>axis.col</code>	colour for axes; default is "white" for black background, "black" for white background.
<code>surface.col</code>	vector of colours for regression planes, used in the order specified by <code>fit</code> .
<code>neg.res.col, pos.res.col</code>	colours for lines representing negative and positive residuals.
<code>point.col</code>	colour of points.
<code>text.col</code>	colour of axis labels.
<code>grid.col</code>	colour of grid lines on the regression surface(s).
<code>fogtype</code>	type of fog effect; one of "exp2", "linear", "exp", "none".
<code>residuals</code>	plot residuals (TRUE or FALSE); available only when there is one surface plotted.
<code>surface</code>	plot surface(s) (TRUE or FALSE).
<code>grid</code>	plot grid lines on the regression surface(s) (TRUE or FALSE).
<code>grid.lines</code>	number of lines (default, 26) forming the grid, in each of the x and y directions.
<code>df.smooth</code>	degrees of freedom for the two-dimensional smooth regression surface; if NULL (the default), the <code>gam</code> function will select the degrees of freedom for a smoothing spline by generalized cross-validation; if a positive number, a fixed regression spline will be fit with the specified degrees of freedom.
<code>df.additive</code>	degrees of freedom for each explanatory variable in an additive regression; if NULL (the default), the <code>gam</code> function will select degrees of freedom for the smoothing splines by generalized cross-validation; if a positive number or a vector of two positive numbers, fixed regression splines will be fit with the specified degrees of freedom for each term.
<code>sphere.size</code>	relative sizes of spheres representing points; the actual size is dependent on the number of observations.
<code>threshold</code>	if the actual size of the spheres is less than the threshold, points are plotted instead.
<code>speed</code>	relative speed of revolution of the plot.
<code>fov</code>	field of view (in degrees); controls degree of perspective.
<code>fit</code>	one or more of "linear", "quadratic", "smooth", "additive"; to display fitted surface(s); partial matching is supported – e.g., <code>c("lin", "quad")</code> .
<code>groups</code>	if NULL (the default), no groups are defined; if a factor, a different surface or set of surfaces is plotted for each level of the factor; in this event, the colours in <code>plane.col</code> are used successively for the points, surfaces, and residuals corresponding to each level of the factor.

<code>parallel</code>	when plotting surfaces by groups, should the surfaces be constrained to be parallel? A logical value, with default TRUE.
<code>model.summary</code>	print summary or summaries of the model(s) fit (TRUE or FALSE).
<code>col</code>	colours for the point labels, given by group. There must be at least as many colours as groups; if there are no groups, the first colour is used. Normally, the colours would correspond to the <code>plane.col</code> argument to <code>scatter3d</code> .
<code>squares</code>	logical. If TRUE, the residuals are plotted as squares. The sum of the area of the squares is the "residual sum of squares". If FALSE, the residuals are plotted as vertical lines.
<code>square.color</code>	color for the squares.
<code>coef.ratio</code>	number, defaults to 1. Setting <code>coef.ratio</code> to a number other than 1 is a primitive way of plotting squared pseudo-residuals that are not the least-squares residuals. The reason for displaying non-least-squares residuals is to compare them to the least-squares residuals and thus get a 3d visual image of what minimizing the sum of squares actually means.
<code>...</code>	other arguments are ignored.

Value

`scatter3d` not return a useful value; it is used for its side-effect of creating a 3D scatterplot. `Identify3d` returns the labels of the identified points.

Note

You have to install the `rgl` and `mgcv` packages to produce 3D plots.

Author(s)

John Fox <jfox@mcmaster.ca>. Squared residuals added by Richard M. Heiberger <rmh@temple.edu>.

See Also

[rgl.open](#), [gam](#)

Examples

```
## Not run:
State.x77 <- as.data.frame(state.x77)
with(State.x77, scatter3d(Income, Murder, Illiteracy))
with(State.x77, Identify3d(Income, Murder, Illiteracy, labels=row.names(State.x77)))
with(State.x77, scatter3d(Income, Murder, Illiteracy, fit=c("linear", "quadratic")))

## End(Not run)
```

scatterPlot.HH *Scatterplot menu with different defaults than Rcmdr.*

Description

Alternate menu into the scatterplot in the car package. This menu by default uses solid dots, larger fonts, and turns off marginal boxplots and smoother lines. Otherwise it is identical to the Rcmdr Scatterplot menu item.

Usage

```
scatterPlot.HH()
```

Author(s)

Richard M. Heiberger <rmh@temple.edu>

scatterPlotMatrix.HH *Scatterplot Matrices*

Description

This is variation of the Rcmdr interface to the car package scatterplotMatrix. The revision uses row1atop=FALSE to force the main diagonal of the scatterplot matrix to go uphill from southwest to northeast.

Usage

```
scatterPlotMatrix.HH()
```

Author(s)

John Fox <jfox@mcmaster.ca>. row1atop=FALSE added by Richard M. Heiberger <rmh@temple.edu>.

See Also

[scatterplotMatrix](#)

twoWayTable.HH	<i>Rcmdr menu interface to chisq.test</i>
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Description

Pearson's Chi-squared Test for Count Data

twoWayTable.HH is an original Rcmdr.HH function. It reads the active dataset and constructs the table using xtabs.

enterTable.HH is an original Rcmdr.HH function. It opens a window where the user may enter a table manually.

analyzeTwoWayTable.HH is an additional function. It uses the active dataset as the table.

All three produce identical output, a two-way table, row and column summaries, and the chi square test.

Usage

```
twoWayTable.HH()
```

```
enterTable.HH()
```

```
analyzeTwoWayTable.HH()
```

Author(s)

John Fox <jfox@mcmaster.ca>. additional entry options by Richard M. Heiberger <rmh@temple.edu>.

See Also

[chisq.test](#)

Xyplot.HH	<i>Rcmdr Menu function to specify xyplot, other lattice plots, and likert plots.</i>
-----------	--

Description

These are enhancements of the Rcmdr Xyplot function (which I wrote) to include layout parameters and plot type, to force solid dots, and to distinguish between conditioning variables in the formula and group variables. Xyplot.HH is an interface to the xyplot function. Xyplot.HH2 is an interface to many of the lattice functions (xyplot, bwplot, splom, barchart, dotplot) and to the formula method for likert in the HH package. When either barchart or panel.barchart is selected, then the argument origin=0 is automatically set. When panel.barchart, the user must manually specify the limits (xlim or ylim) to include zero for the effect of origin=0 to be visible.

Usage

```
Xyplot.HH()  
Xyplot.HH2()
```

Author(s)

Richard M. Heiberger <rmh@temple.edu>

See Also

[xyplot](#), [likert](#)

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