

Package ‘meboot’

July 10, 2021

Version 1.4-9.1

Date 2021-07-10

Title Maximum Entropy Bootstrap for Time Series

Author Hrishikesh D. Vinod <Vinod@fordham.edu>, Javier López-de-Lacalle <javlacalle@yahoo.es>, and Fred Violle

Maintainer Fred Violle <fviolle@fordham.edu>

Encoding UTF-8

Depends R (>= 3.5.0), dynlm, nlme, tdigest, hrcrde

Suggests boot, car, ConvergenceConcepts, geopack, lmtest, strucchange, plm, zoo

Description Maximum entropy density based dependent data bootstrap. An algorithm is provided to create a population of time series (ensemble) without assuming stationarity. The reference paper (Vinod, H.D., 2004 <[DOI:10.1016/j.jempfin.2003.06.002](https://doi.org/10.1016/j.jempfin.2003.06.002)>) explains how the algorithm satisfies the ergodic theorem and the central limit theorem.

License GPL (>= 2)

NeedsCompilation yes

RoxygenNote 7.1.1

Repository CRAN

Date/Publication 2021-07-10 18:20:02 UTC

R topics documented:

| | |
|--------------------|----|
| checkConv | 2 |
| elapsedtime | 3 |
| expand.sd | 4 |
| flexMeboot | 5 |
| force.clt | 6 |
| meboot | 7 |
| meboot.part | 9 |
| meboot.pdata.frame | 10 |

| | |
|-----------------------|----|
| mebootSpear | 11 |
| null.ci | 14 |
| olsHALL.b | 15 |
| ullwan | 16 |
| USconsum | 17 |
| USfygt | 18 |
| zero.ci | 19 |

| | |
|--------------|-----------|
| Index | 20 |
|--------------|-----------|

| | |
|-----------|--------------------------|
| checkConv | <i>Check Convergence</i> |
|-----------|--------------------------|

Description

This function generates a 3D array giving $(X_n - X)$ in the notation of the ConvergenceConcepts package by Lafaye de Micheaux and Liquet for sample paths with dimensions = n999 as first dimension, nover = range of n values as second dimension and number of items in key as the third dimension. It is intended to be used for checking convergence of meboot in the context of a specific real world time series regression problem.

Usage

```
checkConv (y, bigx, trueb = 1, n999 = 999, nover = 5,
  seed1 = 294, key = 0, trace = FALSE)
```

Arguments

| | |
|-------|---|
| y | vector of data containing the dependent variable. |
| bigx | vector of data for all regressor variables in a regression or ts object. bigx should not include column of ones for the intercept. |
| trueb | true values of regressor coefficients for simulation. If trueb=0 then use OLS coefficient values rounded to 2 digits as true values of beta for simulation purposes, to be close to but not exactly equal to OLS. |
| n999 | number of replicates to generate in a simulation. |
| nover | number of values of n over which convergence calculated. |
| seed1 | seed for the random number generator. |
| key | the subset of key regression coefficient whose convergence is studied if key=0 all coefficients are studied for convergence. |
| trace | logical. If TRUE, tracing information on the process is printed. |

Details

Use this only when lagged dependent variable is absent.

Warning: `key=0` might use up too much memory for large regression problems.

The algorithm first creates data on the dependent variable for a simulation using known true values denoted by `trueb`. It proceeds to create `n999` regression problems using the seven-step algorithm in `meboot` creating `n999` time series for all variable in the simulated regression. It then creates sample paths over a range of `n` values for coefficients of interest denoted as `key` (usually a subset of original coefficients). For each `key` coefficient there are `n999` paths as `n` increases. If `meboot` algorithm is converging to true values, the value of $(X_n - X)$ based criteria for "convergence in probability" and "almost sure convergence" in the notation of the `ConvergenceConcepts` package should decline. The decline can be plotted and/or tested to check if it is statistically significant as sample size increases. This function permits the user of `meboot` working with a short time series to see if the `meboot` algorithm is working in his or her particular situation.

Value

A 3 dimensional array giving $(X_n - X)$ for sample paths with dimensions = `n999` as first dimension, `nover` = range of `n` values as second dimension and number of items in `key` as the third dimension ready for use in `ConvergenceConcepts` package.

References

- Lafaye de Micheaux, P. and Liquet, B. (2009), Understanding Convergence Concepts: a Visual-Minded and Graphical Simulation-Based Approach, *The American Statistician*, **63**(2) pp. 173-178.
- Vinod, H.D. (2006), Maximum Entropy Ensembles for Time Series Inference in Economics, *Journal of Asian Economics*, **17**(6), pp. 955-978
- Vinod, H.D. (2004), Ranking mutual funds using unconventional utility theory and stochastic dominance, *Journal of Empirical Finance*, **11**(3), pp. 353-377.

See Also

[meboot](#), [criterion](#).

elapsedtime

Internal Function

Description

Internal function.

Usage

```
elapsedtime(ptm1, ptm2)
```

Arguments

ptm1 `proc.time` object.
ptm2 `proc.time` object.

Value

List giving the elapsed time between ptm1 and ptm2.

| | |
|-----------|---|
| expand.sd | <i>Expand the Standard Deviation of Resamples</i> |
|-----------|---|

Description

This function expands the standard deviation of the simulated data. Expansion is needed since some of the ratios of the actual standard deviation to that of the original data are lower than 1 due to attenuation.

Usage

```
expand.sd(x, ensemble, fiv=5)
```

Arguments

x a vector of data or a time series object.
ensemble a matrix or mts object containing resamples of the original data x.
fiv reference value for the upper limit of a uniform distribution used in expansion. For example, if equal to 5 the standard deviation of each resample is expanded through a value from a uniform random distribution with lower limit equal to 1 and upper limit equal to $1+(5/100)=1.05$.

Value

Resamples (by columns) with expanded standard deviations.

Examples

```
set.seed(345)  
out <- meboot(x=AirPassengers, reps=100, trim=0.10, reachbnd=FALSE, elaps=TRUE)  
exp.ens <- expand.sd(x=AirPassengers, out$ensemble)
```

flexMeboot

*Flexible Extension of the Maximum Entropy Bootstrap Procedure***Description**

This function extends the maximum entropy bootstrap procedure implemented in `meboot` to allow for for a flexible trend up, flat or down.

Usage

```
flexMeboot (x, reps = 9, segment = 5, forc = FALSE, myseq = seq(-1, 1, by = 1))
```

Arguments

| | |
|----------------------|---|
| <code>x</code> | vector of data, ts object. |
| <code>reps</code> | number of replicates to generate. |
| <code>segment</code> | block size. |
| <code>forc</code> | logical. If TRUE the ensemble is forced to satisfy the central limit theorem. See force.clt . |
| <code>myseq</code> | directions for trend within a block of data is chosen randomly with the user's choice limited by the range of values given by <code>myseq</code> . For example, <code>myseq=seq(-1, 1, by=0.5)</code> provides five options for direction changes. If the user specifies any single number instead of a sequence, (e.g., <code>myseq=1</code>) then <code>flexMeboot</code> will not change the directions of trends at all, but will modify the original <code>meboot</code> function to re-sample separately within several non-overlapping blocks, before joining them into resampled time series. This may be desirable for long series and for some applications. |

Details

`flexMeboot` uses non-overlapping blocks having only m observations. A trend $a + bt$ is replaced by $a + Bt$, where $B = \text{sample}(\text{myseq}) * b$.

Its steps are as follows:

1. Choose block size `segment` denoted here as m (default equal to $m = 5$) and divide the original time series x of length T into $k = \text{floor}(T/m)$ blocks or subsets. Note that when T/m is not an integer the k -th block will have a few more than m items. Hence let us denote the number of observations in each block as m which equals m for most blocks, except the k -th.
2. Regress each block having m observations as subsets of x on the set $\tau = 1, 2, \dots, m$, and store the intercept b_0 , the slope b_1 of τ and the residuals r .
3. Note that the positive (negative) sign of the slope b_1 in this regression determines the up (down) direction of the time series in that block. Hence the next step of the algorithm replaces b_1 by $B_1 = b_1 * w$, defined by a randomly chosen weight $\text{win}(-1, 0, 1)$. For example, when the random choice yields $w = -1$, the sign of b_1 is reversed. Our weighting independently injects some limited flexibility to the directions of values block segments of the original time series.

4. Reconstruct all time series blocks as: $b_0 + b_1 * w * \tau + r$, by adding back the residual r of the regression on τ .
5. The next step applies the function `meboot` to each block of time series—now having a modified trend—and create a large number, J , of resampled time series for each of the k blocks.
6. Sequentially join the J replicates of all k blocks or subsets together.

Value

A matrix containing by columns the bootstrapped replicated of the original data x .

References

Vinod, H.D. (2012), Constructing Scenarios of Time Heterogeneous Series for Stress Testing, Available at SSRN: <https://www.ssrn.com/abstract=1987879>.

See Also

`meboot`.

Examples

```
set.seed(235)
myseq <- seq(-1, 1, by = 0.5)
xx <- flexMeboot(x = AirPassengers, myseq = myseq, reps = 3)
matplot(cbind(AirPassengers, xx), type = "l")
```

force.clt

Enforce Central Limit Theorem

Description

Function to enforce the maximum entropy bootstrap resamples to satisfy the central limit theorem.

Usage

```
force.clt (x, ensemble)
```

Arguments

`x` a vector of data or a time series object.
`ensemble` a matrix or `mts` object containing resamples of the original data x .

Value

Revised matrix satisfying the central limit theorem.

Examples

```

set.seed(345)
out <- meboot(x=AirPassengers, reps=100, trim=0.10, reachbnd=FALSE, elaps=TRUE)
cm1 <- colMeans(out$ensemb)
# Note that the column means are somewhat non-normal
qqnorm(cm1)

clt.ens <- force.clt(x=AirPassengers, ensemble=out$ensemble)
cm2 <- colMeans(clt.ens)
# Note that the columns are closer to being normal
qqnorm(cm2)

```

meboot

*Generate Maximum Entropy Bootstrapped Time Series Ensemble***Description**

Generates maximum entropy bootstrap replicates for dependent data. (See details.)

Usage

```

meboot(x, reps=999, trim=list(trim=0.10, xmin=NULL, xmax=NULL), reachbnd=TRUE,
expand.sd=TRUE, force.clt=TRUE, scl.adjustment = FALSE, sym = FALSE,
elaps=FALSE, colsubj, coldata, coltimes, ...)

```

Arguments

| | |
|-----------------------------|---|
| <code>x</code> | vector of data, ts object or pdata.frame object. |
| <code>reps</code> | number of replicates to generate. |
| <code>trim</code> | a list object containing the elements: <code>trim</code> , the trimming proportion; <code>xmin</code> , the lower limit for left tail and <code>xmax</code> , the upper limit for right tail. |
| <code>reachbnd</code> | logical. If TRUE potentially reached bounds (<code>xmin</code> = smallest value - trimmed mean and <code>xmax</code> =largest value + trimmed mean) are given when the random draw happens to be equal to 0 and 1, respectively. |
| <code>expand.sd</code> | logical. If TRUE the standard deviation in the ensemble is expanded. See expand.sd . |
| <code>force.clt</code> | logical. If TRUE the ensemble is forced to satisfy the central limit theorem. See force.clt . |
| <code>scl.adjustment</code> | logical. If TRUE scale adjustment is performed to ensure that the population variance of the transformed series equals the variance of the data. |
| <code>sym</code> | logical. If TRUE an adjustment is performed to ensure that the ME density is symmetric. |
| <code>elaps</code> | logical. If TRUE elapsed time during computations is displayed. |

| | |
|----------|---|
| colsubj | the column in <code>x</code> that contains the individual index. It is ignored if the input data <code>x</code> is not a <code>pdata.frame</code> object. |
| coldata | the column in <code>x</code> that contains the data of the variable to create the ensemble. It is ignored if the input data <code>x</code> is not a <code>pdata.frame</code> object. |
| coltimes | an optional argument indicating the column that contains the times at which the observations for each individual are observed. It is ignored if the input data <code>x</code> is not a <code>pdata.frame</code> object. |
| ... | possible argument <code>fi.v</code> to be passed to <code>expand.sd</code> . |

Details

Seven-steps algorithm:

1. Sort the original data in increasing order and store the ordering index vector.
2. Compute intermediate points on the sorted series.
3. Compute lower limit for left tail (`xmin`) and upper limit for right tail (`xmax`). This is done by computing the `trim` (e.g. 10
4. Compute the mean of the maximum entropy density within each interval in such a way that the *mean preserving constraint* is satisfied. (Denoted as m_t in the reference paper.) The first and last interval means have distinct formulas. See Theil and Laitinen (1980) for details.
5. Generate random numbers from the $[0,1]$ uniform interval and compute sample quantiles at those points.
6. Apply to the sample quantiles the correct order to keep the dependence relationships of the observed data.
7. Repeat the previous steps several times (e.g. 999).

The scale and symmetry adjustments are described in Vinod (2013) referenced below.

In some applications, the ensembles must be ensured to be non-negative. Setting `trim$xmin = 0` ensures positive values of the ensembles. It also requires `force.clt = FALSE` and `expand.sd = FALSE`. These arguments are set to `FALSE` if `trim$xmin = 0` is defined and a warning is returned to inform that the value of those arguments were overwritten. Note: The choice of `xmin` and `xmax` cannot be arbitrary and should be cognizant of `range(x)` in data. Otherwise, if there are observations outside those bounds, the limits set by these arguments may not be met. If the user is concerned only with the trimming proportion, then it can be passed as argument simply `trim = 0.1` and the default values for `xmin` and `xmax` will be used.

Value

| | |
|-----------------------|--|
| <code>x</code> | original data provided as input. |
| <code>ensemble</code> | maximum entropy bootstrap replicates. |
| <code>xx</code> | sorted order stats (<code>xx[1]</code> is minimum value). |
| <code>z</code> | class intervals limits. |
| <code>dv</code> | deviations of consecutive data values. |
| <code>dvtrim</code> | trimmed mean of <code>dv</code> . |
| <code>xmin</code> | data minimum for <code>ensemble=xx[1]-dvtrim</code> . |

| | |
|----------|---|
| xmax | data x maximum for ensemble=xx[n]+dvtrim. |
| desintxb | desired interval means. |
| ordxx | ordered x values. |
| kappa | scale adjustment to the variance of ME density. |
| elaps | elapsed time. |

References

Vinod, H.D. (2013), Maximum Entropy Bootstrap Algorithm Enhancements. <https://www.ssrn.com/abstract=2285041>.

Vinod, H.D. (2006), Maximum Entropy Ensembles for Time Series Inference in Economics, *Journal of Asian Economics*, **17**(6), pp. 955-978

Vinod, H.D. (2004), Ranking mutual funds using unconventional utility theory and stochastic dominance, *Journal of Empirical Finance*, **11**(3), pp. 353-377.

Examples

```
## Ensemble for the AirPassenger time series data
set.seed(345)
out <- meboot(x=AirPassengers, reps=100, trim=0.10, elaps=TRUE)

## Ensemble for T=5 toy time series used in Vinod (2004)
set.seed(345)
out <- meboot(x=c(4, 12, 36, 20, 8), reps=999, trim=0.25, elaps=TRUE)
mean(out$sens) # ensemble mean should be close to sample mean 16
```

meboot.part

meboot Internal Function

Description

Internal function.

Usage

```
meboot.part (x, n, z, xmin, xmax, desintxb, reachbnd)
```

Arguments

| | |
|----------|---|
| x | vector of data, ts object or pdata.frame object. |
| n | length of x. |
| z | class intervals limits. |
| xmin | lower limit in the left tail. |
| xmax | upper limit in the left tail |
| desintxb | desired interval means. |
| reachbnd | logical. If TRUE potentially reached bounds (xmin = smallest value - trimmed mean and xmax=largest value + trimmed mean) are given when the random draw happens to be equal to 0 and 1, respectively. |

Value

A vector of resampled data.

See Also

[meboot.](#)

meboot.pdata.frame *Maximum Entropy Bootstrap for Panel Time Series Data*

Description

This function applies the maximum entropy bootstrapped in a panel of time series data.

Usage

```
meboot.pdata.frame(x, reps=999, trim=0.10, reachbnd=TRUE,
  expand.sd=TRUE, force.clt=TRUE, scl.adjustment = FALSE, sym = FALSE, elaps=FALSE,
  colsubj, coldata, coltimes, ...)
```

Arguments

| | |
|----------------|---|
| x | a pdata.frame object containing by columns: the individual index, an optional time index and a panel of time series data. |
| reps | number of replicates to generate for each subject in the panel. |
| trim | the trimming proportion. |
| reachbnd | logical. If TRUE potentially reached bounds (xmin = smallest value - trimmed mean and xmax=largest value + trimmed mean) are given when the random draw happens to be equal to 0 and 1, respectively. |
| expand.sd | logical. If TRUE the standard deviation in the ensemble is expanded. See expand.sd . |
| force.clt | logical. If TRUE the ensemble is forced to satisfy the central limit theorem. See force.clt . |
| scl.adjustment | logical. If TRUE scale adjustment is performed to ensure that the population variance of the transformed series equals the variance of the data. |
| sym | logical. If TRUE an adjustment is performed to ensure that the ME density is symmetric. |
| elaps | logical. If TRUE elapsed time during computations is displayed. |
| colsubj | the column in x that contains the individual index. |
| coldata | the column in x that contains the data of the variable to create the ensemble. |
| coltimes | an optional argument indicating the column that contains the times at which the observations for each individual are observed. |
| ... | possible argument fiv to be passed to expand.sd . |

Details

The observations in `x` should be arranged by individuals. The observations for each individual must be sorted by time.

The argument `colsubj` can be either a numeric or a character index indicating the individual or the time series to which each observation is related.

Only one variable can be replicated at a time, `coldata` must be of length one.

If the times at which observations are observed is provided specifying the column with the times through the argument `coltimes`, these times are used only to label the rows of the `data.frame` returned as output.

Value

A `data.frame` object of dimension: number of rows of `x` times number of replicates indicated in `reps`. The replicates for the panel of data are arranged by columns. Each replicate in each column is sorted with the same order established in the input `x`.

See Also

[meboot](#).

Examples

```
## Ensemble for a panel of series of stock prices
data("ullwan")
out <- meboot(ullwan, reps=99, colsubj=2, coldata=4)
```

mebootSpear

*Generate Maximum Entropy Bootstrapped Time Series Ensemble
Specifying Rank Correlation*

Description

Generates maximum entropy bootstrap replicates for dependent data specifying Spearman rank correlation coefficient between replicates series. (See details.)

Usage

```
mebootSpear(x, reps=999, setSpearman=1, drift=TRUE, trim=0.10,  
xmin=NULL, xmax=NULL, reachbnd=TRUE, expand.sd=TRUE, force.clt=TRUE,  
scl.adjustment = FALSE, sym = FALSE, elaps=FALSE, colsubj, coldata, coltimes,...)
```

Arguments

| | |
|-----------------------------|--|
| <code>x</code> | vector of data, <code>ts</code> object or <code>pdata.frame</code> object. |
| <code>reps</code> | number of replicates to generate. |
| <code>setSpearman</code> | The default setting <code>setSpearman=1</code> assumes that the user wants to generate replicates that are perfectly dependent on original time series. <code>setSpearman<1</code> admits less perfect (more realistic for some purposes) dependence. |
| <code>drift</code> | logical; TRUE default preserves the drift of the original series. |
| <code>trim</code> | the trimming proportion. |
| <code>xmin</code> | the lower limit for left tail. |
| <code>xmax</code> | the upper limit for right tail. |
| <code>reachbnd</code> | logical. If TRUE potentially reached bounds (<code>xmin</code> = smallest value - trimmed mean and <code>xmax</code> =largest value + trimmed mean) are given when the random draw happens to be equal to 0 and 1, respectively. |
| <code>expand.sd</code> | logical. If TRUE the standard deviation in the ensemble is expanded. See expand.sd . |
| <code>force.clt</code> | logical. If TRUE the ensemble is forced to satisfy the central limit theorem. See force.clt . |
| <code>scl.adjustment</code> | logical. If TRUE scale adjustment is performed to ensure that the population variance of the transformed series equals the variance of the data. |
| <code>sym</code> | logical. If TRUE an adjustment is performed to ensure that the ME density is symmetric. |
| <code>elaps</code> | logical. If TRUE elapsed time during computations is displayed. |
| <code>colsubj</code> | the column in <code>x</code> that contains the individual index. It is ignored if the input data <code>x</code> is not a <code>pdata.frame</code> object. |
| <code>coldata</code> | the column in <code>x</code> that contains the data of the variable to create the ensemble. It is ignored if the input data <code>x</code> is not a <code>pdata.frame</code> object. |
| <code>coltimes</code> | an optional argument indicating the column that contains the times at which the observations for each individual are observed. It is ignored if the input data <code>x</code> is not a <code>pdata.frame</code> object. |
| <code>...</code> | possible argument <code>fv</code> to be passed to expand.sd . |

Details

Seven-steps algorithm:

1. Sort the original data in increasing order and store the ordering index vector.
2. Compute intermediate points on the sorted series.
3. Compute lower limit for left tail (`xmin`) and upper limit for right tail (`xmax`). This is done by computing the `trim` (e.g. 10)
4. Compute the mean of the maximum entropy density within each interval in such a way that the *mean preserving constraint* is satisfied. (Denoted as m_t in the reference paper.) The first and last interval means have distinct formulas. See Theil and Laitinen (1980) for details.
5. Generate random numbers from the $[0,1]$ uniform interval and compute sample quantiles at those points.

6. Apply to the sample quantiles the correct order to keep the dependence relationships of the observed data.
7. Repeat the previous steps several times (e.g. 999).

The scale and symmetry adjustments are described in Vinod (2013) referenced below.

In some applications, the ensembles must be ensured to be non-negative. Setting `trim$xmin = 0` ensures positive values of the ensembles. It also requires `force.clt = FALSE` and `expand.sd = FALSE`. These arguments are set to `FALSE` if `trim$xmin = 0` is defined and a warning is returned to inform that the value of those arguments were overwritten. Note: The choice of `xmin` and `xmax` cannot be arbitrary and should be cognizant of `range(x)` in data. Otherwise, if there are observations outside those bounds, the limits set by these arguments may not be met. If the user is concerned only with the trimming proportion, then it can be passed as argument simply `trim = 0.1` and the default values for `xmin` and `xmax` will be used.

`setSpearman<1` is implemented with grid search near the desired value of the rank correlation coefficient, suggested by Fred Violen, a Ph.D. student at Fordham University and author of an R package NNS.

Value

| | |
|-----------------------|--|
| <code>x</code> | original data provided as input. |
| <code>ensemble</code> | maximum entropy bootstrap replicates. |
| <code>xx</code> | sorted order stats (<code>xx[1]</code> is minimum value). |
| <code>z</code> | class intervals limits. |
| <code>dv</code> | deviations of consecutive data values. |
| <code>dvtrim</code> | trimmed mean of <code>dv</code> . |
| <code>xmin</code> | data minimum for <code>ensemble=xx[1]-dvtrim</code> . |
| <code>xmax</code> | data x maximum for <code>ensemble=xx[n]+dvtrim</code> . |
| <code>desintxb</code> | desired interval means. |
| <code>ordxx</code> | ordered x values. |
| <code>kappa</code> | scale adjustment to the variance of ME density. |
| <code>elaps</code> | elapsed time. |

References

- Vinod, H.D. and Violen, F. (2020), Maximum Entropy Bootstrap and Improved Monte Carlo Simulations. https://papers.ssrn.com/sol3/papers.cfm?abstract_id=3621614.
- Vinod, H.D. (2013), Maximum Entropy Bootstrap Algorithm Enhancements. <https://www.ssrn.com/abstract=2285041>.
- Vinod, H.D. (2006), Maximum Entropy Ensembles for Time Series Inference in Economics, *Journal of Asian Economics*, **17**(6), pp. 955-978
- Vinod, H.D. (2004), Ranking mutual funds using unconventional utility theory and stochastic dominance, *Journal of Empirical Finance*, **11**(3), pp. 353-377.

Examples

```
## Ensemble for the AirPassenger time series data
set.seed(345)
out <- mebootSpear(x=AirPassengers, reps=100, xmin=0, setSpearman = 0)
cor(out$rowAvg, AirPassengers, method = "spearman") # rank-correlation should be close to 0
```

null.ci

Get Confidence Interval Around Specified NullZero Total

Description

Function to get two sided confidence interval around zero as the true value. Confidence interval is adjusted so that it covers the true zero $(1-\text{level})*100$ times. Symmetry is not assumed.

Usage

```
null.ci (x, level=0.95, null.value=0, type=8, ...)
```

Arguments

| | |
|------------|--|
| x | a vector of data. |
| level | confidence level. |
| null.value | a specified value of the null, e.g., 0. |
| type | type of quantile, a number between 1 and 9. See quantile . |
| ... | further arguments passed to or from other methods. |

Value

Lower limit and upper limit of the confidence interval.

Examples

```
x <- runif(25, 0, 1)
null.ci(x)
```

`olsHALL.b`*OLS regression model for consumption*

Description

Compute OLS coefficients in a regression model for the consumption variable. See details.

Usage

```
olsHALL.b (y, x)
```

Arguments

| | |
|----------------|---|
| <code>y</code> | dependent variable (consumption). |
| <code>x</code> | regressor variable (disposable income). |

Details

The regression model is: $c(t) = b_1 + b_2*c(t) + b_3*y(t-1) + u(t)$, where 'c' is consumption and 'y' is disposable income.

This function is intended to speed up the ME bootstrap procedure for inference. Instead of using the `lm` or `dynlm` interfaces the function calls directly to the Fortran procedure 'dqrls'.

Value

Coefficient estimates by OLS.

Examples

```
data("USconsum")
USconsum <- log(USconsum)

# lm interface
lmcf1 <- lm(USconsum[-1,1] ~ USconsum[-51,1] + USconsum[-51,2])
coefficients(lmcf1)

# dynlm interface
library("dynlm")
lmcf2 <- dynlm(consum ~ L(consum, 1) + L(dispinc, 1), data=USconsum)
coefficients(lmcf2)

# olsHALL.b
olsHALL.b(y=USconsum[,1], x=USconsum[,2])
```

ullwan

Data about Some of the S&P 500 Stock Prices

Description

This data set collects information about seven S&P 500 stocks whose market capitalization exceeds \$27 billion. The seven companies are labelled as ABT, AEG, ATI, ALD, ALL, AOL and AXP. For each company data from May 1993 to November 1997 are available (469 observations).

Usage

```
data (ullwan)
```

Format

The data are stored in an object of class `pdata.frame` (a `data.frame` class with further attributes useful for panel data from the `plm` package) and `data.frame`.

Details

The following information is contained by columns:

- `Subj`: Company index.
- `Tim`: Times at which the data were observed (on a monthly basis).
- `MktVal`: Market capitalization.
- `Price`: Stock prices.
- `Pupdn`: Binary variable, takes the value 1 if there is a turning point (a switch from a bull to a bear market or vice versa) and 0 otherwise.
- `Tb3`: Interest on 3-month Treasury bills.

Source

Compustat database.

References

Yves Croissant (2005). `plm`: Linear models for panel data. R package version 0.1-2.

USconsum

Consumption and Disposable Income Data (Annual 1948-1998)

Description

Data set employed in Murray (2006, pp.46-47, 799-801) to discuss the Keynesian consumption function on the basis of the Friedman's permanent income hypothesis and Robert Hall's model.

Usage

data (USconsum)

Format

A .rda file storing the data as an mts object.

Details

Annual data. Available time series: (Each corresponding label in the list object appears in quotes.)

- `consum`: consumption per capita in thousands of dollars (1948-1998). Log of this variable is the dependent variable and its lagged value is a regressor in Murray's Table 18.1.
- `dispinc`: disposable income per capita in thousands of dollars (1948-1998). Lagged value of the Log of this variable is the second regressor (proxy for permanent income) in Murray's Table 18.1.

Source

Robert Hall's data for 1948 to 1977 extended by Murray to 1998 by using standard sources for US macroeconomic data from government publications (U.S. Bureau of Labor Statistics for population data, U.S. Bureau of Economic Analysis for income data, U.S. Bureau of Labor Statistics for consumption data).

References

Murray, M.P. (2006), *Econometrics. A modern introduction*, New York: Pearson Addison Wesley.

| | |
|--------|---|
| USfygt | <i>Long-term Treasury Bond Rates and Deficit Data Set (Annual 1948-200)</i> |
|--------|---|

Description

Data set employed in Murray (2006, pp.795-797) to test the null hypothesis that per capita federal deficits explain long-term Treasury bond interest rates based on the Stock and Watson's dynamic OLS model.

Usage

```
data (USfygt)
```

Format

A .rda file storing the data as an `mts` object.

Details

Annual data. Available time series: (Each corresponding label in the list object appears in quotes.)

- "dy": mean changes in real per capita income (1949-1998).
- "fygt1": shorth-term (one-year) Treasury bond interest rates (1953-1998).
- "fygt10": long-term (ten-year) Treasury bond interest rates (1953-2000).
- "infl": inflation (1949-2000).
- "usdef": per capita real federal deficit (1948-2000).
- "reallir": real long term interest rates (not used in Murray's Table 18.12).
- "realsir": real short term interest rates (not used in Murray's Table 18.12).

Source

Data was made available by James Stock and Mark Watson to readers of their famous *Econometrica* paper, 1993, 61, pp 783-820, who in turn used standard sources for US macroeconomic data from government publications.

References

Murray, M.P. (2006), *Econometrics. A modern introduction*, New York: Pearson Addison Wesley.

`zero.ci`*Get Confidence Interval Around Zero*

Description

Function to get two sided confidence interval around zero as the true value. Confidence interval is adjusted so that it covers the true zero $(1-\text{conf1})*100$ times. Symmetry is not assumed.

Usage

```
zero.ci (x, conf1=0.05)
```

Arguments

| | |
|--------------------|-------------------|
| <code>x</code> | a vector of data. |
| <code>conf1</code> | confidence level. |

Value

| | |
|--------------------|---|
| <code>bnlo</code> | count of number of items below lower limit. |
| <code>bnup</code> | count of number of items above upper limit. |
| <code>lolim</code> | lower limit of the confidence interval. |
| <code>uplim</code> | upper limit of the confidence interval. |

Examples

```
x <- runif(25, 0, 1)
zero.ci(x)
```

Index

* datagen

ullwan, 16
USconsum, 17
USfygt, 18

* ts

checkConv, 2
elapsedtime, 3
expand.sd, 4
flexMeboot, 5
force.clt, 6
meboot, 7
meboot.part, 9
meboot.pdata.frame, 10
mebootSpear, 11
null.ci, 14
zero.ci, 19

checkConv, 2
criterion, 3

elapsedtime, 3
expand.sd, 4, 7, 8, 10, 12

flexMeboot, 5
force.clt, 5, 6, 7, 10, 12

meboot, 3, 5, 6, 7, 10, 11
meboot.part, 9
meboot.pdata.frame, 10
mebootSpear, 11

null.ci, 14

olsHALL.b, 15

proc.time, 4

quantile, 14

ullwan, 16
USconsum, 17
USfygt, 18

zero.ci, 19