

Package: scaleboot (via r-universe)

October 16, 2024

Type Package

Title Approximately Unbiased P-Values via Multiscale Bootstrap

Version 1.0-1

Date 2019-12-03

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Depends R (>= 3.5.0)

Imports pvcclus (>= 2.2-0), mvtnorm

Description Calculating approximately unbiased (AU) p-values from multiscale bootstrap probabilities. See Shimodaira (2004) <doi:10.1214/009053604000000823>, Shimodaira (2008) <doi:10.1016/j.jspi.2007.04.001>, Terada and Shimodaira (2017) <arXiv:1711.00949>, and Shimodaira and Terada (2019) <doi.org/10.3389/fevo.2019.00174>.

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URL <http://stat.sys.i.kyoto-u.ac.jp/prog/scaleboot/>

NeedsCompilation no

Date/Publication 2019-12-04 16:00:02 UTC

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

RemoteUrl <https://github.com/cran/scaleboot>

RemoteRef HEAD

RemoteSha c0e40bf17cff7c554f92a68b7b9aadb385b8b60d

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scaleboot-package *Approximately Unbiased P-values via Multiscale Bootstrap*

Description

Calculating approximately unbiased (AU) p-values from multiscale bootstrap probabilities.

Details

For a complete list, use `library(help="scaleboot")`.

The methodology is described in Shimodaira (2008). For the use of `scaleboot`, Shimodaira (2008) may be referenced.

Further information is available in the following vignette:

`usesb` Multiscale Bootstrap using Scaleboot Package

Author(s)

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I thank Paul A. Sheridan for his comments to improve the manual pages.

References

Shimodaira, H. (2008). Testing Regions with Nonsmooth Boundaries via Multiscale Bootstrap, *Journal of Statistical Planning and Inference*, 138, 1227-1241. (<http://dx.doi.org/10.1016/j.jspi.2007.04.001>).

`coef`*Extract Model Coefficients*

Description

Extract the estimated parameters from "scaleboot" or "scalebootv" objects.

Usage

```
## S3 method for class 'scaleboot'  
coef(object, sd=FALSE, ...)  
  
## S3 method for class 'scalebootv'  
coef(object, ...)
```

Arguments

<code>object</code>	an object used to select a method.
<code>...</code>	further arguments passed to or from other methods.
<code>sd</code>	logical. Should standard errors be returned as well?

Value

The `coef` method for the class "scaleboot" returns a matrix consisting of row vectors of beta's for models. If `sd=TRUE`, it returns a list with components `estimate` and `sd` for the beta matrix and its standard error respectively.

Author(s)

Hidetoshi Shimodaira

See Also

[sbfit](#)

Examples

```
data(mam15)  
a <- mam15.relltest[["t4"]] # an object of class "scaleboot"  
coef(a) # print the estimated beta values  
coef(a, sd=TRUE) # with sd
```

interface *Interface to External Packages*

Description

Interface for other packages such as CONSEL (phylogenetic inference), and pvclust (hierarchical clustering)

Usage

```
read.mt(file, tlab="t")

read.ass(file, identity=TRUE, tlab="t", elab="e")

read.cnt(file)

## S3 method for class 'pvclust'
sbfit(x, ...)

sbpvclust(x, mbs, k=3, select="average", ...)
```

Arguments

file	character of a file name to be read.
identity	logical. Should the identity association be included?
tlab	character for basename of tree labels.
elab	character for basename of edge labels.
x	an object of class "pvclust".
mbs	an object of class "scalebootv".
k	numeric of k for a AU p-value.
select	character of model name (such as "poly.3") or one of "average" and "best".
...	further arguments passed to or from other methods.

Details

CONSEL is a program package consisting of small programs written in the C language for assessing the confidence of phylogenetic tree selection. Some functions for interfacing with CONSEL are: `read.mt`, `read.ass`, and `read.cnt` for reading, respectively, `mt`, `ass`, and `cnt` format. Once `mt` file is read, we can calculate improved versions of approximately unbiased p-values by `relltestin` `scaleboot` instead of CONSEL.

`pvclust` is a R package for hierarchical clustering with p-values. Functions for interface to `pvclust` are: `sbfit` method for an object of class "pvclust" to convert it to "scalebootv" class, and `sbpvclust` for writing back the result to a "pvclust" object with additional class "sbclust". `plot` method for class "sbclust" overwrites that for "pvclust".

Value

`read.mt` returns a matrix of dimension sequence-length by tree-number. If `identity=FALSE`, then `read.ass` returns a list containing components `x` for edge→tree associations and `y` for tree→edge associations. If `identity=TRUE`, `read.ass` returns a list vector of edge→tree associations, where the identity associations for tree→tree are included. `read.cnt` returns a list containing components `bps`, `nb`, and `sa` to be used for `sbfit`. The list also contains components `cnt`, `id`, and `val`.

`sbfit.pvclust` returns an object of class "scalebootv". `sbpvclust` returns an object of class "sbclust" added to the class "pvclust".

Author(s)

Hidetoshi Shimodaira

References

Shimodaira, H. and Hasegawa, M. (2001). CONSEL: for assessing the confidence of phylogenetic tree selection, *Bioinformatics*, 17, 1246-1247 (software is available from <http://stat.sys.i.kyoto-u.ac.jp/prog/consel/>).

Suzuki, R. and Shimodaira, H. (2006). pvclust: An R package for hierarchical clustering with p-values, *Bioinformatics*, 22, 1540-1542 (software is available from CRAN or <http://stat.sys.i.kyoto-u.ac.jp/prog/pvclust/>).

See Also

[lung73](#), [mam15](#), [relltest](#)

Examples

```
## replace au/bp entries in pvclust object
## see help(lung73) for details
library(pvclust)
data(lung73)
plot(lung73.pvclust) # draw dendrogram of pvclust object
pvrect(lung73.pvclust)
lung73.new <- sbpvclust(lung73.pvclust, lung73.sb) # au <- k.3
plot(lung73.new) # redraw dendrogram with the new au/bp values
pvrect(lung73.new)

## Not run:
## reading CONSEL files
## sample files are found in mam15 subdirectory
## see help(mam15) for details
mam15.mt <- read.mt("mam15.mt")
mam15.ass <- read.ass("mam15.ass")
mam15.cnt <- read.cnt("mam15.cnt")

## End(Not run)
```

lung73

Clustering of 73 Lung Tumors

Description

Bootstrapping hierarchical clustering of the DNA microarray data set of 73 lung tissue samples each containing 916 observed genes.

Usage

```
data(lung73)
```

Format

lung73.pvclust and lung.pvclust are objects of class "pvclust" defined in **pvclust** of Suzuki and Shimodaira (2006).

lung73.sb and lung.sb are an object of class "scalebootv" of length 72.

Details

The microarray dataset of Garber et al. (2001) is reanalyzed in Suzuki and Shimodaira (2006), and is found in data(lung) of the **pvclust** package. We reanalyze it, again, by the script shown in Examples. The result of pvclust is stored in lung73.pvclust and lung.pvclust, and model fitting to bootstrap probabilities by the **scaleboot** package is stored in lung73.sb and lung.sb. A wide scale range is used in lung73.pvclust and lung73.sb, and the default scale range of pvclust is used in lung.pvclust and lung.sb. The microarray dataset is not included in data(lung73), but it is found in data(lung) of the **pvclust** package.

Source

Garber, M. E. et al. (2001) Diversity of gene expression in adenocarcinoma of the lung, *Proceedings of the National Academy of Sciences*, 98, 13784-13789 (dataset is available from http://genome-www.stanford.edu/lung_cancer/adeno/).

References

Suzuki, R. and Shimodaira, H. (2006). pvclust: An R package for hierarchical clustering with p-values, *Bioinformatics*, 22, 1540-1542 (software is available from CRAN or <http://stat.sys.i.kyoto-u.ac.jp/prog/pvclust/>).

See Also

[sbpvclust](#), [sbfite.pvclust](#)

Examples

```

## Not run:
## Parallel setup
library(parallel)
length(cl <- makeCluster(detectCores()))
## script to create lung73.pvclust and lung73.sb
## multiscale bootstrap resampling of hierarchical clustering
library(pvclust)
data(lung)
### default pvclust scales
lung.pvclust <- pvclust(lung, nboot=10000, parallel=cl)
lung.sb <- sbfit(lung.pvclust,cluster=cl) # model fitting
### wider range of scales than pvclust default
sa <- 9^seq(-1,1,length=13)
lung73.pvclust <- pvclust(lung,r=1/sa,nboot=10000,parallel=cl)
lung73.sb <- sbfit(lung73.pvclust,cluster=cl) # model fitting

## End(Not run)

## replace si/au/bp entries in pvclust object
library(pvclust)
data(lung73) # loading the previously computed bootstrap

### the original pvclust result
plot(lung.pvclust, print.pv = c("si", "au", "bp"), cex=0.5, cex.pv=0.5)
pvrect(lung.pvclust, pv="si") # (default pvclust uses pv="au")

### default pvclust scales with p-values of k=2
lung.k2 <- sbpvclust(lung73.pvclust,lung73.sb, k=2)
plot(lung.k2, print.pv = c("si", "au", "bp"), cex=0.5, cex.pv=0.5)
pvrect(lung.k2, pv="si")

### wider scales with p-values of k=3 (default of scaleboot)
lung73.k3 <- sbpvclust(lung73.pvclust,lung73.sb)
plot(lung73.k3, print.pv = c("si", "au", "bp"), cex=0.5, cex.pv=0.5)
pvrect(lung73.k3, pv="si")

## diagnostics of fitting

### diagnose edges 61,...,69
lung73.sb[61:69] # print fitting details
plot(lung73.sb[61:69]) # plot curve fitting
summary(lung73.sb[61:69]) # print raw(=bp)/si/au p-values

### diagnose edge 67
lung73.sb[[67]] # print fitting
plot(lung73.sb[[67]],legend="topleft") # plot curve fitting
summary(lung73.sb[[67]]) # print au p-values

```

mam15

*Mammal Phylogenetic Analysis for 15 trees***Description**

Phylogenetic analysis of six mammal species for 15 trees and 105 trees.

Usage

```
data(mam15)
```

Format

mam15.mt is a matrix of size 3414 * 15. The (i,j) element is the site-wise log-likelihood value at site-i for tree-j for i=1,...,3414, and j=1,...,15. They are constrained trees with clade (cow, seal) being fixed.

mam15.ass is a list of length 25 for association vectors. The components are t1, t2, ..., t15 for trees, and e1, e2, ..., e10 for edges.

mam15.relltest is an object of class "relltest" of length 25.

mam15.aux is a list of tree topologies (tpl), clade patterns (cld), taxa names(tax).

mam105.mt, mam105.ass, mam105.rellst, mam105.aux are those for 105 unconstrained trees.

mam26.mt, mam26.ass, mam26.aux are those for 26 trees including the 15 constrained trees, 10 partially resolved trees corresponding to the 10 internal edges, and the star topology.

Details

An example of phylogenetic analysis of six mammal species: Homo sapiens (human), Phoca vitulina (harbor seal), Bos taurus (cow), Oryctolagus cuniculus (rabbit), Mus musculus (mouse), Didelphis virginiana (opossum). The data is stored in the file 'mam15.aa', which contains amino acid sequences of length N=3414 for the six species obtained from mtDNA (see Note below). Here we fix (Phovi,Bosta) as a group of taxa. With this constraint, we consider 15 tree topologies of the six mammals as stored in the file 'mam15.tpl';

```
((Homsa, (Phovi, Bosta)), Orycu, (Musmu, Didvi)); t1
(Homsa, Orycu, ((Phovi, Bosta), (Musmu, Didvi))); t2
(Homsa, ((Phovi, Bosta), Orycu), (Musmu, Didvi)); t3
(Homsa, (Orycu, Musmu), ((Phovi, Bosta), Didvi)); t4
((Homsa, (Phovi, Bosta)), (Orycu, Musmu), Didvi); t5
(Homsa, ((Phovi, Bosta), (Orycu, Musmu)), Didvi); t6
(Homsa, (((Phovi, Bosta), Orycu), Musmu), Didvi); t7
(((Homsa, (Phovi, Bosta)), Musmu), Orycu, Didvi); t8
(((Homsa, Musmu), (Phovi, Bosta)), Orycu, Didvi); t9
(Homsa, Orycu, (((Phovi, Bosta), Musmu), Didvi)); t10
(Homsa, (((Phovi, Bosta), Musmu), Orycu), Didvi); t11
((Homsa, ((Phovi, Bosta), Musmu)), Orycu, Didvi); t12
```



```
(Homsa,Orycu,(((Phovi,Bosta),Didvi),Musmu)); t13
((Homsa,Musmu),Orycu,((Phovi,Bosta),Didvi)); t14
((Homsa,Musmu),((Phovi,Bosta),Orycu),Didvi); t15
```

The log-likelihood values are calculated using the PAML software (Ziheng 1997) for phylogenetic inference. The two files ‘mam15.aa’ and ‘mam15.tpl’ are fed into PAML to generate the file ‘mam15.lnf’ of site-wise log-likelihood values.

Using the CONSEL software (Shimodaira and Hasegawa 2001), we convert ‘mam15.lnf’ and ‘mam15.tpl’ to a format suitable for the **scaleboot** package. We do not use CONSEL for calculating AU p-values, but use it only for file conversion. We type

```
seqmt --paml mam15.lnf
treeass --outgroup 6 mam15.tpl > mam15.log
```

The first line above generates ‘mam15.mt’, which is a simple text file containing a matrix of site-wise log-likelihood values. The second line above generates ‘mam15.ass’ and ‘mam15.log’, which contain information regarding which edges are included in a tree. A part of ‘mam15.log’ is as follows.

```
# leaves: 6
6
 1 Homsa
 2 Phovi
 3 Bosta
 4 Orycu
 5 Musmu
 6 Didvi

# base edges: 10
10 6

    123456
 1 +++--- ;
 2 ++++-- ;
 3 +---+- ;
 4 -++++- ;
 5 ---++- ;
 6 +---+- ;
 7 -++++- ;
 8 +++-+- ;
 9 +----+ ;
10 -++-+- ;
```

The above defines edges named e1,...e10 (base edges) as clusters for six mammal species. For example, e1 = +++— = (Homsa, Phovi, Bosta).

The converted files are read by the **scaleboot** package in R:

```
mam15.mt <- read.mt("mam15.mt")
mam15.ass <- read.ass("mam15.ass")
```

mam15.mt is a matrix of size 3414 * 6 for the site-wise log-likelihood values. For testing trees, we need only mam15.mt. mam15.ass is used for testing edges, and it is a list of length 25 for association vectors for t1,t2,...,t15, and e1,e2,...,e10. For example, mam15.ass\$t1 = 1, indicating tree "t1" is included in tree "t1", and mam15.ass\$e1 = c(1, 5, 8), indicating edge "e1" is included in trees "t1", "t5", and "t8".

Multiscale bootstrap resampling is performed by the function `relltest`. The simplest way to get AU p-values for trees is:

```
mam15.trees <- relltest(mam15.mt) # resampling and fitting
summary(mam15.trees) # calculates AU p-values
```

The `relltest` returns an object of class "relltest". It calls the function `scaleboot` internally with the number of bootstrap replicates `nb=10000`, and takes about 20 mins. Typically, `nb=10000` is large enough, but it would be safe to use larger value, say `nb=100000` as in the examples below.

Note that the default value of scales in `relltest` has a much wider range than that of CONSEL. It is `sa=9^seq(-1,1,length=13)` for `relltest`, and is `sa=1/seq(from=0.5,to=1.4,by=0.1)` for CONSEL.

The `mam15.relltest` object in `data(mam15)` is similar to `mam15.trees` above, but is also calculated for edges using `mam15.ass`. We can extract the result for trees by

```
mam15.trees <- mam15.relltest[1:15]
```

The results for trees stored in the `mam15.trees` object above are in the order specified in the columns of `mam15.mt`. To sort it by increasing order of the log-likelihood difference, we can type

```
stat <- attr(mam15.trees,"stat") # the log-likelihood differences
o <- order(stat) # sort it in increasing order
mam15.trees <- mam15.trees[o] # same as mam15.trees in Examples
```

Results of the fitting are shown by using the print method.

```
> mam15.trees
```

Test Statistic, and Shimodaira-Hasegawa test:

```
  stat  shtest
t1  -2.66 94.51 (0.07)
t3   2.66 80.25 (0.13)
t2   7.40 57.85 (0.16)
t5  17.57 17.30 (0.12)
t6  18.93 14.32 (0.11)
t7  20.11 11.49 (0.10)
t4  20.60 10.98 (0.10)
t15 22.22  7.34 (0.08)
t8  25.38  3.31 (0.06)
t14 26.32  3.29 (0.06)
```

t13 28.86 1.71 (0.04)
 t9 31.64 0.61 (0.02)
 t11 31.75 0.57 (0.02)
 t10 34.74 0.20 (0.01)
 t12 36.25 0.12 (0.01)

Multiscale Bootstrap Probabilities (percent):

	1	2	3	4	5	6	7	8	9	10	11	12	13
t1	86	81	77	73	68	63	58	52	46	41	36	31	28
t3	14	19	23	27	30	32	32	31	30	27	25	22	20
t2	0	0	0	0	1	2	4	5	7	9	10	11	11
t5	0	0	0	0	0	1	1	2	3	5	6	6	7
t6	0	0	0	0	1	2	3	5	6	7	8	9	9
t7	0	0	0	0	0	0	0	1	2	3	4	5	5
t4	0	0	0	0	0	1	2	3	4	4	5	6	6
t15	0	0	0	0	0	0	0	0	1	1	2	2	3
t8	0	0	0	0	0	0	0	0	0	0	1	1	1
t14	0	0	0	0	0	0	0	1	1	2	3	4	4
t13	0	0	0	0	0	0	0	0	0	0	1	1	2
t9	0	0	0	0	0	0	0	0	0	0	0	1	1
t11	0	0	0	0	0	0	0	0	0	0	0	1	1
t10	0	0	0	0	0	0	0	0	0	0	0	0	0
t12	0	0	0	0	0	0	0	0	0	0	0	0	0

Numbers of Bootstrap Replicates:

1	2	3	4	5	6	7	8	9	10	11	12	13
1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05

Scales (Sigma Squared):

1	2	3	4	5	6	7	8	9	10	11	12	13
0.1111	0.1603	0.2311	0.3333	0.4808	0.6933	1	1.442	2.080	3	4.327	6.241	9.008

AIC values of Model Fitting:

	poly.1	poly.2	poly.3	sing.3
t1	89483.40	964.33	964.75	966.33
t3	75434.97	1750.22	1306.50	1752.22
t2	29361.29	403.41	36.33	-6.21
t5	23893.19	260.44	-0.22	-14.11
t6	35791.26	330.50	4.31	-2.49
t7	15221.10	93.59	-10.33	-12.04
t4	29790.60	453.95	5.22	-7.57
t15	6874.98	46.16	-10.48	-17.08
t8	1747.13	-6.88	-12.39	-13.68
t14	10905.94	131.48	2.65	-10.79
t13	3411.26	27.66	-8.30	-15.14
t9	1494.58	19.46	-13.78	-15.86
t11	914.42	-19.65	-19.71	-19.61
t10	259.68	-14.79	-17.27	-16.76

```
t12 178.79 -19.19 -19.61 -19.30
```

The AU p-values are shown by the summary method.

```
> summary(mam15.trees)
```

Corrected P-values (percent):

	raw	k.1	k.2	k.3	model	aic
t1	57.58 (0.16)	56.16 (0.04)	74.55 (0.05)	74.55 (0.05)	poly.2	964.33
t3	31.86 (0.15)	30.26 (0.05)	46.41 (0.09)	45.33 (0.13)	poly.3	1306.50
t2	3.68 (0.06)	3.68 (0.03)	12.97 (0.20)	16.12 (0.45)	sing.3	-6.21
t5	1.34 (0.04)	1.33 (0.02)	7.92 (0.25)	10.56 (0.56)	sing.3	-14.11
t6	3.18 (0.06)	3.15 (0.02)	13.15 (0.21)	15.86 (0.44)	sing.3	-2.49
t7	0.49 (0.02)	0.52 (0.01)	3.66 (0.21)	4.75 (0.42)	sing.3	-12.04
t4	1.55 (0.04)	1.53 (0.02)	10.54 (0.27)	14.84 (0.66)	sing.3	-7.57
t15	0.08 (0.01)	0.07 (0.00)	1.11 (0.19)	1.85 (0.48)	sing.3	-17.08
t8	0.00 (0.00)	0.00 (0.00)	0.04 (0.03)	0.07 (0.07)	sing.3	-13.68
t14	0.22 (0.01)	0.23 (0.01)	2.76 (0.26)	4.59 (0.71)	sing.3	-10.79
t13	0.02 (0.00)	0.01 (0.00)	0.50 (0.20)	1.30 (0.83)	sing.3	-15.14
t9	0.00 (0.00)	0.00 (0.00)	0.23 (0.05)	1.41 (0.29)	sing.3	-15.86
t11	0.00 (0.00)	0.00 (0.00)	0.00 (0.00)	0.00 (0.00)	poly.3	-19.71
t10	0.00 (0.00)	0.00 (0.00)	0.00 (0.00)	0.00 (0.00)	poly.3	-17.27
t12	0.00 (0.00)	0.00 (0.00)	0.00 (0.00)	0.00 (0.00)	poly.3	-19.61

The p-values for 15 trees are shown above. "raw" is the ordinary bootstrap probability, "k.1" is equivalent to "raw" but calculated from the multiscale bootstrap, "k.2" is equivalent to the third-order AU p-value of CONSEL, and finally "k.3" is an improved version of AU p-value.

The details for each tree are shown by extracting the corresponding element. For example, details for the seventh largest tree in the log-likelihood value ("t4") is obtained by

```
> mam15.trees[[7]] # same as mam15.trees$t4
```

Multiscale Bootstrap Probabilities (percent):

1	2	3	4	5	6	7	8	9	10	11	12	13
0.00	0.00	0.01	0.08	0.27	0.80	1.55	2.55	3.58	4.42	5.22	6.00	6.38

Numbers of Bootstrap Replicates:

1	2	3	4	5	6	7	8	9	10	11	12	13
1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05	1e+05

Scales (Sigma Squared):

1	2	3	4	5	6	7	8	9	10	11	12	13
0.1111	0.1603	0.2311	0.3333	0.4808	0.6933	1	1.442	2.080	3	4.327	6.241	9.008

Coefficients:

	beta0	beta1	beta2
poly.1	2.8388 (0.0048)		
poly.2	1.8556 (0.0061)	0.3259 (0.0019)	
poly.3	1.7157 (0.0085)	0.4508 (0.0061)	-0.0152 (0.0007)

```
sing.3 1.6178 (0.0153) 0.5435 (0.0143) 0.3261 (0.0201)
```

```
Model Fitting:
```

	rss	df	pfit	aic
poly.1	29814.60	12	0.0000	29790.60
poly.2	475.95	11	0.0000	453.95
poly.3	25.22	10	0.0050	5.22
sing.3	12.43	10	0.2571	-7.57

```
Best Model: sing.3
```

```
> summary(mam15.trees[[7]])
```

```
Raw Bootstrap Probability: 1.55 (0.04)
```

```
Corrected P-values (percent):
```

	k.1	k.2	k.3	aic
poly.1	0.23 (0.00)	0.23 (0.00)	0.23 (0.00)	29790.60
poly.2	1.46 (0.02)	6.30 (0.09)	6.30 (0.09)	453.95
poly.3	1.57 (0.02)	9.50 (0.21)	10.57 (0.27)	5.22
sing.3	1.53 (0.02)	10.54 (0.27)	14.84 (0.66)	-7.57

```
Best Model: sing.3
```

```
> plot(mam15.trees[[7]],legend="topleft")
```

The plot diagnostics found in the bottom line are especially useful for confirming which model is fitting best.

See other examples below.

Note

Dataset files for phylogenetic inference are found at <http://github.com/shimo-lab/scaleboot>. Look at the subdirectory 'dataset/mam15-files'. This dataset was originally used in Shimodaira and Hasegawa (1999).

Source

H. Shimodaira and M. Hasegawa (1999). Multiple comparisons of log-likelihoods with applications to phylogenetic inference, *Molecular Biology and Evolution*, 16, 1114-1116.

References

Yang, Z. (1997). PAML: a program package for phylogenetic analysis by maximum likelihood, *Computer Applications in BioSciences*, 13:555-556 (software is available from <http://abacus.gene.ucl.ac.uk/software/paml.html>).

Shimodaira, H. and Hasegawa, M. (2001). CONSEL: for assessing the confidence of phylogenetic tree selection, *Bioinformatics*, 17, 1246-1247 (software is available from <http://stat.sys.i.kyoto-u.ac.jp/prog/consel/>).

See Also

[mam105](#), [relltest](#), [summary.scalebootv](#), [read.mt](#), [read.ass](#).

Examples

```
data(mam15)

## show the results for trees and edges
mam15.relltest # print stat, shtest, bootstrap probabilities, and AIC
summary(mam15.relltest) # print AU p-values

## Not run:
## simpler script to create mam15.trees
mam15.mt <- read.mt("mam15.mt")
mam15.ass <- read.ass("mam15.ass")
mam15.trees <- relltest(mam15.mt,nb=100000)

## End(Not run)

## Not run:
## script to create mam15.relltest
mam15.mt <- read.mt("mam15.mt")
mam15.ass <- read.ass("mam15.ass")
mam15.relltest <- relltest(mam15.mt,nb=100000,ass=mam15.ass)

## End(Not run)

## Not run:
## Parallel version of the above script (but different in random seed)
## It took 13 mins (40 cpu's of Athlon MP 2000+)
mam15.mt <- read.mt("mam15.mt")
mam15.ass <- read.ass("mam15.ass")
library(parallel)
cl <- makeCluster(40)
mam15.relltest <- relltest(mam15.mt,nb=100000,ass=mam15.ass,cluster=cl)

## End(Not run)
```

Description

plot method for class "scaleboot".

Usage

```
## S3 method for class 'scaleboot'
plot(x, models=NULL, select=NULL, sort.by=c("aic","none"),
     k=NULL, s=NULL, sp=NULL, lambda=NULL, bpk=NULL,
     xval = c("square", "inverse","sigma"),
     yval = c("psi", "zvalue", "pvalue"), xlab = NULL,
     ylab = NULL,log.xy = "", xlim = NULL, ylim = NULL,
     add = F, length.x = 300, main=NULL,
     col =1:6, lty = 1:5, lwd = par("lwd"), ex.pch=2:7,
     pch = 1, cex = 1, pt.col = col[1],pt.lwd = lwd[1],
     legend.x = NULL, inset = 0.1, cex.legend=1,...)

## S3 method for class 'summary.scaleboot'
plot(x, select="average",
     k=x$parex$k,s=x$parex$s,sp=x$parex$sp,lambda=x$parex$lambda, ...)

## S3 method for class 'scalebootv'
plot(x,models=attr(x,"models"),sort.by="none",...)

## S3 method for class 'summary.scalebootv'
plot(x, select="average",...)

## S3 method for class 'scaleboot'
lines(x,z,models=names(x$fi), k=NULL,s=NULL,sp=NULL,lambda=NULL,
      bpk=NULL, length.x=z$length.x, col=z$col,lty=z$lty,lwd=z$lwd,... )

sblegend(x="topright",y=NULL,z,inset=0.1,...)

sbplotbeta(beta, p=0.05, col.contour=c("blue","red","green"),
            drawcontours = TRUE, drawlabels = TRUE,
            labcex=1,length=100, cex=1, col="black",
            xlim=NULL, ylim=NULL, lim.countourexpend=0 )
```

Arguments

x	an object used to select a method. For sblegend, x is a numeric or character such as "lefttop" or "righttop", which is passed to legend.
models	character vector of model names. Numeric is also allowed.
select	"average", "best", or one of the fitted models.
sort.by	"aic" or "none".
k	k for extrapolation.
s	s for extrapolation.
sp	sp for extrapolation.

lambda	a numeric of specifying the type of p-values; Bayesian (lambda=0) Frequentist (lambda=1).
bpk	(experimental for 2-step bootstrap)
xval	specifies x-axis. "square" for σ^2 , "inverse" for $1/\sigma$, "sigma" for σ .
yval	specifies y-axis. "zvalue" for $\psi(\sigma^2 \beta)/\sigma$ or $qnorm(1-bp[i])$, "pvalue" for $1 - \Phi(\psi(\sigma^2 \beta)/\sigma)$ or $bp[i]$, "psi" for $\psi(\sigma^2 \beta)$ or $\sqrt{sa[i]}*qnorm(1-bp[i])$.
xlab	label for x-axis.
ylab	label for y-axis.
log.xy	character to specify log-scale. "", "x", "y", or "xy".
xlim	range for x-axis.
ylim	range for y-axis.
add	logical for adding another plot.
length.x	the number of segments to draw curves.
main	for title.
col	color for model curves.
lty	lty for model curves.
lwd	lwd for model curves.
ex.pch	pch for extrapolation.
pch	pch for bp points.
cex	cex for bp points.
pt.col	col for bp points.
pt.lwd	lwd for bp points.
legend.x	passed to sblegend as the first argument.
...	further arguments passed to or from other methods.
z	output from previous plot.scaleboot.
y	numeric passed to legend.
inset	inset distance from the margins, which is passed to legend.
cex.legend	cex for legend
beta	matrix of beta values. beta[,1] is beta0, beta[,2] is beta1.
p	significance level for drawing contour lines.
col.contour	colors for SI, AU, BP.
drawcontours	draw contours when TRUE.
drawlabels	draw labels at contours when TRUE.
labcex	cex for contours.
length	grid size for drawing contours.
lim.countourexpend	expand contour plotting region

Details

The plot method plots bootstrap probabilities and calls the lines method, which draws fitted curves for models.

Author(s)

Hidetoshi Shimodaira

See Also

[sbfite](#), [sbphylo](#).

Examples

```
data(mam15)
## a single plot
a <- mam15.relltest[["t4"]] # an object of class "scaleboot"
plot(a,legend="topleft") # x=sigma^2, y=psi
plot(a,xval="inverse",yval="zvalue",
     legend="topleft") # x=1/sigma, y=z-value
plot(a,xval="sigma",log="x",yval="pvalue",
     legend="topleft") # x=log(sigma), y=probability
## plot of extrapolation
plot(summary(a),legend="topleft")
## multiple plots
b <- mam15.relltest[1:15] # an object of class "scalebootv"
plot(b) # x=sigma^2, y=psi
```

relltest

RELL Test for Phylogenetic Inference

Description

Performs the RELL test for finding the largest item. This calculates AU p-values for each item via the multiscale bootstrap resampling. This is particularly useful for testing tree topologies in phylogenetic analysis.

Usage

```
relltest(dat,nb=10000,sa=9^seq(-1,1,length=13),ass=NULL,
        cluster=NULL,nofit=FALSE,models=NULL,seed=100)
```

Arguments

dat	a matrix. Row vectors are to be resampled. Each column vector gives score values to be evaluated for an item. For the phylogenetic analysis, <code>dat[i, j]</code> is the site-wise log-likelihood value at site- <i>i</i> for tree- <i>j</i> , and we are to find the tree with the largest expected value of <code>sum(dat[, j])</code> .
nb	Number of replicates for each scale.
sa	Scales in sigma squared (σ^2).
ass	A list of association vectors for testing edges as well as trees. If <code>ass=NULL</code> , then only the results for trees are returned.
cluster	parallel cluster object which may be generated by function <code>makeCluster</code> .
nofit	logical. Passed to <code>sbfit</code> .
models	character vectors. Passed to <code>sbfit</code> .
seed	If non <code>NULL</code> , then a random seed is set. Specifying a seed is particularly important when <code>cluster</code> is non <code>NULL</code> , in this case <code>seed + seq(along=cluster)</code> are set to cluster nodes.

Details

`relltest` performs the resampling of estimated log-likelihoods (RELL) method of Kishino et al. (1990). For resampling indices stored in a vector `i`, the resampled log-likelihood for a tree-*j* is approximately calculated by `sum(dat[i, j])`. This approximation avoids time-consuming recalculation of the maximum likelihood estimates of tree parameters, which are to be calculated by an external phylogenetic software such as PAML as described in [mam15](#). In the implementation of `relltest`, the resampled log-likelihood is calculated by `sum(dat[i, j])*nrow(dat)/length(i)` so that the statistic is comparable to the case when $n' = n$.

`relltest` first calls `scaleboot` internally for multiscale bootstrap resampling, and then `scaleboot` calls `sbfit` for fitting models to the bootstrap probabilities. The AU p-values (named "k.3") produced by the summary method are improvements of the third-order p-values calculated by CONSEL software (Shimodaira and Hasegawa 2001). In addition, `relltest` calls `scaleboot` with `sa=1` for calculating p-values via the Shimodaira-Hasegawa test (SH-test) of Shimodaira and Hasegawa (1999).

See [mam15](#) for details through an example.

Value

`relltest` returns an object of class "`relltest`" that is inherited from the class "`scalebootv`" by adding two extra components called "`stat`" and "`shtest`". "`stat`" is a vector of the test statistics from the SH-test (i.e., the log-likelihood differences), and "`shtest`" is a list with two components: "`pv`", a vector of SH-test p-values, and "`pe`", a vector of standard errors of the p-values. The results of multiscale bootstrap resampling are stored in the "`scalebootv`" components returned by a call to `sbfit`.

Author(s)

Hidetoshi Shimodaira

References

- Kishino, H., Miyata, T. and Hasegawa, M. (1990). Maximum likelihood inference of protein phylogeny and the origin of chloroplasts., *J. Mol. Evol.*, 30, 151-160.
- Shimodaira, H. and Hasegawa, M. (1999). Multiple comparisons of log-likelihoods with applications to phylogenetic inference, *Molecular Biology and Evolution*, 16, 1114-1116.
- Shimodaira, H. and Hasegawa, M. (2001). CONSEL: for assessing the confidence of phylogenetic tree selection, *Bioinformatics*, 17, 1246-1247 (software is available from <http://stat.sys.i.kyoto-u.ac.jp/prog/consel/>).
- Luke Tierney, A. J. Rossini, Na Li and H. Sevcikova. snow: Simple Network of Workstations. R package version 0.2-1.

See Also

[sbfite](#), [scaleboot](#), [mam15](#).

Examples

```
## Not run:
## a quick example
data(mam15) # loading mam15.mt
mam15.trees <- relltest(mam15.mt,nb=10000) # nb=10000 is default
mam15.trees # SH-test p-values and result of fitting
summary(mam15.trees) # AU p-values

## End(Not run)

## Not run:
## An example from data(mam15).
## It may take 20 mins to run relltest below.
mam15.mt <- read.mt("mam15.mt") # site-wise log-likelihoods
mam15.trees <- relltest(mam15.mt) # resampling and fitting
summary(mam15.trees) # AU p-values

## End(Not run)
```

Description

Extract or modify the AIC values for models.

Usage

```
sbaic(x,...)
## S3 method for class 'scaleboot'
sbaic(x,k,...)
## S3 method for class 'scalebootv'
sbaic(x,...)

sbaic(x) <- value
## S3 replacement method for class 'scaleboot'
sbaic(x) <- value
## S3 replacement method for class 'scalebootv'
sbaic(x) <- value
```

Arguments

x	an object used to select a method.
k	numeric, the penalty per parameter to be used.
value	numeric vector of AIC values for models.
...	further arguments passed to and from other methods.

Details

sbaic can be used to modify the aic components for models in x as shown in the examples below.

Value

For an object of class "scaleboot", sbaic returns a numeric vector of AIC values for models. If k is missing, then the aic components in the fi vector of x are returned. If k is specified, $rss - k * df$ is calculated for each model. For the usual AIC, $k=2$. For the BIC (Schwarz's Bayesian information criterion), $k = \log(\sum(x\$nb))$.

Author(s)

Hidetoshi Shimodaira

References

Sakamoto, Y., Ishiguro, M., and Kitagawa G. (1986). *Akaike Information Criterion Statistics*. D. Reidel Publishing Company.

See Also

[sbfitt](#).

Examples

```
data(mam15)
a <- mam15.relltest[["t4"]] # an object of class "scaleboot"
sbaic(a) # print AIC for models
sbaic(a,k=log(sum(a$nb))) # print BIC for models
sbaic(a) <- sbaic(a,k=log(sum(a$nb))) # set BIC
sbaic(a) # print BIC for models
```

sbconf

Bootstrap Confidence Intervals

Description

A confidence interval for a scalar parameter is obtained by inverting the approximately unbiased p-value. This function is very slow, and it is currently experimental.

Usage

```
sbconf(x, ...)
```

Default S3 method:

```
sbconf(x, sa, probs=c(0.05,0.95), model="poly.2",
       k=2,s=1,sp=-1, cluster=NULL,...)
```

S3 method for class 'sbconf'

```
sbconf(x, probs=x$probs,model=x$model,
       k=x$k,s=x$s,sp=x$sp, nofit=FALSE, ...)
```

S3 method for class 'sbconf'

```
plot(x,model=x$model,k=x$k,s=x$s,sp=x$sp,
     models = attr(x$fits,"models"), log.xy = "",
     xlab="test statistic",ylab=NULL, type.plot = c("p","l","b"),
     yval=c("aic","zvalue","pvalue"), sd=2,add=FALSE, col=1:6,
     pch=NULL,lty=1:5,lwd=par("lwd"), mk.col=col[1],
     mk.lwd=lwd[1], mk.lty=lty[1], ...)
```

Arguments

x	an object used to select a method. For <code>sbconf.default</code> , x is a list vector of size <code>length{sa}</code> with each element being a vector of bootstrap replicates of a statistic or a list vector of a scalar component.
...	further arguments passed to or from other methods.
sa	vector of scales in sigma squared (σ^2).
probs	a vector of probabilities at which p-values are inverted.

model	a character to specify a model for an AU p-value. This should be included in <code>sboptions("models")</code> , for which model fitting is made internally.
k	a numeric to specify an order of AU p-value.
s	σ_0^2
sp	σ_p^2
cluster	parallel cluster object which may be generated by function <code>makeCluster</code> .
nofit	logical. No further calls to <code>sbfit</code> are made.
models	AIC values are plotted for these models.
log.xy	character string to be passed to <code>plot.default</code> .
xlab	a label for the x axis.
ylab	a label for the y axis.
type.plot	a character to be passed to <code>plot.default</code> .
yval	determines y-axis. "aic" for AIC values of models, "zvalue" for AU corrected z-values, and "pvalue" for AU corrected p-values.
sd	If positive, draws curves \pm sd*standard error for z-values and p-values.
add	logical. Should not the frame be drawn?
col	vector of colors of plots.
pch	vector of pch's of plots.
lty	vector of lty's of plots.
lwd	numeric of lwd of plots.
mk.col	color for crosses drawn at probs.
mk.lwd	lwd for crosses drawn at probs.
mk.lty	lty for crosses drawn at probs.

Details

Let $x[[i]]$ be a vector of bootstrap replicates for a statistic with scale $sa[i]$. For a threshold value y , the bootstrap probability is $bp[i]=\text{sum}(x[[i]]<y)/\text{length}(x[[i]])$. `sbconf` computes bp for several y values, and finds a value y at which the AU p-value, given by `sbfit`, equals a probability value specified in `probs`. In this manner, AU p-values are inverted to obtain bootstrap confidence intervals.

See the examples below for details.

Value

`sbconf` method returns an object of class "sbconf".

The `print` method for an object of class "sbconf" prints the confidence intervals.

Author(s)

Hidetoshi Shimodaira

See Also

[scaleboot](#).

Examples

```
## An example to calculate confidence intervals
## The test statistic is that for "t4" in data(mam15)

data(mam15) # load mam15.mt
sa <- 10^seq(-2,2,length=13) # parameter for multiscale bootstrap

## Definition of a test statistic of interest.
## "myfun" returns the maximum difference of log-likelihood value
## for a tree named a.
myfun <- function(x,w,a) maxdif(wsumrow(x,w))[[a]]
maxdif <- function(x) {
  i1 <- which.max(x) # the largest element
  x <- -x + x[i1]
  x[i1] <- -min(x[-i1]) # the second largest value
  x
}
wsumrow <- function(x,w) {
  apply(w*x,2,sum)*nrow(x)/sum(w)
}

## Not run:
## a quick example with nb=1000 (fairely fast in 2017)
## Compute multiscale bootstrap replicates
nb <- 1000 # nb = 10000 is better but slower
# the following line takes some time (less than 1 minute in 2017)
sim <- scaleboot(mam15.mt,nb,sa,myfun,"t4",count=FALSE,onlyboot=TRUE)

## show 90
## each tail is also interpreted as 95
(conf1 <- sbconf(sim$stat,sim$sa,model="sing.3",k=1)) # with k=1
(conf2 <- sbconf(conf1,model="sing.3",k=2)) # with k=2
(conf3 <- sbconf(conf2,model="sing.3",k=3)) # with k=3

## plot diagnostics for computing the confidence limits
plot(conf3) # AIC values for models v.s. test statistic value
plot(conf3,yval="zval",type="l") # corrected "k.3" z-value

## End(Not run)

## Not run:
## a longer example with nb=10000 (it was slow in 2010)
## In the following, we used 40 cpu's.
nb <- 10000

library(parallel)
cl <- makeCluster(40)
```

```

clusterExport(cl,c("maxdif","wsumrow"))

## Compute multiscale bootstrap replicates
## (It took 80 secs using 40 cpu's)
sim <- scaleboot(mam15.mt,nb,sa,myfun,"t4",count=FALSE,
                 cluster=cl,onlyboot=TRUE)

## Modify option "probs0" to a fine grid with 400 points
## default: 0.001 0.010 0.100 0.900 0.990 0.999
## NOTE: This modification is useful only when cl != NULL,
##       in which case calls to sbfit for the grid points
##       are made in parallel, although iterations seen later
##       are made sequentially.
sboptions("probs0",pnorm(seq(qnorm(0.001),qnorm(0.999),length=400)))

## Calculate bootstrap confidence intervals using "k.1" p-value.
## (It took 70 secs using 40 cpu's)
## First, sbfit is applied to bp's determined by option "probs0"
## Then, additional fitting is made only twice for iteration.
## p[1]=0.05 iter=1 t=4.342723 e=0.0003473446 r=0.0301812
## p[2]=0.95 iter=1 t=42.76558 e=0.002572495 r=0.1896809
conf1 <- sbconf(sim$stat,sim$sa,model="sing.3",k=1,cluster=cl)

## The confidence interval with "k.1" is printed as
##      0.05      0.95
## 4.342723 42.765582
conf1

## Calculate bootstrap confidence intervals
##           using "k.2" and "k.3" p-values.
## (It took only 10 secs)
## p[1]=0.05 iter=1 t=-2.974480 e=0.003729190 r=0.04725755
## p[2]=0.95 iter=1 t=39.51767 e=0.001030929 r=0.06141937
##      0.05      0.95
## -2.974480 39.517671
conf2 <- sbconf(conf1,model="sing.3",k=2)
conf2
## p[1]=0.05 iter=1 t=-3.810157 e=0.01068678 r=0.08793868
## p[2]=0.95 iter=1 t=39.32669 e=0.001711107 r=0.09464663
##      0.05      0.95
## -3.810157 39.326686
conf3 <- sbconf(conf2,model="sing.3",k=3)
conf3

## plot diagnostics
plot(conf3) # AIC values for models v.s. test statistic value
plot(conf3,yval="zval",type="l") # corrected "k.3" z-value

stopCluster(cl)

## End(Not run)

```


Description

sbfit is used to fit parametric models to multiscale bootstrap probabilities by the maximum likelihood method.

Usage

```
sbfit(x, ...)

## Default S3 method:
sbfit(x,nb,sa,models=NULL,nofit=FALSE,bpm=NULL,sam=NULL,...)

## S3 method for class 'matrix'
sbfit(x,nb,sa,models=NULL,names.hp=rownames(x),
      bpm=NULL,sam=NULL,nofit=FALSE,cluster=NULL,...)

## S3 method for class 'data.frame'
sbfit(x,...)

## S3 method for class 'scaleboot'
sbfit(x,models=names(x$fi),...)

## S3 method for class 'scalebootv'
sbfit(x,models=attr(x,"models"),...)

## S3 method for class 'scaleboot'
print(x,sort.by=c("aic","none"),...)

## S3 method for class 'scalebootv'
print(x,...)
```

Arguments

x	an object used to select a method. For <code>sbfit.default</code> , x is denoted by <code>nb</code> and is a vector of bootstrap probabilities for a hypothesis. For <code>sbfit.matrix</code> , x is denoted by <code>bps</code> and is a matrix with row vectors of <code>bp</code> for several hypotheses.
nb	vector of numbers of bootstrap replicates. A short vector (or scalar) is cyclically extended to match the size of <code>bp</code> .
sa	vector of scales in sigma squared (σ^2). Should be the same size as <code>bp</code> .
models	character vector of model names. Valid model names are <code>poly.m</code> for $m \geq 1$ and <code>sing.m</code> for $m \geq 3$. The default is set by <code>sboptions()\$models</code> , whose default is <code>c("poly.1","poly.2","poly.3","sing.3","sphe.3")</code> . If <code>models</code> is an integer value, <code>sbmodelnames(m=models)</code> is used.

<code>nofit</code>	logical. If TRUE, fitting is not performed.
<code>bpm</code>	(experimental: bootstrap probabilities for 2-step bootstrap)
<code>sam</code>	(experimental: scales for 2-step bootstrap)
<code>bpms</code>	(experimental: bootstrap probabilities for 2-step bootstrap)
<code>names.hp</code>	character vector of hypotheses names.
<code>cluster</code>	parallel cluster object which may be generated by function <code>makeCluster</code> .
<code>sort.by</code>	sort key.
<code>...</code>	further arguments passed to and from other methods.

Details

`sbfit.default` fits parametric models to `bp` by maximizing the log-likelihood value of a binomial model. A set of multiscale bootstrap resampling should be performed before a call to `sbfit` for preparing `bp`, where `bp[i]` is a bootstrap probability of a hypothesis calculated with a number of bootstrap replicates `nb[i]` and a scale $\sigma^2=sa[i]$. The scale is defined as $\sigma^2 = n/n'$, where n is the sample size of data, and n' is the sample size of replicated data for bootstrap resampling.

Each model specifies a $\psi(\beta, s) = \psi(\sigma^2 | \beta)$ function with a parameter vector β . The model may describe how the bootstrap probability changes along the scale. Let `cnt[i]=bp[i]*nb[i]` be the frequency indicating how many times the hypothesis of interest is observed in bootstrap replicates at scale `sa[i]`. Then we assume that `cnt[i]` is binomially distributed with number of trials `nb[i]` and success probability $1 - pnorm(\psi(\beta, s=sa[i]) / \sqrt{sa[i]})$. Currently, `sbpsi.poly` and `sbpsi.sing` are available as ψ functions. The estimated model parameters are accessed by the `coef.scaleboot` method.

The model fitting is performed in the order specified in `models`, and the initial values for numerical optimization of the likelihood function are prepared by using previously estimated model parameters. Thus, "poly.(m-1)" should be specified before "poly.m", and "poly.(m-1)" and "sing.(m-1)" should be specified before "sing.m".

`sbfit.matrix` calls `sbfit.default` repeatedly, once for each row vector `bp` of the matrix `bps`. Parallel computing is performed when `cluster` is non NULL.

`sbfit.scaleboot` calls `sbfit.default` with `bp`, `nb`, and `sa` components in `x` object for refitting by giving another `models` argument. It discards the previous result of fitting, and recomputes the model parameters.

`sbfit.scalebootv` calls `sbfit.matrix` with the `bps`, `nb`, and `sa` components in the attributes of `x`.

Value

`sbfit.default` and `sbfit.scaleboot` return an object of class "scaleboot", and `sbfit.matrix` and `sbfit.scalebootv` return an object of class "scalebootv".

An object of class "scaleboot" is a list containing at least the following components:

<code>bp</code>	the vector of bootstrap probabilities used.
<code>nb</code>	the <code>rep(nb, length=length(bp))</code> used.
<code>sa</code>	the <code>sa</code> used.

`fi` list vector of fitted results for models used. Each list consists of components "par" (estimated parameter), "mag" (magnification factor for "par" to make the actual parameter vector $\text{beta} = \text{par} * \text{mag}$), "value" (maximum log-likelihood), "hessian" (hessian matrix), "var" (variance estimate of "par"), "mask" (logical vector indicating parameter elements which are not at boundaries), "init" (initial values used for optimization), "psi" (psi function name of the model), "df" (degrees of freedom), "rss" (equivalent to the residual sum of squares, but actually defined as $2 * (\text{lik0} - \text{lik})$ where `lik0` and `lik` are the log-likelihood function of the non-restricted model and the model of interest, respectively), "pfit" (p-value for "rss"), "aic" (aic value of the model relative to the non-restricted model).

An object of class "scalebootv" is a vector of "scaleboot" objects, and in addition, it has attributes "models", "bps", "nb", and "sa".

Author(s)

Hidetoshi Shimodaira <shimo@i.kyoto-u.ac.jp>

References

- Shimodaira, H. (2002). An approximately unbiased test of phylogenetic tree selection, *Systematic Biology*, 51, 492-508.
- Shimodaira, H. (2004). Approximately unbiased tests of regions using multistep-multiscale bootstrap resampling, *Annals of Statistics*, 32, 2616-2641.
- Shimodaira, H. (2008). Testing Regions with Nonsmooth Boundaries via Multiscale Bootstrap, *Journal of Statistical Planning and Inference*, 138, 1227-1241. (<http://dx.doi.org/10.1016/j.jspi.2007.04.001>).

See Also

[sbpsi](#), [summary.scaleboot](#), [plot.scaleboot](#), [coef.scaleboot](#), [sbaic](#).

Examples

```
## Testing a hypothesis
## Examples of fitting models to a vector of bp's
## mam15.relltest$t4 of data(mam15), but
## using a different set of scales (sigma^2 values).
## In the below, sigma^2 ranges 0.01 to 100 in sa[i]
## This very large range is only for illustration.
## Typically, the range around 0.1 to 10
## is recommended for much better model fitting.
## In other examples, we have used
## sa = 9^seq(-1,1,length=13).

cnt <- c(0,0,0,0,6,220,1464,3565,5430,6477,6754,
        6687,5961) # observed frequencies at scales
nb <- 100000 # number of replicates at each scale
bp <- cnt/nb # bootstrap probabilities (bp's)
```

```

sa <- 10^seq(-2,2,length=13) # scales (sigma squared)
## model fitting to bp's
f <- sbfit(bp,nb,sa) # model fitting ("scaleboot" object)
f # print the result of fitting
plot(f,legend="topleft") # observed bp's and fitted curves
## approximately unbiased p-values
summary(f) # calculate and print p-values
## refitting with models up to "poly.4" and "sing.4"
f <- sbfit(f,models=1:4)
f # print the result of fitting
plot(f,legend="topleft") # observed bp's and fitted curves
summary(f) # calculate and print p-values

## Not run:
## Testing multiple hypotheses (only two here)
## Examples of fitting models to vectors of bp's
## mam15.relltest[c("t1,t2")]
cnt1 <- c(85831,81087,76823,72706,67946,62685,57576,51682,
          45887,41028,35538,31232,27832) # cnt for "t1"
cnt2 <- c(2,13,100,376,975,2145,3682,5337,7219,8559,
          10069,10910,11455) # cnt for "t2"
cnts <- rbind(cnt1,cnt2)
nb <- 100000 # number of replicates at each scale
bps <- cnts/nb # row vectors are bp's
sa <- 9^seq(-1,1,length=13) # scales (sigma squared)
fv <- sbfit(bps,nb,sa) # returns a "scalebootv" object
fv # print the result of fitting
plot(fv) # multiple plots
summary(fv) # calculate and print p-values

## End(Not run)

```

sboptions

Options for Multiscale Bootstrap

Description

To set and examine global options for scaleboot package.

Usage

```
sboptions(x, value)
```

Arguments

x	character of an option name.
value	When specified, this value is set.

Details

Invoking `sboptions()` with no arguments returns a list with the current values of the options. Otherwise it returns option(s) with name(s) specified by `x`. When value is specified, it is set to the option named `x`.

Author(s)

Hidetoshi Shimodaira

Examples

```
sboptions() # show all the options
sboptions("models") # show the default model names
new.models <- sbmodelnames(m=1:2) # character vector c("poly.1","poly.2")
old.models <- sboptions("models",new.models) # set the new model names
sboptions("models") # show the default model names
sboptions("models",old.models) # set back the default value
sboptions("models") # show the default model names
```

sbphylo

Tables for phylogenetic inference

Description

Creating tables of p-values and tree/edge associations for phylogenetic inference. Trees and edges are sorted by the likelihood value.

Usage

```
sbphylo(relltest,ass,trees,edges,edge2tree,
        treename=NULL,edgename=NULL,taxaname=NULL,mt=NULL,sort=TRUE)
```

```
## S3 method for class 'sbphylo'
summary(object, k = 2,...)
```

```
## S3 method for class 'sbphylo'
print(x,...)
```

```
## S3 method for class 'summary.sbphylo'
print(x,...)
```

Arguments

<code>relltest</code>	relltest output.
<code>ass</code>	read.ass output.
<code>trees</code>	relltest output for trees.

edges	relltest output for edges.
edge2tree	read.ass output for "edge to tree" association.
treename	character vector for tree descriptions.
edgename	character vector for edge descriptions.
taxaname	character vector for taxa names.
mt	read.mt output for the site-wise log-likelihood values.
sort	sorting trees and edges by likelihood when TRUE.
object	output of sbphylo.
k	integer of k for calculating p-values.
x	sbphylo or summary.sbphylo objects.
...	further arguments passed to and from other methods.

Details

First, apply sbphylo to consel results, and summary will make tables. Output tables are suitable for publication. For the input of sbphylo, you should specify either of (relltest, ass) or (trees, edges, edge2tree).

Value

sbphylo returns a list of several information of multiscale bootstrap. It does not do actual computation, but only sort trees and edges in decreasing order of likelihood values. The compiled information is then passed to summary method, which returns a list containing character tables and its numerical values of p-values.

Author(s)

Hidetoshi Shimodaira

See Also

[relltest](#), [read.ass](#), [read.mt](#)

Examples

```
## working with CONSEL outputs
data(mam15)
mam15.trees <- mam15.relltest[attr(mam15.ass,"trees")] # 15 trees
mam15.edges <- mam15.relltest[attr(mam15.ass,"edges")] # 10 edges
mam15.edge2tree <- mam15.ass[attr(mam15.ass,"edges")] # 10 edges
mam15 <- sbphylo(trees=mam15.trees,edges=mam15.edges,
                edge2tree=mam15.edge2tree) # sort trees and edges by likelihood
mam15 # print method for sbphylo
tab <- summary(mam15) # summary method for sbphylo
tab # prints character table

## plot (beta0,beta1)
a1 <- attr(summary(mam15$trees,k=2),"table")
```

```
a2 <- attr(summary(mam15$edges,k=2),"table")
beta <- rbind(a1$value,a2$value)[,c("beta0","beta1")]
sbplotbeta(beta) # for diagnostics of p-values
```

sbpsi *Model Specification Functions*

Description

sbpsi.poly and sbpsi.sing are ψ functions to specify a polynomial model and a singular model, respectively.

Usage

```
sbpsi.poly(beta,s=1,k=1,sp=-1,lambda=NULL,aux=NULL,check=FALSE)
sbpsi.sing(beta,s=1,k=1,sp=-1,lambda=NULL,aux=NULL,check=FALSE)
sbpsi.sphe(beta,s=1,k=1,sp=-1,lambda=NULL,aux=NULL,check=FALSE)
sbpsi.generic(beta,s=1,k=1,sp=-1,lambda=NULL,aux=NULL,check=FALSE,zfun,eps=0.01)
sbmodelnames(m=1:3,one.sided=TRUE,two.sided=FALSE,rev.sided=FALSE,
  poly,sing,poa,pob,poc,pod,sia,sib,sic,sid,sphe,pom,sim)
```

Arguments

beta	numeric vector of parameters; $\beta_0=\text{beta}[1]$, $\beta_1=\text{beta}[2]$,... $\beta_{m-1}=\text{beta}[m]$, where m is the number of parameters.
s	σ_0^2 .
k	numeric to specify the order of derivatives.
sp	σ_p^2 .
lambda	a numeric of specifying the type of p-values; Bayesian (lambda=0) Frequentist (lambda=1).
aux	auxiliary parameter. Currently not used.
check	logical for boundary check.
zfun	z-value function with (s,beta) as parameters.
eps	delta for numerical computation of derivatives.
m	numeric vector to specify the numbers of parameters.
one.sided	logical to include poly and sing models.
two.sided	logical to include poa and sia models.
rev.sided	logical to include pob and sib models.

poly	maximum number of parameters in poly models.
sing	maximum number of parameters in sing models.
sphe	maximum number of parameters in sphe models.
poa	maximum number of parameters in poa models.
pob	maximum number of parameters in pob models.
poc	maximum number of parameters in poc models.
pod	maximum number of parameters in pod models.
sia	maximum number of parameters in sia models.
sib	maximum number of parameters in sib models.
sic	maximum number of parameters in sic models.
sid	maximum number of parameters in sid models.
pom	maximum number of parameters in pom models.
sim	maximum number of parameters in sim models.

Details

For $k = 1$, the sbpsi functions return their ψ function values at $\sigma^2 = \sigma_0^2$. Currently, four types of sbpsi functions are implemented. `sbpsi.poly` defines the polynomial model;

$$\psi(\sigma^2|\beta) = \sum_{j=0}^{m-1} \beta_j \sigma^{2j}$$

for $m \geq 1$. `sbpsi.sing` defines the singular model;

$$\psi(\sigma^2|\beta) = \beta_0 + \sum_{j=1}^{m-2} \frac{\beta_j \sigma^{2j}}{1 + \beta_{m-1}(\sigma - 1)}$$

for $m \geq 3$ and $0 \leq \beta_{m-1} \leq 1$. `sbpsi.sphe` defines the spherical model; currently the number of parameters must be $m=3$. `sbpsi.generic` is a generic sbpsi function for specified `zfun`.

For $k > 1$, the sbpsi functions return values extrapolated at $\sigma^2 = \sigma_p^2$ using derivatives up to order $k - 1$ evaluated at $\sigma^2 = \sigma_0^2$;

$$q_k = \sum_{j=0}^{k-1} \frac{(\sigma_p^2 - \sigma_0^2)^j}{j!} \left. \frac{d^j \psi(x|\beta)}{dx^j} \right|_{\sigma_0^2},$$

which reduces to $\psi(\sigma_0^2|\beta)$ for $k = 1$. In the [summary.scaleboot](#), the AU p-values are defined by $p_k = 1 - \Phi(q_k)$ for $k \geq 1$.

Value

`sbpsi.poly` and `sbpsi.sing` are examples of a sbpsi function; users can develop their own sbpsi functions for better model fitting by preparing `sbpsi.foo` and `sbini.foo` functions for model `foo`. If `check=FALSE`, a sbpsi function returns the ψ function value or the extrapolation value. If `check=TRUE`, a sbpsi function returns NULL when all the elements of beta are included in the their valid intervals. Otherwise, a sbpsi function returns a list with components `beta` for the parameter value being modified to be on a boundary of the interval and `mask`, a logical vector indicating which elements are not on the boundary.

`sbmodelnames` returns a character vector of model names.

Author(s)

Hidetoshi Shimodaira

See Also[sbfit](#).

sbpval

*Extract P-values***Description**

sbpval extracts p-values from "summary.scaleboot" or "summary.scalebootv" objects.

Usage

```
sbpval(x, ...)

## S3 method for class 'summary.scaleboot'
sbpval(x,select=c("average","best","all"),...)

## S3 method for class 'summary.scalebootv'
sbpval(x,...)
```

Arguments

x an object used to select a method.

select character. If "average" or "best", only the p-values of corresponding models are returned. If "all", then p-values of all the models are returned.

... further arguments passed to or from other methods.

Details

This method is used only to extract previously calculated p-values from the summary object.

Value

The sbpval method for the class "summary.scaleboot" returns a list of three components (pvalue, sd, hypothesis). pvalue is a vector of pvalues p_k for k as specified in the summary method. sd is a vector of their standard errors. hypothesis is either "null" or "alternative" for the selective inference.

The sbpval method for the class "summary.scalebootv" returns a list of the three components, where pvalue and sd are matrices and hypothesis is a vector.

Author(s)

Hidetoshi Shimodaira

See Also

[summary.scaleboot.](#)

Examples

```
data(mam15)
a <- mam15.relltest[["t4"]] # an object of class "scaleboot"
b <- summary(a) # calculate p-values
b # print the p-values
sbpval(b) # extract a vector of p-values which are averaged by Akaike weights.
sbpval(b,select="all") # extract a matrix of p-values
```

scaleboot

Multiscale Bootstrap Resampling

Description

Performs multiscale bootstrap resampling for a specified statistic.

Usage

```
scaleboot(dat,nb,sa,fun,parm=NULL,count=TRUE,weight=TRUE,
          cluster=NULL,onlyboot=FALSE,seed=NULL,...)
```

```
countw.assmax(x,w,ass)
```

```
countw.shtest(x,w,obs)
```

```
countw.shtestass(x,w,assobs)
```

Arguments

dat	data matrix or data-frame. Row vectors are to be resampled.
nb	vector of the numbers of bootstrap replicates.
sa	vector of scales in sigma squared (σ^2).
fun	function for a statistic.
parm	parameter to be passed to fun above.
count	logical. Should only the accumulative counts be returned? Otherwise, raw statistic vectors are returned.
weight	logical. In fun above, resampling is specified by a weight vector. Otherwise, resampling is specified by a vector of indices.
cluster	parallel cluster object which may be generated by function makeCluster.
onlyboot	logical. Should only bootstrap resampling be performed? Otherwise, sbfit or sbconf is called internally.

seed	If non NULL, random seed is set. Specifying a seed is particularly important when <code>cluster</code> is non NULL, in which case <code>seed + seq(along=cluster)</code> are set to cluster nodes.
...	further arguments passed to and from other methods.
x	data matrix or data-frame passed from <code>scaleboot</code> .
w	weight vector for resampling.
ass	a list of association vectors. An example of <code>parm</code> above.
obs	a vector of observed test statistics. An example of <code>parm</code> above.
assobs	a list of <code>ass</code> and <code>obs</code> above. An example of <code>parm</code> above.

Details

These functions are used internally by `relltest` for computing raw bootstrap probabilities of phylogenetic inference. Alternatively, we used `pvclust` to get raw bootstrap probabilities of hierarchical clustering. In other cases, users may utilize `scaleboot` function or prepare their own functions.

`scaleboot` performs multiscale bootstrap resampling for a statistic defined by `fun`, which should be one of the two possible forms `fun(x, w, parm)` and `fun(x, i, parm)`. The former is used when `weight=TRUE`, and the weight vector `w` is generated by a multinomial distribution. The latter is used when `weight=FALSE`, and the index vector `i` is generated by resampling n' elements from $\{1, \dots, n\}$. When `count=TRUE`, `fun` should return a logical, or a vector of logicals.

Examples of `fun(x, w, parm)` are `countw.assmax` for AU p-values, `countw.shtest` for SH-test of trees, and `countw.shtestass` for SH-test of both trees and edges. The definitions are given below.

```
countw.assmax <- function(x,w,ass) {
  y <- maxdif(wsumrow(x,w)) <= 0 # countw.max
  if(is.null(ass)) y
  else {
    z <- vector("logical",length(ass))
    for(i in seq(along=ass)) z[i] <- any(y[ass[[i]]])
    z
  }
}

countw.shtest <- function(x,w,obs) maxdif(wsumrow(x,w)) >= obs

countw.shtestass <- function(x,w,assobs)
  unlist(assmaxdif(wsumrow(x,w),assobs$ass)) >= assobs$obs

### weighted sum of row vectors
##
## x = matrix (array of row vectors)
## w = weight vector (for rows)
##
wsumrow <- function(x,w) {
  apply(w*x,2,sum)*nrow(x)/sum(w)
}
```

```

### calc max difff
##
## y[i] := max_{j neq i} x[j] - x[i]
##
maxdif <- function(x) {
  i1 <- which.max(x) # the largest element
  x <- -x + x[i1]
  x[i1] <- -min(x[-i1]) # the second largest value
  x
}

### calc assmaxdif
##
## y[[i]][j] := max_{k neq ass[[i]]} x[k] - x[ass[[i]][j]]
##
assmaxdif <- function(x,a) {
  y <- vector("list",length(a))
  names(y) <- names(a)
  for(i in seq(along=a)) y[[i]] <- max(x[-a[[i]])] - x[a[[i]]]
  y
}

```

When `count=TRUE`, the summation of outputs from `fun` is calculated. This gives the frequencies for how many times the hypotheses are supported by the bootstrap replicates.

Value

If `onlyboot=TRUE`, then a list of raw results from the multiscale bootstrap resampling is returned. The components are "stat" for list vectors of outputs from `fun` (only when `count=FALSE`), "bps" for a matrix of multiscale bootstrap probabilities (only when `count=FALSE`), "nb" for the number of bootstrap replicates used, and "sa" for the scales used. Note that scales are redefined by `sa <- nsize/round(nsize/sa)`, where `nsize` is the sample size.

If `onlyboot=FALSE`, then the result of a call to `sbfit` is returned when `count=TRUE`, otherwise the result of `sbconf` is returned when `count=FALSE`.

Author(s)

Hidetoshi Shimodaira

See Also

[sbfit](#), [sbconf](#), [relltest](#).

Examples

```

## Not run:
## An example to calculate AU p-values for phylogenetic trees
## See also the Examples of "sbconf"

```

```

data(mam15) # load mam15.mt
sa <- 9^seq(-1,1,length=13) # parameter for multiscale bootstrap
nb <- 1000 # nb=10000 is better but slower
# Now compute bootstrap probabilities and fit models to them
sim <- scaleboot(mam15.mt,nb,sa,countw.assmax) # takes some time (< 1 min)
sim # show bootstrap probabilities and model fitting
summary(sim) # show AU p-values

## End(Not run)

## Not run:
## The following lines are only for illustration purpose
## a line from the definition of relltest
scaleboot(dat,nb,sa,countw.assmax,ass,cluster=cluster,
          names.hp=na,nofit=noft,models=models,seed=seed)

## two lines from rell.shtest (internal function)
scaleboot(z,nb,1,countw.shtest,tobs,cluster=cluster,
          onlyboot=TRUE,seed=seed)
scaleboot(z,nb,1,countw.shtestass,pa,cluster=cluster,
          onlyboot=TRUE,seed=seed)

## End(Not run)

```

summary.scaleboot *P-value Calculation for Multiscale Bootstrap*

Description

summary method for class "scaleboot" and "scalebootv".

Usage

```

## S3 method for class 'scaleboot'
summary(object,models=names(object$fi),k=3,sk=k,s=1,sp=-1,
        hypothesis=c("auto","null","alternative"),
        type=c("Frequentist","Bayesian"),...)

## S3 method for class 'scalebootv'
summary(object,models=attr(object,"models"),k=3,sk=k,
        hypothesis="auto",type="Frequentist", select="average",...)

## S3 method for class 'summary.scaleboot'
print(x,sort.by=c("aic","none"),verbose=FALSE,...)

## S3 method for class 'summary.scalebootv'
print(x,...)

```

Arguments

object	an object used to select a method.
models	character vector of model names. If numeric, names(object\$fi)[models] is used for each "scaleboot" object.
k	numeric vector of k for calculating p-values.
sk	numeric vector of k for calculating selective inference p-values.
s	σ_0^2
sp	σ_p^2
hypothesis	specifies type of selective inference. "null" takes the region as null hypothesis, and "alternative" takes the region as alternative hypothesis. "auto" determines it by the sign of beta0. The selective pvalues (sk. 1, sk. 2, ...) are selective pvalues when "null", and they are one minus selective pvalues when "alternative".
type	If numeric, it is passed to sbpsi functions as lambda to specify p-value type. If "Frequentist" or "Bayesian", then equivalent to specifying lambda = 1 or 0, respectively.
select	character of model name (such as "poly.3") or one of "average" and "best". If "average" or "best", then the averaging by Akaike weights or the best model is used, respectively.
x	object.
sort.by	sort key.
verbose	logical.
...	further arguments passed to and from other methods.

Details

For each model, a class of approximately unbiased p-values, indexed by $k = 1, 2, \dots$, is calculated. The p-values are named k.1, k.2, ..., where $k = 1$ (k.1) corresponds to the ordinary bootstrap probability, and $k = 2$ (k.2) corresponds to the third-order accurate p-value of Shimodaira (2002). As the k value increases, the bias of testing decreases, although the p-value becomes less stable numerically and the monotonicity of rejection regions becomes worse. Typically, $k = 3$ provides a reasonable compromise. The `sbpval` method is available to extract p-values from the "summary.scaleboot" object.

The p-value is defined as

$$p_k = 1 - \Phi \left(\sum_{j=0}^{k-1} \frac{(\sigma_p^2 - \sigma_0^2)^j}{j!} \frac{d^j \psi(x|\beta)}{dx^j} \Big|_{\sigma_0^2} \right),$$

where $\psi(\sigma^2|\beta)$ is the model specification function, σ_0^2 is the evaluation point for the Taylor series, and σ_p^2 is an additional parameter. Typically, we do not change the default values $\sigma_0^2 = 1$ and $\sigma_p^2 = -1$.

The p-values are justified only for good fitting models. By default, the model which minimizes the AIC value is selected. We can modify the AIC value by using the `sbaic` function. We also diagnose the fitting by using the `plot` method.

Now includes selective inference p-values. The method is described in Terada and Shimodaira (2017; arXiv:1711.00949) "Selective inference for the problem of regions via multiscale bootstrap".

Value

summary.scaleboot returns an object of the class "summary.scaleboot", which is inherited from the class "scaleboot". It is a list containing all the components of class "scaleboot" and the following components:

pv	matrix of p-values of size length(models) * length(k) with elements p_k .
pe	matrix of standard errors of p-values.
spv	matrix of selective inference p-values of size length(models) * length(sk) with elements sp_k .
spe	matrix of standard errors of selective inference p-values.
betapar	list array containing (beta0, beta1) and its covariance matrix for each model. They are obtained by linear extrapolation. This will be used for interpreting the fitting in terms of signed distance and curvature.
best	a list consisting of components model for the best fitting model name, aic for its AIC value, pv and spv for vector of p-values, and pe and spe for vectors of standard errors. Also includes betapar for the best model.
average	a list of results for the average model computed by Akaike weight.
parex	a list of components k, s, and sp.

Author(s)

Hidetoshi Shimodaira

See Also

[sbfit](#), [sbpsi](#), [sbpval](#), [sbaic](#).

Examples

```
data(mam15)
## For a single hypothesis
a <- mam15.relltest[["t4"]] # an object of class "scaleboot"
summary(a) # calculate and print p-values (k=3)
summary(a,k=2) # calculate and print p-values (k=2)
summary(a,k=1:4) # up to "k.4" p-value.

## For multiple hypotheses
b <- mam15.relltest[1:15] # an object of class "scalebootv"
summary(b) # calculate and print p-values (k=3)
summary(b,k=1:4) # up to "k.4" p-value.
```

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