Package ‘paleotree’

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Description Provides tools for transforming, a posteriori time-scaling, and modifying phylogenies containing extinct (i.e. fossil) lineages. In particular, most users are interested in the functions timePaleoPhy(), bin_timePaleoPhy(), cal3TimePaleoPhy() and bin_cal3TimePaleoPhy(), which date cladograms of fossil taxa using stratigraphic data. This package also contains a large number of likelihood functions for estimating sampling and diversification rates from different types of data available from the fossil record (e.g. range data, occurrence data, etc). paleotree users can also simulate diversification and sampling in the fossil record using the function simFossilRecord(), which is a detailed simulator for branching birth-death-sampling processes composed of discrete taxonomic units arranged in ancestor-descendant relationships. Users can use simFossilRecord() to simulate diversification in incompletely sampled fossil records, under various models of morphological differentiation (i.e. the various patterns by which morphotaxa originate from one another), and with time-dependent, longevity-dependent and/or diversity-dependent rates of diversification, extinction and sampling. Additional functions allow users to translate simulated ancestor-descendant data from simFossilRecord() into standard time-scaled phylogenies or unscaled cladograms that reflect the relationships among taxon units.

License CC0
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Description

Analyzes, time-scales and simulates phylogenies of extinct/fossil lineages, along with calculation of diversity curves. Also fits likelihood models to estimate sampling rates from stratigraphic ranges.
Details

Package: paleotree
Type: Package
License: CC0

This package contains functions for analyzing sampling rates given ranges of fossil taxa, in both continuous and discrete time, functions for a posteriori time-scaling phylogenies of fossil taxa and functions for simulating the fossil record in both taxic and phylogenetic varieties.

Author(s)

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References


See Also

This package relies extensively on the phylogenetic toolkit and standards offered by the ape package, and hence lists this package as a depends, so it is loaded simultaneously.

Examples

```r
# get the package version of paleotree
closestVersion("paleotree")

# get the citation for paleotree
citation("paleotree")

# Simulate some fossil ranges with simFossilRecord
set.seed(444);
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
ntaxa = c(30,40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
# let's see what the 'true' diversity curve looks like in this case
plot the FADs and LADs with taxicDivCont()
taxicDivCont(taxa)
```
"#simulate a fossil record with imperfect sampling with sampleRanges
rangesCont <- sampleRanges(taxa, r = 0.5)
#plot the diversity curve based on the sampled ranges
layout(1:2)
taxicDivCont(rangesCont)
#Now let's use binTimeData to bin in intervals of 10 time units
rangesDisc <- binTimeData(rangesCont, int.length = 10)
#plot with taxicDivDisc
taxicDivDisc(rangesDisc)
#compare to the continuous time diversity curve
layout(1)

taxa2phylo assumes we know speciation events perfectly... what if we don't?
#first, let's use taxa2cladogram to get the 'ideal' cladogram of the taxa
cladogram <- taxa2cladogram(taxa, plot = TRUE)
#Now let's try timePaleoPhy using the continuous range data
tree <- timePaleoPhy(cladogram, rangesCont, type = "basic", plot = TRUE)
#plot diversity curve
phyloDiv(tree, drop.ZLB = TRUE)

#that tree lacked the terminal parts of ranges (tips stops at the taxon FADs)
#let's add those terminal ranges back on with add.term
tree <- timePaleoPhy(cladogram, rangesCont, type = "basic", add.term = TRUE, plot = TRUE)
#plot diversity curve
phyloDiv(tree)

---

**binTimeData**

*Bin Simulated Temporal Ranges in Discrete Intervals*

**Description**

Converts a matrix of simulated continuous-time first occurrences and last occurrences for fossil taxa into first and last occurrences given in some set of simulated/placed discrete-time intervals, which is output along with information of the dates of the given intervals.

**Usage**

```
binTimeData(timeData, int.length = 1, start = NA, int.times = NULL)
```

**Arguments**

- **timeData**: Two-column matrix of simulated first and last occurrences in absolute continuous time
- **int.length**: Time interval length, default is 1 time unit
- **start**: Starting time for calculating the intervals.
- **int.times**: A two column matrix with the start and end times of the intervals to be used.
Details

This function takes a simulated matrix of per-taxon first and last occurrences and, by dividing the time-scale into time intervals of non-zero length, lists taxon occurrences within those intervals. By default, a set of sequential non-overlapping time-interval of equal non-zero length are used, with the length controlled by the argument int.length.

Alternatively, a two column matrix of interval start and end times to be used can be input via the argument int.times. None of these intervals can have a duration (temporal length) greater than zero. If a first or last appearance in the input range data could fit into multiple intervals (i.e. the input discrete time intervals are overlapping), then the appearance data is placed in the interval of the shortest duration. When output, the interval times matrix (see below) will be sorted from first to last.

This function is SPECIFICALLY for simulating the effect of having a discrete time-scale for analyses using simulations. This function should not be used for non-simulations uses, such as binning temporal occurrences for analyses of real data. In those case, the temporal ranges (which, in real data, will probably be given as discrete time intervals) should already be tabulated within discrete intervals prior to use in paleotree. The user should place the temporal information in a list object, as described for the output of binTimeData (see below).

As with many functions in the paleotree library, absolute time is always decreasing, i.e. the present day is zero. However, the numbering of intervals giving in the output increases with time, as these are numbered relative to each other, from first to last.

As of version 1.7 of paleotree, taxa which are extant as indicated in timeData as being in a time interval bounded (0,0), unless time-bins are preset using argument int.times (prior to version 1.5 they were erroneously listed as NA).

Value

A list containing:

- **int.times** A 2 column matrix with the start and end times of the intervals used; time decreases relative to the present.
- **taxon.times** A 2 column matrix with the first and last occurrences of taxa in the intervals listed in int.times, with numbers referring to the row of int.times.

Author(s)

David W. Bapst

See Also

- `simFossilRecord`
- `sampleRanges`
- `taxicDivCont`

Examples

```r
#Simulate some fossil ranges with simFossilRecord
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(30,40), nExtant = 0)
```
branchClasses

```r
taxa <- fossilRecord2FossilTaxa(record)
# simulate a fossil record with imperfect sampling with sampleRanges()
rangesCont <- sampleRanges(taxa, r = 0.5)
# Now let's use binTimeData() to bin in intervals of 1 time unit
rangesDisc <- binTimeData(rangesCont, int.length = 1)
# plot with taxicDivDisc()
equalDiscInt <- taxicDivDisc(rangesDisc)

# example with pre-set intervals input (including overlapping)
presetIntervals <- cbind(c(1000,990,990,940), c(980,970,950,930))
rangesDisc1 <- binTimeData(rangesCont, int.times = presetIntervals)
taxicDivDisc(rangesDisc1)
# now let's plot diversity with (different) equal length intervals used above
taxicDivDisc(rangesDisc1, int.times = equalDiscInt[,1:2])

# example with extant taxa
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(30,40))
taxa <- fossilRecord2FossilTaxa(record)
# simulate a fossil record with imperfect sampling with sampleRanges()
rangesCont <- sampleRanges(taxa, r = 0.5, modern.samp.prob = 1)
# Now let's use binTimeData() to bin in intervals of 1 time unit
rangesDisc <- binTimeData(rangesCont, int.length = 1)
# plot with taxicDivDisc()
taxicDivDisc(rangesDisc)

# example with pre-set intervals input (including overlapping)
presetIntervals <- cbind(c(40,30,20,10), c(30,20,10,0))
rangesDisc1 <- binTimeData(rangesCont, int.times = presetIntervals)
taxicDivDisc(rangesDisc1)
```

---

**branchClasses**

Partitions the branch lengths of a tree into several classes based on their placement.

**Description**

Partitions the branch lengths of a tree into several classes based on their placement.

**Usage**

```r
branchClasses(tree, whichExtant = NULL, tol = 0.01)
```

**Arguments**

- `tree` A time-scaled phylogeny to be analyzed, as an object of class phylo, ideally with a $root.time$ element as is typical for paleotree output phylogenies. If $root.time$ is not present, the most recent tips will be interpreted as being at the modern day (i.e. 0 time-units before present).
whichExtant

A logical vector with length equal to number of tips in the tree. TRUE indicates a taxon that is extant at the modern day, while FALSE equates to extinction. If present, this vector is used for determining which taxa are extant, and which are extinct.

tol

Tolerance used to distinguish extant taxa, if whichExtant is not provided, to avoid issues with number rounding. Taxa within tol of the modern day will be considered extant.

Details

This function will partition the internode (node to node, including internal node to terminal tip) branch lengths of a tree into four separate classes: all (all the internode branches of a tree), int (internal branches which run from one internode to another), live (terminal branches which run from an internal node to a terminal tip representing an extinction event before the present) and dead (terminal branches which run from an internal node to a terminal tip at the modern day, reflecting a still-living taxon).

The depths of the internal 'mother' node (i.e. time of origin, before the modern day of each branch length are included as the /codenames of the branch length vectors.

This function is mainly of use for modeling internode branch lengths in a phylogeny including fossil taxa.

Value

The output is a list consisting of four vectors, with the /codenames of the vectors being their corresponding time of origin. See details.

Examples

# simulated example
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1, nTotalTaxa = c(30, 40), nExtant = c(10, 20))
taxa <- fossilRecord2FossilTaxa(record)
tree <- taxa2phylo(taxa)
brlenRes <- branchClasses(tree)

# see frequency histograms of branch lengths
layout(1:4)
for(x in 1:length(brlenRes)){
hist(brlenRes[[x]],main = "Branch Lengths",xlab = names(brlenRes)[x])
}

# see frequency histograms of branch depths
layout(1:4)
for(x in 1:length(brlenRes)){
hist(as.numeric(names(brlenRes[[x]])),main = "Branch Depths",xlab = names(brlenRes)[x])
}

layout(1)
Three Rate Calibrated 'a posteriori' Time-Scaling of Paleo-Phylogenies

Description

Time-scales an unscaled cladogram of fossil taxa, using information on their ranges and estimates of the instantaneous rates of branching, extinction and sampling. The output is a sample of a posteriori time-scaled trees, as resulting from a stochastic algorithm which samples observed gaps in the fossil record with weights calculated based on the input rate estimates. This function also uses the three-rate calibrated time-scaling algorithm to stochastically resolve polytomies and infer potential ancestor-descendant relationships, simultaneous with the time-scaling treatment.

Usage

cal3TimePaleoPhy(tree, timeData, brRate, extRate, sampRate, ntrees = 1, 
anctwt = 1, node.mins = NULL, dateTreatment = "firstLast", 
FAD.only = FALSE, adj.obs.wt = TRUE, root.max = 200, step.size = 0.1,  
randres = FALSE, noisyDrop = TRUE, verboseWarnings = TRUE,  
diagnosticMode = FALSE, tolerance = 1e-04, plot = FALSE)

bin_cal3TimePaleoPhy(tree, timeList, brRate, extRate, sampRate, ntrees = 1, 
anctwt = 1, node.mins = NULL, dateTreatment = "firstLast",  
FAD.only = FALSE, sites = NULL, point.occurs = FALSE,  
nonstoch.bin = FALSE, adj.obs.wt = TRUE, root.max = 200,  
step.size = 0.1, randres = FALSE, noisyDrop = TRUE,  
verboseWarnings = TRUE, tolerance = 1e-04, diagnosticMode = FALSE,  
plot = FALSE)

Arguments

tree An unscaled cladogram of fossil taxa, of class phylo. Tip labels must match the taxon labels in the respective temporal data.

timeData Two-column matrix of first and last occurrences in absolute continuous time, with row names as the taxon IDs used on the tree. This means the first column is very precise FADs (first appearance dates) and the second column is very precise LADs (last appearance dates), reflect the precise points in time when taxa first and last appear. If there is stratigraphic uncertainty in when taxa appear in the fossil record, it is preferable to use the 'bin' time-scaling functions; however, see the argument dateTreatment.

brRate Either a single estimate of the instantaneous rate of branching (also known as the 'per-capita' origination rate, or speciation rate if taxonomic level of interest is species) or a vector of per-taxon estimates

extRate Either a single estimate of the instantaneous extinction rate (also known as the 'per-capita' extinction rate) or a vector of per-taxon estimates
sampRate
Either a single estimate of the instantaneous sampling rate or a vector of per-
taxon estimates.

ntrees
Number of time-scaled trees to output.

anc.wt
Weighting against inferring ancestor-descendant relationships. The argument
anc.wt allows users to change the default consideration of anc-desc relationships. This value is used as a multiplier applied to the probability of choosing any node position which would infer an ancestor-descendant relationship. By default, anc.wt = 1, and thus these probabilities are unaltered. If anc.wt is less than 1, the probabilities decrease and at anc.wt = 0, no ancestor-descendant relationships are inferred at all. Can be a single value or a vector of per-taxon values, such as if a user wants to only allow plesiomorphic taxa to be ancestors.

node.mins
The minimum dates of internal nodes (clades) on a phylogeny can be set using node.mins. This argument takes a vector of the same length as the number of nodes, with dates given in the same order as nodes are ordered in the tree$edge matrix. Note that in tree$edge, terminal tips are given the first set of numbers (1:ntip(tree)), so the first element of node.mins is the first internal node (the node numbered ntip(tree)+1, which is generally the root for most phylo objects read by read.tree). Not all nodes need be given minimum dates; those without minimum dates can be given as NA in node.mins, but the vector must be the same length as the number of internal nodes in tree. These are minimum date constraints, such that a node will be ‘frozen’ by the cal3 algorithm so that constrained nodes will always be at least as old as this date, but the final date may be even older depending on the taxon dates used, the parameters input for the cal3 algorithm and any other minimum node dates given (e.g. if a clade is given a very old minimum date, this will (of course) over-ride any minimum dates given for clades that that node is nested within). If the constrained nodes include a polytomy, this polytomy will still be resolved with respect to the cal3 algorithm, but the first divergence will be ‘frozen’ so that it is at least as old as the minimum age, while any additional divergences will be allowed to occur after this minimum age.

dateTreatment
This argument controls the interpretation of timeData. The default setting ‘first-
Last’ treats the dates in timeData as a column of precise first and last appearances. A second option, added by great demand, is ‘minMax’ which treats these dates as minimum and maximum bounds on single point dates. Under this option, all taxa in the analysis will be treated as being point dates, such that the first appearance is also the last. These dates will be pulled under a uniform distribution. If ‘minMax’ is used, add.term becomes meaningless, and the use of it will return an error message. A third option is ‘randObs’. This assumes that the dates in the matrix are first and last appearance times, but that the desired time of observation is unknown. Thus, this is much like ‘firstLast’ except the effective time of observation (the taxon’s LAD under ‘firstLast’) is treated an uncertain date, and is randomly sampled between the first and last appearance times. The FAD still is treated as a fixed number, used for dating the nodes. In previous versions of paleotree, this was called in cal3timePaleoPhy using the argument rand.obs, which has been removed for clarity. This temporal uncertainty in times of observation might be useful if a user is interested in applying phylogeny-based approaches to studying trait evolution, but have per-taxon
measurements of traits that come from museum specimens with uncertain temporal placement. With both arguments 'minMax' and 'randObs', the sampling of dates from random distributions should compel users to produce many time-scaled trees for any given analytical purpose. Note that 'minMax' returns an error in 'bin' time-scaling functions; please use 'points.occur' instead.

**FAD.only** Should the tips represent observation times at the start of the taxon ranges? If TRUE, result is similar to when terminal ranges are no added on with timePaleoPhy. If FAD.only is TRUE and dateTreatment is 'minMax' or 'randObs', the function will stop and a warning will be produced, as these combinations imply contradictory sets of times of observation.

**adj.obs.wt** If the time of observation are before the LAD of a taxon, should the weight of the time of observation be adjusted to account for the known observed history of the taxon which occurs AFTER the time of observation? Should only have an effect if time of observation IS NOT the LAD, if times of observation for a potential ancestor are earlier than the first appearance of their potential descendants and if the ancestral weights for taxa are not set to zero (so there can be potential ancestors).

**root.max** Maximum time before the first FAD that the root can be pushed back to.

**step.size** Step size of increments used in zipper algorithm to assign node ages.

**randres** Should polytomies be randomly resolved using multiRdi in ape rather than using the cal3 algorithm to weight the resolution of polytomies relative to sampling in the fossil record?

**noisyDrop** If TRUE (the default), any taxa dropped from tree due to not having a matching entry in the time data will be listed in a system message.

**verboseWarnings** if TRUE (the default), then various warnings and messages regarding best practices will be issued to the console about the analysis. If FALSE, the function will run as quietly as possible.

**diagnosticMode** If TRUE, cal3TimePaleoPhy will return to the console and to graphic devices an enormous number of messages, plots and ancilliary information that may be useful or entirely useless to figuring out what is going wrong.

**tolerance** Acceptable amount of shift in tip dates from dates listed in timeData. Shifts outside of the range of tolerance will cause a warning message to be issued that terminal tips appear to be improperly aligned.

**plot** If true, plots the input, "basic" time-scaled phylogeny (an intermediary step in the algorithm) and the output cal3 time-scaled phylogeny.

**timeList** A list composed of two matrices giving interval times and taxon appearance dates. The rownames of the second matrix should be the taxon IDs, identical to the tip.labels for tree. See details.

**sites** Optional two column matrix, composed of site IDs for taxon FADs and LADs. The sites argument allows users to constrain the placement of dates by restricting multiple fossil taxa whose FADs or LADs are from the same very temporarily restricted sites (such as fossil-rich Lagerstatten) to always have the same date, across many iterations of time-scaled trees. To do this, simply give a matrix
where the "site" of each FAD and LAD for every taxon is listed, as corresponding to the second matrix in timeList. If no sites matrix is given (the default), then it is assumed all fossil come from different "sites" and there is no shared temporal structure among the events.

point.occur
If true, will automatically produce a 'sites' matrix which forces all FADs and LADs to equal each other. This should be used when all taxa are only known from single 'point occurrences', i.e. each is only recovered from a single bed/horizon, such as a Lagerstatten.

nonstoch.bin
If true, dates are not stochastically pulled from uniform distributions. See below for more details.

Details

The three-rate calibrated ("cal3") algorithm time-scales trees a posteriori by stochastically picking node divergence times relative to a probability distribution of expected waiting times between speciation and first appearance in the fossil record. This algorithm is extended to apply to resolving polytomies and designating possible ancestor-descendant relationships. The full details of this method are provided in Bapst (2013, MEE).

Briefly, cal3 time-scaling is done by examining each node separately, moving from the root upwards. Ages of branching nodes are constrained below by the ages of the nodes below them (except the root; hence the need for the root.max argument) and constrained above by the first appearance dates (FADs) of the daughter lineages. The position of the branching event within this constrained range implies different amounts of unobserved evolutionary history. cal3 considers a large number of potential positions for the branching node (the notation in the code uses the analogy of viewing the branching event as a 'zipper') and calculates the summed unobserved evolutionary history implied by each branching time. The probability density of each position is then calculated under a gamma distribution with a shape parameter of 2 (implying that it is roughly the sum of two normal waiting times under an exponential) and a rate parameter which takes into account both the probability of not observing a lineage of a certain duration and the 'twiginess' of the branch, i.e. the probability of having short-lived descendants which went extinct and never were sampled (ala Friedman and Brazeau, 2011). These densities calculated under the gamma distribution are then used as weights to stochastically sample the possible positions for the branching node. This basic framework is extended to polytomies by allowing a branching event to fall across multiple potential lineages, adding each lineage one by one, from earliest appearing to latest appearing (the code notation refers to this as a 'parallel zipper').

As with many functions in the paleotree library, absolute time is always decreasing, i.e. the present day is zero.

These functions will intuitively drop taxa from the tree with NA for range or that are missing from timeData.

The sampling rate used by cal3 methods is the instantaneous sampling rate, as estimated by various other function in the paleotree package. See make_durationFreqCont for more details. If you have the per-time unit sampling probability ('R' as opposed to 'r') look at the sampling parameter conversion functions also included in this package (e.g. sProb2sRate). Most datasets will probably use make_durationFreqDisc and sProb2sRate prior to using this function, as shown in an example below.

The branching and extinction rate are the 'per-capita' instantaneous origination/extinction rates for the taxic level of the tips of the tree being time-scaled. Any user of the cal3 time-scaling
method has multiple options for estimating these rates. One is to separately calculate
the per-capita rates (following the equations in Foote, 2001) across multiple intervals and

take the mean for each rate. A second, less preferred option, would be to use the extinction rate
calculated from the sampling rate above (under ideal conditions, this should be very close
to the mean 'per-capita' rate calculated from by-interval FADs and LADs). The branching rate
in this case could be assumed to be very close to the extinction rate, given the tight relationship
observed in general between these two (Stanley, 1976; see Foote et al., 1999, for a defence of
this approach), and thus the extinction rate estimate could be used also for the branching rate estimate.
(This is what is done for the examples below.) A third option for calculating all three rates simultaneously
would be to apply likelihood methods developed by Foote (2002) to forward and reverse survivorship
curves. Note that only one of these three suggested methods is implemented in
cal3TimePaleoPhy.

By default, the cal3 functions will consider that ancestor-descendant relationships may exist among
the given taxa, under a budding cladogenetic or anagenetic modes. Which tips are designated as
which is given by two additional elements added to the output tree, $budd.tips (taxa designated as
ancestors via budding cladogenesis) and $anag.tips (taxa designated as ancestors via anagenesis).
This can be turned off by setting anc.wt = 0. As this function may infer anagenetic relationships
during time-scaling, this can create zero-length terminal branches in the output. Use dropZLB to get
rid of these before doing analyses of lineage diversification.

Unlike timePaleoPhy, cal3 methods will always resolve polytomies. In general, this is done using
the rate calibrated algorithm, although if argument randres = TRUE, polytomies will be randomly
resolved with uniform probability, ala multi2di from ape. Also, cal3 will always add the terminal
ranges of taxa. However, because of the ability to infer potential ancestor-descendant relationships,
the length of terminal branches may be shorter than taxon ranges themselves, as budding may have
occurred during the range of a morphologically static taxon. By resolving polytomies with the cal3
method, this function allows for taxa to be ancestral to more than one descendant taxon. Thus, users
who believe their dataset may contain indirect ancestors are encouraged by the package author to
try cal3 methods with their consensus trees, as opposed to using the set of most parsimonious trees.
Comparing the results of these two approaches may be very revealing.

Like timePaleoPhy, cal3TimePaleoPhy is designed for direct application to datasets where taxon
first and last appearances are precisely known in continuous time, with no stratigraphic uncertainty.
This is an uncommon form of data to have from the fossil record, although not an impossible
form (micropaleontologists often have very precise range charts, for example). This means that
most users should not use cal3TimePaleoPhy directly, unless they have written their own code
to deal with stratigraphic uncertainty. For some groups, the more typical 'first' and 'last' dates
represent the minimum and maximum absolute ages for the fossil collections that a taxon is known
is known from. Presumably, the first and last appearances of that taxon in the fossil record is at
unknown dates within these bounds. These should not be mistaken as the FADs and LADs desired
by cal3TimePaleoPhy, as cal3TimePaleoPhy will use the earliest dates provided to calibrate node
ages, which is either an overly conservative approach to time-scaling or fairly nonsensical.

Alternatively to using cal3TimePaleoPhy, bin_cal3TimePaleoPhy is a wrapper of cal3TimePaleoPhy
which produces time-scaled trees for datasets which only have interval data available. For each output
tree, taxon first and last appearance dates are placed within their listed intervals under a uniform
distribution. Thus, a large sample of time-scaled trees will approximate the uncertainty in the actual
timing of the FADs and LADs.

The input timeList object can have overlapping (i.e. non-sequential) intervals, and intervals of
uneven size. Taxa alive in the modern should be listed as last occurring in a time interval that begins at time 0 and ends at time 0. If taxa occur only in single collections (i.e. their first and last appearance in the fossil record is synchronous, the argument point.occur will force all taxa to have instantaneous durations in the fossil record. Otherwise, by default, taxa are assumed to first and last appear in the fossil record at different points in time, with some positive duration. The sites matrix can be used to force only a portion of taxa to have simultaneous first and last appearances.

By setting the argument nonstoch.bin to TRUE in bin_cal3TimePaleoPhy, the dates are NOT stochastically pulled from uniform bins but instead FADs are assigned to the earliest time of whichever interval they were placed in and LADs are placed at the most recent time in their placed interval. This option may be useful for plotting. The sites argument becomes arbitrary if nonstoch.bin is TRUE.

If timeData or the elements of timelist are actually data frames (as output by read.csv or read.table), these will be coerced to a matrix.

A tutorial for applying the time-scaling functions in paleotree, particularly the cal3 method, along with an example using real (graptolite) data, can be found at the following link:
http://nemagruptus.blogspot.com/2013/06/a-tutorial-to-cal3-time-scaling-using.html

Value

The output of these functions is a time-scaled tree or set of time-scaled trees, of either class phylo or multiphylo, depending on the argument ntrees. All trees are output with an element $\$root.time. This is the time of the root on the tree and is important for comparing patterns across trees.

Additional elements are sampledLogLike and $\$sumLogLike which respectively record a vector containing the 'log-densities' of the various node-ages selected for each tree by the 'zipper' algorithm, and the sum of those log-densities. Although they are very similar to log-likelihood values, they may not be true likelihoods, as node ages are conditional on the other ages selected by other nodes. However, these values may give an indication about the relative optimality of a set of trees output by the cal3 functions.

Trees created with bin_cal3TimePaleoPhy will output with some additional elements, in particular $\$ranges.used, a matrix which records the continuous-time ranges generated for time-scaling each tree. (Essentially a pseudo-timeData matrix.)

Note

Most importantly, please note the stochastic element of the three rate-calibrated time-scaling methods. These do not use traditional optimization methods, but instead draw divergence times from a distribution defined by the probability of intervals of unobserved evolutionary history. This means analyses MUST be done over many cal3 time-scaled trees for analytical rigor! No one tree is correct.

Similarly, please account for stratigraphic uncertainty in your analysis. Unless you have exceptionally resolved data, use a wrapper with the cal3 function, either the provided bin_cal3TimePaleoPhy or code a wrapper function of your own that accounts for stratigraphic uncertainty in your dataset. Remember that the FADs (earliest dates) given to timePaleoPhy will *always* be used to calibrate node ages!
Author(s)
David W. Bapst

References


See Also

timePaleoPhy, make_durationFreqCont, pqrPs, sProb2sRate, multi2di

Examples

```r
#Simulate some fossil ranges with simFossilRecord
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1, nTotalTaxa = c(30, 40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)

#simulate a fossil record with imperfect sampling with sampleRanges
rangesCont <- sampleRanges(taxa, r = 0.5)

#let's use taxa2cladogram to get the 'ideal' cladogram of the taxa
cladogram <- taxa2cladogram(taxa, plot = TRUE)

#this library allows one to use rate calibrated type time-scaling methods (Bapst, in prep.)
#to use these, we need an estimate of the sampling rate (we set it to 0.5 above)
likFun <- make_durationFreqCont(rangesCont)
srRes <- optim(parInit(likFun), likFun, lower = parLower(likFun), upper = parUpper(likFun), method = "L-BFGS-B", control = list(maxit = 1000000))
sRate <- srRes[[1]][2]

# we also need extinction rate and branching rate
# we can get extRate from getSampRateCont too
# we'll assume extRate = brRate (ala Foote et al., 1999); may not always be a good assumption
divRate <- srRes[[1]][1]

#now let's try cal3TimePaleoPhy, which time-scales using a sampling rate to calibrate
#this can also resolve polytomies based on sampling rates, with some stochastic decisions
ttree <- cal3TimePaleoPhy(cladogram, rangesCont, brRate = divRate, extRate = divRate, sampRate = sRate, ntrees = 1, plot = TRUE)

#notice the warning it gives!
phyloDiv(ttree)
```
# by default, cal3TimePaleoPhy may predict indirect ancestor-descendant relationships
# can turn this off by setting anc.wt = 0
tree <- cal3TimePaleoPhy(cladogram, rangesCont, brRate = divRate, extRate = divRate,
        sampRate = sRate, ntrees = 1, anc.wt = 0, plot = TRUE)

# let's look at how three trees generated with very different time of obs. look
FAD <- cal3TimePaleoPhy(cladogram, rangesCont, brRate = divRate, extRate = divRate,
        FAD.only = TRUE, dateTreatment = "firstLast", sampRate = sRate, ntrees = 1, plot = TRUE)
treeRand <- cal3TimePaleoPhy(cladogram, rangesCont, brRate = divRate, extRate = divRate,
        FAD.only = FALSE, dateTreatment = "randObs", sampRate = sRate, ntrees = 1, plot = TRUE)

# by default the time of observations are the LADs
treeLAD <- cal3TimePaleoPhy(cladogram, rangesCont, brRate = divRate, extRate = divRate,
        FAD.only = FALSE, dateTreatment = "randObs", sampRate = sRate, ntrees = 1, plot = TRUE)

layout(1:3)
parOrig <- par(no.readonly = TRUE)
par(mar = c(0, 0, 0, 0))
plot(ladderize(treeFAD)); text(5, 5, "time.obs = FAD", cex = 1.5, pos = 4)
plot(ladderize(treeRand)); text(5, 5, "time.obs = Random", cex = 1.5, pos = 4)
plot(ladderize(treeLAD)); text(5, 5, "time.obs = LAD"), cex = 1.5, pos = 4)
layout(1); par(parOrig)

# to get a fair sample of trees, let's increase ntrees
ntrees <- cal3TimePaleoPhy(cladogram, rangesCont, brRate = divRate, extRate = divRate,
        sampRate = sRate, ntrees = 9, plot = FALSE)

# let's compare nine of them at once in a plot
layout(matrix(1:9, 3, 3))
parOrig <- par(no.readonly = TRUE)
par(mar = c(0, 0, 0, 0))
for(i in 1:9)(plot(ladderize(ntrees[[i]]), show.tip.label = FALSE))
layout(1)
par(parOrig)

# they are all a bit different!

# can plot the median diversity curve with multiDiv
multiDiv(ntrees)

# using node.mins
# let's say we have (molecular??) evidence that node #5 is at least 1200 time-units ago
# to use node.mins, first need to drop any unshared taxa
droppers <- cladogram$tip.label[is.na(
        match(cladogram$tip.label, names(which(is.na(rangesCont[, 1])))))]
cladodrop <- drop.tip(cladogram, droppers)
# now make vector same length as number of nodes
nodeDates <- rep(NA, Nnode(cladodrop))
nodeDates[5] <- 1200
tree <- cal3TimePaleoPhy(cladodrop, rangesCont, brRate = divRate, extRate = divRate,
        sampRate = sRate, ntrees = 1, node.mins = nodeDates, plot = TRUE)

# example with time in discrete intervals
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
        nTotalTaxa = c(30, 40), nExtant = 0)
taxa <- fossilRecord2fossilTaxa(record)
# simulate a fossil record with imperfect sampling with sampleRanges()
rangesCont <- sampleRanges(taxa,r = 0.5)
# let's use taxa2cladogram to get the 'ideal' cladogram of the taxa
cladogram <- taxa2cladogram(taxa,plot = TRUE)
# Now let's use binTimeData to bin in intervals of 1 time unit
rangesDisc <- binTimeData(rangesCont,int.length = 1)
# we can do something very similar for the discrete time data (can be a bit slow)
likFun <- make_durationFreqDisc(rangesDisc)
spRes <- optim(parInit(likFun),likFun,lower = parLower(likFun),upper = parUpper(likFun),
method = "L-BFGS-B",control = list(maxit = 1000000))
sProb <- spRes[[1]][[2]]
# but that's the sampling PROBABILITY per bin, not the instantaneous rate of change
# we can use sProb2sRate() to get the rate. We'll need to also tell it the int.length
sRate1 <- sProb2sRate(sProb,int.length = 1)
# we also need extinction rate and branching rate (see above)
# need to divide by int.length...
divRate <- spRes[[1]][[1]]/1
# estimates that r = 0.3... kind of low (simulated sampling rate is 0.5)
# Note: for real data, you may need to use an average int.length (no constant length)
ttree <- bin_cal3TimePaleoPhy(cladogram,rangesDisc,brRate = divRate,extRate = divRate,
sampRate = sRate1,ntrees = 1,plot = TRUE)
phylodiv(ttree)
# can also force the appearance timings not to be chosen stochastically
ttree1 <- bin_cal3TimePaleoPhy(cladogram,rangesDisc,brRate = divRate,extRate = divRate,
sampRate = sRate1,ntrees = 1,nonstoch.bin = TRUE,plot = TRUE)
phylodiv(ttree1)

# testing node.mins in bin_cal3TimePaleoPhy
ttree <- bin_cal3TimePaleoPhy(cladogram,rangesDisc,brRate = divRate,extRate = divRate,
sampRate = sRate1,ntrees = 1,node.mins = nodeDates,plot = TRUE)
# with randres = TRUE
ttree <- bin_cal3TimePaleoPhy(cladogram,rangesDisc,brRate = divRate,extRate = divRate,
sampRate = sRate1,ntrees = 1,randres = TRUE,node.mins = nodeDates,plot = TRUE)

# example with multiple values of anc.wt
ancWt <- sample(0:1,nrow(rangesDisc[[2]]),replace = TRUE)
names(ancWt) <- rownames(rangesDisc[[2]])
ttree1 <- bin_cal3TimePaleoPhy(cladogram,rangesDisc,brRate = divRate,extRate = divRate,
sampRate = sRate1,ntrees = 1,anc.wt = ancWt,plot = TRUE)

cladogeneticTraitCont  Simulate Cladogenetic Trait Evolution

Description

This function simulates trait evolution at each speciation/branching event in a matrix output from simFossilRecord, after transformation with fossilRecord2fossilTaxa.
Usage

cladogeneticTraitCont(taxa, rate = 1, meanChange = 0, rootTrait = 0)

Arguments

taxa A five-column matrix of taxonomic data, as output by fossilRecord2fossilTaxa after simulation with simFossilRecord (or via the depreciated function simFossil1Taxa)
rate rate of trait change; variance of evolutionary change distribution per speciation event
meanChange Mean change per speciation event. Default is 0; change to simulate 'active' speciational trends, where the expected change at each speciational event is non-zero.
rootTrait The trait value of the first taxon in the dataset; set to 0 by default.

Details

This function simulates continuous trait evolution where change occurs under a Brownian model, but only at events that create new distinct morphotaxa (i.e. species as recognized in the fossil record), either branching events or anagenesis (pseudospeciation). These are the types of morphological differentiation which can be simulated in the function simFossilRecord. This is sometimes referred to as cladogenetic or speciation trait evolution and is related to Punctuated Equilibrium theory. Anagenetic shifts are not cladogenetic events per se (no branching!), so perhaps the best way to think of this function is it allows traits to change anytime simFossilRecord created a new 'morphotaxon' in a simulation.

Importantly, trait changes only occur at the base of 'new' species, thus allowing cladogenetic trait evolution to be asymmetrical at branching points: i.e. only one branch actually changes position in trait-space, as expected under a budding cladogenesis model. This distinction is important as converting the taxa matrix to a phylogeny and simulating the trait changes under a 'speciational' tree-transformation would assume that divergence occurred on both daughter lineages at each node. (This has been the standard approach for simulating cladogenetic trait change on trees).

Cryptic taxa generated with prop.cryptic in simFossilRecord will not differ at all in trait values. These species will all be identical.

See this link for additional details: http://nemagraptus.blogspot.com/2012/03/simulating-budding-cladogenetictrait.html

Value

Returns a vector of trait values for each taxon, with value names being the taxa IDs (column 1 of the input) with a 't' pasted (as with rtree in the ape library).

Author(s)

David W. Bapst
See Also

simFossilRecord,

This function is similar to Brownian motion simulation functions such as rTraitCont in ape, 
sim.char in geiger and fastBM in phytools.

See also unitLengthTree in this package and speciationnalTree in the package geiger. These 
are tree transformation functions; together with BM simulation functions, they would be expected 
to have a similar effect as this function (when cladogenesis is 'bifurcating' and not 'budding'; see 
above).

Examples

```r
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(30,1000), plot = TRUE)
taxa <- fossilRecord2FossilTaxa(record)
trait <- cladogeneticTraitCont(taxa)
tree <- taxa2phylo(taxa)
plotTraitgram(trait,tree,conf.int = FALSE)

#with cryptic speciation
record <- simFossilRecord(p = 0.1, q = 0.1, prop.cryptic = 0.5,
nruns = 1, nTotalTaxa = c(30,1000), plot = TRUE)
taxa <- fossilRecord2FossilTaxa(record)
trait <- cladogeneticTraitCont(taxa)
tree <- taxa2phylo(taxa)
plotTraitgram(trait,tree,conf.int = FALSE)
```

---

**Description**

This is just a small collection of miscellaneous functions that may be useful, primarily for commu-
nity ecology analyses, particularly for paleoecological data. They are here mainly for pedagogical 
reasons (i.e. for students) as they don’t appear to be available in other ecology-focused packages.

**Usage**

```r
pairwiseSpearmanRho(x, dropAbsent = “bothAbsent”, asDistance = FALSE,
diag = NULL, upper = NULL, na.rm = FALSE)
HurlbertPIE(x, nAnalyze = Inf)
```
Arguments

x The community abundance matrix. Must be a matrix with two dimensions for `pairwiseSpearmanRho`; if a vector is supplied to `HurlbertPIE`, then it is treated as if it was matrix with a single row and number of columns equal to its length. Taxonomic units are assumed to be the columns and sites (samples) are assumed to be the rows, for both functions.

dropAbsent Should absent taxa be dropped? Must be one of either 'bothAbsent' (drop taxa absent in both sites for a given pairwise comparison), 'eitherAbsent' (drop taxa absent in either site), or 'noDrop' (drop none of the taxa). The default 'bothAbsent' is recommended, see examples.

asDistance Should the rho coefficients be rescaled on a scale similar to dissimilarity metrics, i.e. bounded 0 to 1, with 1 representing maximum dissimilarity (i.e. a Spearman rho correlation of -1)? ( dissimilarity = (1 - rho) / 2 )

diag Should the diagonal of the output distance matrix be included?

upper Should the upper triangle of the output distance matrix be included?

na.rm Should taxa listed with NA values be dropped from a pair-wise site comparison? If FALSE, the returned value for that site pair will be NA if NAs are present.

nanalyze Allows users to select that PIE be calculated only on `nanalyze` most abundant taxa in a site sample. `nanalyze` must be a vector of length 1, consisting of whole-number value at least equal to 2. By default, `nanalyze = Inf` so all taxa are accepted. Note that if there are less taxa in a sample than `nanalyze`, the number present will be used.

Details

pairwiseSpearmanRho returns Spearman rho correlation coefficients based on the rank abundances of taxa (columns) within sites (rows) from the input matrix, by internally wrapping the function `cor.test`. It allows for various options that automatically allow for dropping taxa not shared between two sites (the default), as well as several other options. This allows the rho coefficient to behave like the Bray-Curtis distance, in that it is not affected by the number of taxa absent in both sites.

pairwiseSpearmanRho can also rescale the rho coefficients with (1-rho)/2 to provide a measure similar to a dissimilarity metric, bounded between 0 and 1. This function was written so several arguments would be in a similar format to the vegan library function `vegdist`. If used to obtain rho rescaled as a dissimilarity, the default output will be the lower triangle of a distance matrix object, just as is returned by default by `vegdist`. This behavior can be modified via the arguments for including the diagonal and upper triangle of the matrix. Otherwise, a full matrix is returned (by default) if the asDistance argument is not enabled.

HurlbertPIE provides the 'Probability of Interspecific Encounter' metric for relative community abundance data, a commonly used metric for evenness of community abundance data based on derivations in Hurlbert (1971). An optional argument allows users to apply Hurlbert's PIE to only a subselection of the most abundant taxa.

Value

pairwiseSpearmanRho will return either a full matrix (the default) or (if `asDistance` is true, a distance matrix, with only the lower triangle shown (by default). See details.
HurlbertPIE returns a named vector of PIE values for the input data.

Author(s)

David W. Bapst

References


See Also

twoWayEcologyCluster; example dataset: kanto

Examples

```r
# let's load some example data:
# a classic dataset collected by Satoshi and Okido from the Kanto region
data(kanto)

rhoBothAbsent <- pairwiseSpearmanRho(kanto,dropAbsent = "bothAbsent")

# other dropping options
rhoEitherAbsent <- pairwiseSpearmanRho(kanto,dropAbsent = "eitherAbsent")
rhoNoDrop <- pairwiseSpearmanRho(kanto,dropAbsent = "noDrop")

# compare
layout(1:3)
lim <- c(-1,1)
plot(rhoBothAbsent, rhoEitherAbsent, xlim = lim, ylim = lim)
abline(0,1)
plot(rhoBothAbsent, rhoNoDrop, xlim = lim, ylim = lim)
abline(0,1)
plot(rhoEitherAbsent, rhoNoDrop, xlim = lim, ylim = lim)
abline(0,1)
layout(1)

#using dropAbsent = "eitherAbsent" reduces the number of taxa so much that
# the number of taxa present drops too low to be useful
# dropping none of the taxa restricts the rho measures to high coefficients
# due to the many shared 0s for absent taxa

#################

# Try the rho coefficients as a rescaled dissimilarity
rhoDist <- pairwiseSpearmanRho(kanto,asDistance = TRUE,dropAbsent = "bothAbsent")

# What happens if we use these in typical distance matrix based analyses?
```
# Cluster analysis
clustRes <- hclust(rhoDist)
plot(clustRes)

# Principle Coordinates Analysis
pcoRes <- pcocor(rhoDist, correction = "lingoes")
scores <- pcoRes$vector
#plot the PCO
plot(scores, type = "n")
text(labels = rownames(kanto), scores[,1], scores[,2], cex = 0.5)

# measuring evenness with Hurlbert's PIE
kantoPIE <- HurlbertPIE(kanto)

#histogram
hist(kantoPIE)
#evenness of the kanto data is fairly high

# barplot
parX <- par(mar = c(7,5,3,3))
barplot(kantoPIE, las = 3, cex.names = 0.7,
ylab = "Hurlbert's PIE", ylim = c(0.5,1), xpd = FALSE)
par(parX)

# and we can see that the Tower has extremely low unevenness
# ... over high abundance of ghosts?

# let's look at evenness of 5 most abundant taxa
kantoPIE_5 <- HurlbertPIE(kanto, nAnalyze = 5)

# barplot
parX <- par(mar = c(7,5,3,3))
barplot(kantoPIE_5, las = 3, cex.names = 0.7,
ylab = "Hurlbert's PIE for 5 most abundant taxa", ylim = c(0.5,1), xpd = FALSE)
par(parX)

---

compareTimescaling  
**Comparing the Time-Scaling of Trees**

**Description**

These functions take two trees and calculate the changes in node ages (for compareNodeAges) for shared clades or terminal branch lengths leading to shared tip taxa (for compareTermBranches).

**Usage**

```r
compareNodeAges(tree1, tree2, dropUnshared = FALSE)
```
compareTimescaling

\texttt{compareTermBranches(tree1, tree2)}

**Arguments**

\begin{itemize}
  \item \texttt{tree1} A time-scaled phylogeny of class 'phylo'
  \item \texttt{tree2} A time-scaled phylogeny of class 'phylo'; for \texttt{compareNodeAges}, \texttt{tree2} can also be an object of class 'multiPhylo' composed of multiple phylogenies. See below.
  \item \texttt{dropUnshared} If TRUE, nodes not shared across all input trees are dropped from the final output for \texttt{compareNodeAge}. This argument has no effect if \texttt{tree2} is a single phylogeny (a 'phylo'-class object).
\end{itemize}

**Details**

For their most basic usage, these functions compare the time-scaling of two trees. Any taxa not-shared on both trees are dropped before analysis, based on tip labels.

As with many paleotree functions, calculations relating to time on trees are done with respect to any included \$\text{root.time} elements. If these are not present, the latest tip is assumed to be at the present day (time = 0).

\texttt{compareNodeAges} calculates the changes in the clade ages among those clades shared by the two trees, relative to the first tree in absolute time. For example, a shift of +5 means the clade originates 5 time-units later in absolute time on the second tree, while a shift of -5 means the clade originated 5 time-units prior on the second tree.

For \texttt{compareNodeAges}, if \texttt{tree2} is actually a \texttt{multiPhylo} object composed of multiple phylogenies, the output will be a matrix, with each row representing a different tree and each column a different clade shared between at least some subset of the trees in \texttt{tree2} and the tree in \texttt{tree1}. values in the matrix are the changes in clade ages between from \texttt{tree1} (as baseline) to \texttt{tree2}, with NA values representing a clade that is not contained in the tree represented by that row (but is contained in \texttt{tree1} and at least one other tree in \texttt{tree2}). The matrix can be reduced to only those clades shared by all trees input via the argument \texttt{dropUnshared}. Note that this function distinguishes clades based on their shared taxa, and cannot so infer that two clades might be identical if it were not for single taxon within the crown of one considered clade, despite that such a difference should probably have no effect on compare a node divergence date. Users should consider their dataset for such scenarios prior to application of \texttt{compareNodeAges}, perhaps by dropping all taxa not included in all other trees to be considered (this is NOT done by this function).

\texttt{compareTermBranches} calculates the changes in the terminal branch lengths attached to tip taxa shared by the two trees, relative to the first tree. Thus, a shift of +5 means that this particular terminal taxon is connected to a terminal branch which is five time-units longer.

**Value**

\texttt{compareTermBranches} returns a vector of temporal shifts for terminal branches with the shared tip names as labels.

\texttt{compareNodeAges}, if both \texttt{tree1} and \texttt{tree2} are single trees, outputs a vector of temporal shifts for nodes on \texttt{tree2} with respect to \texttt{tree1}. If \texttt{tree2} is multiple trees, then a matrix is output, with each row representing each tree in \texttt{tree2} (and carrying the name of each tree, if any is given). The values are temporal shifts for each tree in \texttt{tree2} with respect to \texttt{tree1}. For either case, the column names or
element names (for a vector) are the sorted taxon names of the particular clade, the dates of which are given in that column. See above for more details. These names can be very long when large trees are considered.

See Also
datenodes, taxa2phylo, phyloDiv

Examples

```r
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(30,40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
# get the true tree
tree1 <- taxa2phylo(taxa)
# simulate a fossil record with imperfect sampling with sampleRanges()
rangesCont <- sampleRanges(taxa,r = 0.5)
# let's use taxa2cladogram to get the 'ideal' cladogram of the taxa
cladogram <- taxa2cladogram(taxa,plot = TRUE)
# Now let's try timePaleoPhy using the continuous range data
tree2 <- timePaleoPhy(cladogram,rangesCont,type = "basic")
# let's look at the distribution of node shifts
hist(compareNodeAges(tree1,tree2))
# let's look at the distribution of terminal branch lengths
hist(compareTermBranches(tree1,tree2))

# testing ability to compare multiple trees with compareNodeAges
trees <- cal3TimePaleoPhy(cladogram,rangesCont,brRate = 0.1,extRate = 0.1,
sampRate = 0.1,ntrees = 10)
nodeComparison <- compareNodeAges(tree1,trees)
# plot it as boxplots for each node
boxplot(nodeComparison,names = NULL);abline(h = 0)
# plot mean shift in node dates
abline(h = mean(apply(nodeComparison,2,mean),na.rm = TRUE)),lty = 2)

# just shifting a tree back in time
set.seed(444)
tree1 <- rtree(10)
tree2 <- tree1
tree1$root.time <- 10
compareNodeAges(tree1,tree2)
compareTermBranches(tree1,tree2)
```
**Description**

This function constrains a model to make submodels with fewer parameters, using a structure and syntax taken from the function constrain in Rich Fitzjohn’s package `diversitree`.

**Usage**

```r
constrainParPaleo(f, ..., formulae = NULL, names = parnames(f),
       extra = NULL)
```

**Arguments**

- **f**: A function to constrain. This function must be of S3 class `paleotreeFunc` and have all necessary attributes expected of that class, which include parameter names and upper and lower bounds. As I have deliberately not exported the function which creates this class, it should be impossible for regular users to obtain such objects easily without using one of the 'make' functions, which automatically output a function of the appropriate class and attributes.

- **...**: Formulae indicating how the function should be constrained. See details and examples for lengthy discussion.

- **formulae**: Optional list of constraints, possibly in addition to those in `...`.

- **names**: Optional Character vector of names, the same length as the number of parameters in `x`. Use this only if `parnames` does not return a vector for your function. Generally this should not be used. DWB: This argument is kept for purposes of keeping the function as close to the parent as possible but, in general, should not be used because the input function must have all attributes expected of class `paleotreeFunc`, including parameter names.

- **extra**: Optional vector of additional names that might appear on the RHS of constraints but do not represent names in the function's `argnames`. This can be used to set up dummy variables (example coming later).

**Details**

This function is based on (but does not depend on) the function `constrain` from the package `diversitree`. Users should refer to this parent function for more detailed discussion of model constraint usage and detailed examples.

The parent function was forked to add functionality necessary for dealing with the high parameter count models typical to some paleontological analyses, particularly the inverse survivorship method. This necessitated that the new function be entirely separate from its parent. Names of functions involved (both exported and not) have been altered to avoid overlap in the package namespaces. Going forward, the paleotree package maintainer (Bapst) will try to be vigilant with respect to changes in `constrain` in the original package, `diversitree`.

Useful information from the `diversitree` manual (11/01/13):

"If f is a function that takes a vector x as its first argument, this function returns a new function that takes a shorter vector x with some elements constrained in some way; parameters can be fixed to particular values, constrained to be the same as other parameters, or arbitrary expressions of free parameters."
In general, formulae should be of the structure:

\[ LHS \sim RHS \]

...where the LHS is the 'Parameter We Want to Constrain' and the RHS is whatever we are constraining the LHS to, usually another parameter. LHS and RHS are the 'left-hand side' and 'right-hand side' respectively (which I personally find obscure).

Like the original `constrain` function this function is based on, this function cannot remove constraints previously placed on a model object and there may be cases in which the constrained function may not make sense, leading to an error. The original function will sometimes issue nonsensical functions with an incorrect number/names of parameters if the parameters to be constrained are given in the wrong order in formulae.

**Differences from diversitree’s `constrain` Function:**

This forked paleotree version of `constrain` has two additional features, both introduced to aid in constraining models with a high number of repetitive parameters. (I did not invent these models, so don’t shoot the messenger.)

First, it allows nuanced control over the constraining of many parameters simultaneously, using the 'all' and 'match' descriptors. This system depends on parameters being named as such: name.group1.group2.group3 and so on. Each 'group' is reference to a system of groups, perhaps referring to a time interval, region, morphology, taxonomic group or some other discrete characterization among the data (almost all functions envisioned for paleotree are for per-taxon diversification models). The number of group systems is arbitrary, and may be from zero to a very large number; it depends on the 'make' function used and the arguments selected by the user. For example, the parameter 'x.1' would be for the parameter 'x' in the first group of the first group system (generally a time interval for most paleotree functions). For a more complicated exampled, with the parameter 'x.1.3.1', the third group for the second group system (perhaps this taxonomic data point has a morphological feature not seen in some other taxa) and group 1 of the third group system (maybe biogeographic region 1? the possibilities are endless depending on user choices).

The 'all' option work like so: if 'x.all~x.1' is given as a formulae, then all x parameters will be constrained to equal x.1. For example, if there is x.1, x.2, x.3 and x.4 parameters for a model, 'x.all~x.1' would be equivalent to individually giving the formulae 'x.2~x.1', 'x.3~x.1' and 'x.4~x.1'. This means that if there are many parameters of a particular type (say, 50 'x' parameters) it is easy to constrain all with a short expression. It is not necessary that the The 'all' can be used anywhere in the name of the parameter in a formulae, including to make all parameters for a given 'group' the same. Furthermore, the LHS and RHS don’t need to be same parameter group, and both can contain 'all' statements, even multiple 'all' statements. Consider these examples, all of which are legal:

- **x.all ~ y.1** Constrains all values of the parameter x for every group to be equal to the single value for the parameter y for group 1 (note that there’s only a single set of groups).
- **all.1 ~ x.1** Constrains all parameters for the first group to equal each other, here arbitrary named x.1. For example, if there is parameters named x.1, y.1 and z.1, all will be constrained to be equal to a single parameter value.
- **x.all.all ~ y.2.3** Constrains all values for x in any and all groups to equal the single value for y for group 2 (of system 1) and group 3 (of system 2).
- **x.all ~ y.all** Constrains all values of x for every group and y for every group to be equal to a single value, which by default will be reported as y.1

The 'match' term is similar, allowing parameter values from the same group to be quickly matched and made equivalent. These 'match' terms must have a matching (cue laughter) term both in the
corresponding LHS and RHS of the formula. For example, consider ‘x.match~y.match’ where there are six parameters: x.1, x.2, x.3, y.1, y.2 and y.3. This will effectively constrain x.1~y.1, x.2~y.2 and x.3~y.3. This is efficient for cases where we have some parameters that we often treat as equal. For example, in paleontology, we sometimes make a simplifying assumption that birth and death rates are equal in multiple time intervals. Some additional legal examples are:

**x.match.1 ~ y.match.1** This will constrain only parameters of x and y to to equal each other if they both belong to the same group for the first group system AND belong to group 1 of the first group.

**all.match. ~ x.match** This will constrain all named parameters in each group to equal each other; for example, if there are parameters x.1, y.1, z.1, x.2, y.2 and z.2, this will constrain them such that y.1~x.1, z.1~x.1, y.2~x.2 and z.2~x.2, leaving x.1 and x.2 as the only parameters effectively.

There are two less fortunate qualities to the introduction of the above terminology. Unfortunately, this package author apologizes that his programming skills are not good enough to allow more complex sets of constraints, as would be typical with the parent function, when ‘all’ or ‘match’ terms are included. For example, it would not be legal to attempt to constrain ‘y.all ~ x.1 / 2’, where the user presumably is attempting to constrain all y values to equal the x parameter to equal half of the x parameter for group 1. This will not be parsed as such and should return an error. However, there are workarounds, but they require using constrainParPaleo more than once. For the above example, a user could first use ‘y.all ~ y.1’ constraining all y values to be equal. Then a user could constrain with the formula ‘y.1 ~ x.1 / 2’ which would then constrain y.1 (and all the y values constrained to equal it) to be equal to the desired fraction.

Furthermore, this function expects that parameter names don’t already have period-separated terms that are identical to ‘all’ or ‘match’. No function in paleotree should produce such natively. If such were to occur, perhaps by specially replacing parameter names, constrainParPaleo would confuse these terms for the specialty terms described here.

Secondly, this altered version of constrain handles the parameter bounds included as attributes in functions output by the various ’make’ functions. This means that if ‘x.1 ~ y.1’, constrainParPaleo will test if the bounds on x.1 and y.1 are the same. If the bounds are not the same, constrainParPaleo will return an error. This is important, as some models in paleotree may make a parameter a rate (bounded zero to some value greater than one) or a probability (bounded zero to one), depending on user arguments. Users may not realize these differences and, in many cases, constraining a rate to equal a probability is nonsense (absolute poppycock!). If a user really wishes to constrain two parameters with different bounds to be equal (I have no idea why anyone would want to do this), they can use the parameter bound replacement functions described in modelMethods to set the parameter bounds as equal. Finally, once parameters with the same bounds are constrained, the output has updated bounds that reflect the new set of parameters for the new constrained function.

**Value**

Modified from the diversitree manual: This function returns a constrained function that can be passed through to the optimization functions of a user’s choice, such as `optim`, `find.mle` in diversitree or `mcmc`. It will behave like any other function. However, it has a modified `class` attribute so that some methods will dispatch differently: `parnames`, for example, will return the names of the parameters of the constrained function and `parInit` will return the initial values for those same constrained set of parameters. All arguments in addition to `x` will be passed through to the original function `f`.
Additional useful information from the diversitree manual (11/01/13):

For help in designing constrained models, the returned function has an additional argument `parsNonly`, when this is TRUE the function will return a named vector of arguments rather than evaluate the function (see Examples).

**Author(s)**

This function (and even this help file!) was originally written by Rich Fitzjohn for his library diversitree, and subsequently rewritten and modified by David Bapst.

**References**


**See Also**

As noted above, this function is based on (but does not depend on) the function `constrain` from the library diversitree.

**Examples**

```r
#simulation example with make_durationFreqCont, with three random groups
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(30, 40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
rangesCont <- sampleRanges(taxa,r = 0.5)
gr1 <- matrix(sample(1:3,nrow(taxa),replace = TRUE),1) #groupings matrix
likFun <- make_durationFreqCont(rangesCont,groups = gr1)

# can constrain both extinction rates to be equal
constrainFun <- constrainParPaleo(likFun,q=2-q1)

# see the change in parameter names and bounds
parnames(likFun)
parnames(constrainFun)
parbounds(likFun)
parbounds(constrainFun)

# some more ways to constrain stuff!

#constrain all extinction rates to be equal
constrainFun <- constrainParPaleo(likFun,q.all=q1)
parnames(constrainFun)

#constrain all rates for everything to be a single parameter
constrainFun <- constrainParPaleo(likFun,r.all=r.all)
parnames(constrainFun)

#constrain all extinction rates to be equal & all sampling to be equal
constrainFun <- constrainParPaleo(likFun,q.all=q1,r=all=r)
```

---

**constrainParPaleo**

This function (and even this help file!) was originally written by Rich Fitzjohn for his library diversitree, and subsequently rewritten and modified by David Bapst.

**Additional useful information from the diversitree manual (11/01/13):**

For help in designing constrained models, the returned function has an additional argument `parsNonly`, when this is TRUE the function will return a named vector of arguments rather than evaluate the function (see Examples).

**Author(s)**

This function (and even this help file!) was originally written by Rich Fitzjohn for his library diversitree, and subsequently rewritten and modified by David Bapst.

**References**


**See Also**

As noted above, this function is based on (but does not depend on) the function `constrain` from the library diversitree.

**Examples**

```r
#simulation example with make_durationFreqCont, with three random groups
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(30, 40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
rangesCont <- sampleRanges(taxa,r = 0.5)
gr1 <- matrix(sample(1:3,nrow(taxa),replace = TRUE),1) #groupings matrix
likFun <- make_durationFreqCont(rangesCont,groups = gr1)

# can constrain both extinction rates to be equal
constrainFun <- constrainParPaleo(likFun,q=2-q1)

# see the change in parameter names and bounds
parnames(likFun)
parnames(constrainFun)
parbounds(likFun)
parbounds(constrainFun)

# some more ways to constrain stuff!

#constrain all extinction rates to be equal
constrainFun <- constrainParPaleo(likFun,q.all=q1)
parnames(constrainFun)

#constrain all rates for everything to be a single parameter
constrainFun <- constrainParPaleo(likFun,r.all=r.all)
parnames(constrainFun)

#constrain all extinction rates to be equal & all sampling to be equal
constrainFun <- constrainParPaleo(likFun,q.all=q1,r.all=r)
```
createMrBayesConstraints

Transform a Topology into a Set of Constraint Commands for MrBayes

Description

Takes a phylogeny in the form of an object of class phylo and outputs a set of topological constraints for MrBayes as a set of character strings, either printed in the R console or in a named text file, which can be used as commands in the MrBayes block of a NEXUS file for use with (you guessed it!) MrBayes.

Usage

createMrBayesConstraints(tree, partial = TRUE, file = NULL, includeIngroupConstraint = FALSE)

Arguments

tree An object of class phylo.

partial If TRUE (the default), then constraints will be defined as partial constraints with respect to the rest of the taxa in the input tree. If FALSE, constraints will be defined as hard clade membership constraints (i.e. no additional taxa will be allowed to belong to the nodes present in the observed tree. Depending on your analysis, partial = TRUE may require additionally defining an outgroup.

file Filename (possibly with path) as a character string to a file which will be overwritten with the output constraint lines. If NULL, constraint lines are printed in the console.

includeIngroupConstraint When writing the prset line, should a group named 'ingroup' be included, which presumes an ingroup constraint was defined by the user for sake or rooting the tree? This is mainly used for use with paleotree function createMrBayesTIPDatingNexus for automating the construction of tip-dating analyses, which must constrain the ingroup.
createMrBayesTipCalibrations

Details

Partial = TRUE may be useful if the reason for using `createMrBayesConstraints` is to constrain a topology containing some of the taxa in an analysis, while allowing other taxa to freely vary. For example, Slater (2013) constrained an analysis so extant taxon relationships were held constant, using a molecular-based topology, while allowing fossil taxa to freely vary relative to their morphological character data.

Value

If argument `file` is NULL, then the constrain commands are output as a series of character strings.

Author(s)

David W. Bapst, with some inspiration from Graham Slater. This code was produced as part of a project funded by National Science Foundation grant EAR-1147537 to S. J. Carlson.

References


See Also

createMrBayesTipDatingNexus, createMrBayesTipCalibrations

Examples

```r
set.seed(444)
tree <- rtree(10)
createMrBayesConstraints(tree)
createMrBayesConstraints(tree, partial = FALSE)

## Not run:

createMrBayesConstraints(tree, file = "topoConstraints.txt")

## End(Not run)
```

---

`createMrBayesTipCalibrations`

*Construct A Block of Tip Age Calibrations for Use with Tip-Dating Analyses in MrBayes*
createMrBayesTipCalibrations

Description
Takes a set of tip ages (in several possible forms, see below), and outputs a set of tip age calibrations for use with tip-dating analyses (sensu Zhang et al., 2016) in the popular phylogenetics program MrBayes. These calibrations are printed as a set of character strings, as well as a line placing an offset exponential prior on the tree age, either printed in the R console or in a named text file, which can be used as commands in the MrBayes block of a NEXUS file for use with (you guessed it!) MrBayes.

Usage
createMrBayesTipCalibrations(tipTimes, ageCalibrationType, whichAppearance = "first", treeAgeOffset, minTreeAge = NULL, collapseUniform = TRUE, anchorTaxon = TRUE, file = NULL)

Arguments
tipTimes
This input may be either a timeList object (i.e. a list of length 2, composed of a table of interval upper and lower time boundaries (i.e. earlier and latter bounds), and a table of first and last intervals for taxa) or a matrix with rownames for taxa as you want listed in the MrBayes block, with either one, two or four columns containing ages (respectively) for point occurrences with precise dates (for a single column), uncertainty bounds on a point occurrence (for two columns), or uncertainty bounds on the first and last occurrence (for four columns). Note that precise first and last occurrence dates should not be entered as a two column matrix, as this will instead be interpreted as uncertainty bounds on a single occurrence. Instead, either select which you want to use for tip-dates and give a 1-column matrix, or repeat (and collate) the columns, so that the first and last appearances has uncertainty bounds of zero.
ageCalibrationType
This argument decides how age calibrations are defined, and currently allows for four options: "fixedDateEarlier" which fixes tip ages at the earlier (lower) bound for the selected age of appearance (see argument whichAppearance for how that selection is made), "fixedDateLatter" which fixes the date to the latter (upper) bound of the selected age of appearance, "fixedDateRandom" which fixes tips to a date that is randomly drawn from a uniform distribution bounded by the upper and lower bounds on the selected age of appearance, or (the recommended option) "uniformRange" which places a uniform prior on the age of the tip, bounded by the latest and earliest (upper and lower) bounds on the the selected age.
whichAppearance
Which appearance date of the taxa should be used: their 'first' or their 'last' appearance date? The default option is to use the 'first' appearance date. Note that use of the last appearance date means that tips will be constrained to occur before their last occurrence, and thus could occur long after their first occurrence (!).
treeAgeOffset
A parameter given by the user controlling the offset between the minimum and expected tree age prior. mean tree age for the offset exponential prior on tree age will be set to the minimum tree age, plus this offset value. Thus, an offset
of 10 million years would equate to a prior assuming that the expected tree age is around 10 million years before the minimum age.

`minTreeAge`  
if NULL (the default), then `minTreeAge` will be set as the oldest date among the tip age used (those used being determine by user choices (or oldest bound on a tip age). Otherwise, the user can supply their own minimum tree, which must be greater than whatever the oldest tip age used is.

`collapseUniform`  
MrBayes won't accept uniform age priors where the maximum and minimum age are identical (i.e. its actually a fixed age). Thus, if this argument is TRUE (the default), this function will treat any taxon ages where the maximum and minimum are identical as a fixed age, and will override setting `ageCalibrationType = "uniformRange"` for those dates. All taxa with their ages set to fixed by the behavior of `anchorTaxon` or `collapseUniform` are returned as a list within a commented line of the returned MrBayes block.

`anchorTaxon`  
This argument may be a logical (default is TRUE, or a character string of length = 1. This argument has no effect if `ageCalibrationType` is not set to "uniformRange", but the argument may still be evaluated. If `ageCalibrationType = "uniformRange"`, MrBayes will do a tip-dating analysis with uniform age uncertainties on all taxa (if such uncertainties exist; see `collapseUniform`). However, MrBayes does not record how each tree sits on an absolute time-scale, so if the placement of every tip is uncertain, lining up multiple dated trees sampled from the posterior (where each tip’s true age might differ) could be a nightmare to back-calculate, if not impossible. Thus, if `ageCalibrationType = "uniformRange"`, and there are no tip taxa given fixed dates due to `collapseUniform` (i.e. all of the tip ages have a range of uncertainty on them), then a particular taxon will be selected and given a fixed date equal to its earliest appearance time for its respective `whichAppearance`. This taxon can either be indicated by the user or instead the first taxon listed in `tiptimes` will be arbitrary selected. All taxa with their ages set to fixed by the behavior of `anchorTaxon` or `collapseUniform` are returned as a list within a commented line of the returned MrBayes block.

`file`  
Filename (possibly with path) as a character string to a file which will be overwritten with the output tip age calibrations. If NULL, tip calibration commands are output to the console.

**Details**

Beware: some combinations of arguments might not make sense for your data.

(But that’s always true, is it not?)

**Value**

If argument `file` is NULL, then the tip age commands are output as a series of character strings.

All taxa with their ages set to fixed by the behavior of `anchorTaxon` or `collapseUniform` are returned as a list within a commented line of the returned MrBayes block.
Author(s)

David W. Bapst. This code was produced as part of a project funded by National Science Foundation grant EAR-1147537 to S. J. Carlson.

References


See Also

createMrBayesConstraints, createMrBayesTipDatingNexus

Examples

```r
# load retiolitid dataset
data(retiolitiniae)

# uniform prior, with a 10 million year offset for
# the expected tree age from the earliest first appearance
createMrBayesTipCalibrations(tipTimes = retioRanges, whichAppearance = "first", ageCalibrationType = "uniformRange", treeAgeOffset = 10)

# fixed prior, at the earliest bound for the first appearance
createMrBayesTipCalibrations(tipTimes = retioRanges, whichAppearance = "first", ageCalibrationType = "fixedDateEarlier", treeAgeOffset = 10)

# fixed prior, sampled from between the bounds on the last appearance
# you should probably never do this, fyi
createMrBayesTipCalibrations(tipTimes = retioRanges, whichAppearance = "first", ageCalibrationType = "fixedDateRandom", treeAgeOffset = 10)

## Not run:
createMrBayesTipCalibrations(tipTimes = retioRanges, whichAppearance = "first", ageCalibrationType = "uniformRange", treeAgeOffset = 10, file = "tipCalibrations.txt")

## End(Not run)
```
Construct a Fully Formatted NEXUS Script for Performing Tip-Dating Analyses With MrBayes

Description
This function is meant to expedite the creation of NEXUS files formatted for performing tip-dating analyses in the popular phylogenetics software MrBayes, particularly clock-less tip-dating analyses executed with 'empty' morphological matrices (i.e. where all taxa are coded for a single missing character), although a pre-existing morphological matrix can also be input by the user (see argument `origNexusFile`). Under some options, this pre-existing matrix may be edited by this function. The resulting full NEXUS script is output as a set of character strings either printed to the R console, or output to file which is then overwritten.

Usage
```r
createmrbayestipdatingnexus(tipTimes = NULL, 
outgrouptaxa = NULL, 
treeConstraints = NULL, 
ageCalibrationType = "first", 
treeAgeOffset = NULL, 
collapseUniform = TRUE, 
anchorTaxon = TRUE, 
newFile = NULL, 
origNexusFile = NULL, 
parsedOriginalNexus = TRUE, 
createEmptyMorphMat = TRUE, 
orderedChars = NULL, 
morphModel = "strong", 
runName = NULL, 
gen = "10000000", 
doNotRun = FALSE, 
cleanNames = TRUE, 
printExecute = TRUE)
```

Arguments

**tipTimes**  This input may be either a timeList object (i.e. a list of length 2, composed of a table of interval upper and lower time boundaries (i.e. earlier and latter bounds), and a table of first and last intervals for taxa) or a matrix with rownames for taxa as you want listed in the MrBayes block, with either one, two or four columns containing ages (respectively) for point occurrences with precise dates (for a single column), uncertainty bounds on a point occurrence (for two columns), or uncertainty bounds on the first and last occurrence (for four columns). Note that precise first and last occurrence dates should not be entered as a two column matrix, as this will instead be interpreted as uncertainty bounds on a single occurrence. Instead, either select which you want to use for tip-dates and give a 1-column matrix, or repeat (and collate) the columns, so that the first and last appearances has uncertainty bounds of zero.

**outgroupTaxa**  A vector of type 'character', containing taxon names designating the outgroup. All taxa not listed in the outgroup will be constrained to be a monophyletic ingroup, for sake of rooting the resulting dated tree. Either `treeConstraints` or `outgroupTaxa` must be defined, but *not both*. If the outgroup-ingroup split is not present on the supplied `treeConstraints`, add that split to `treeConstraints` manually.
treeConstraints

An object of class phylo, from which (if treeConstraints is supplied) the set topological constraints are derived, as as described for argument tree for function createMrBayesConstraints. Either treeConstraints or outgroupTaxa must be defined, but not both. If the outgroup-ingroup split is not present on the supplied treeConstraints, add that split to treeConstraints manually.

ageCalibrationType

This argument decides how age calibrations are defined, and currently allows for four options: "fixedDateEarlier" which fixes tip ages at the earlier (lower) bound for the selected age of appearance (see argument whichAppearance for how that selection is made), "fixedDateLatter" which fixes the date to the latter (upper) bound of the selected age of appearance, "fixedDateRandom" which fixes tips to a date that is randomly drawn from a uniform distribution bounded by the upper and lower bounds on the selected age of appearance, or (the recommended option) "uniformRange" which places a uniform prior on the age of the tip, bounded by the latest and earliest (upper and lower) bounds on the the selected age.

whichAppearance

Which appearance date of the taxa should be used: their 'first' or their 'last' appearance date? The default option is to use the 'first' appearance date. Note that use of the last appearance date means that tips will be constrained to occur before their last occurrence, and thus could occur long after their first occurrence (!). In addition, createMrBayesTipDatingNexus allows for two options for this argument that are in addition to those offered by createMrBayesTipCalibrations. Both of these options will duplicate the taxa in the inputs multiple times, modifying their OTU labels, thus allowing multiple occurrences of long-lived morphotaxa to be listed as multiple OTUs arrayed across their stratigraphic duration. If whichAppearance = "firstLast", taxa will be duplicated so each taxon is listed as occurring twice: once at their first appearance, and a second time at their last appearance. Note that if a taxon first and last appears in the same interval, and ageCalibrationType = "uniformRange", then the resulting posterior trees may place the OTU assigned to the last occurrence before the first occurrence (but the assignment, in that case, was entirely arbitrary). When whichAppearance = "rangeThrough", each taxon will be duplicated into as many OTUs as each interval that a taxon ranges through (in a timeList format, see other paleotree functions), with the corresponding age uncertainties for those intervals. If the input tipTimes is not a list of length = 2, however, the function will return an error under this option.

treeAgeOffset

A parameter given by the user controlling the offset between the minimum and expected tree age prior. mean tree age for the offset exponential prior on tree age will be set to the minimum tree age, plus this offset value. Thus, an offset of 10 million years would equate to a prior assuming that the expected tree age is around 10 million years before the minimum age.

minTreeAge

if NULL (the default), then minTreeAge will be set as the oldest date among the tip age used (those used being determine by user choices (or oldest bound on a tip age). Otherwise, the user can supply their own minimum tree, which must be greater than whatever the oldest tip age used is.
collapseUniform

MrBayes won’t accept uniform age priors where the maximum and minimum age are identical (i.e. its actually a fixed age). Thus, if this argument is TRUE (the default), this function will treat any taxon ages where the maximum and minimum are identical as a fixed age, and will override setting ageCalibrationType = "uniformRange" for those dates. All taxa with their ages set to fixed by the behavior of anchorTaxon or collapseUniform are returned as a list within a commented line of the returned MrBayes block.

anchorTaxon

This argument may be a logical (default is TRUE, or a character string of length = 1. This argument has no effect if ageCalibrationType is not set to "uniformRange", but the argument may still be evaluated. If ageCalibrationType = "uniformRange", MrBayes will do a tip-dating analysis with uniform age uncertainties on all taxa (if such uncertainties exist; see collapseUniform). However, MrBayes does not record how each tree sits on an absolute time-scale, so if the placement of every tip is uncertain, lining up multiple dated trees sampled from the posterior (where each tip’s true age might differ) could be a nightmare to back-calculate, if not impossible. Thus, if ageCalibrationType = "uniformRange", and there are no tip taxa given fixed dates due to collapseUniform (i.e. all of the tip ages have a range of uncertainty on them), then a particular taxon will be selected and given a fixed date equal to its earliest appearance time for its respective whichAppearance. This taxon can either be indicated by the user or instead the first taxon listed in tipTimes will be arbitrary selected. All taxa with their ages set to fixed by the behavior of anchorTaxon or collapseUniform are returned as a list within a commented line of the returned MrBayes block.

newFile

Filename (possibly with path) as a character string leading to a file which will be overwritten with the output tip age calibrations. If NULL, tip calibration commands are output to the console.

origNexusFile

Filename (possibly with path) as a character string leading to a NEXUS text file, presumably containing a matrix of character date formatted for MrBayes. If supplied (it does not need to be supplied), the listed file is read as a text file, and concatenated with the MrBayes script produced by this function, so as to reproduce the original NEXUS matrix for executing in MrBayes. Note that the taxa in this NEXUS file are NOT checked against the user input tipTimes and treeConstraints, so it is up to the user to ensure the taxa are the same across the three data sources.

parseOriginalNexus

If TRUE (the default), the original NEXUS file is parsed and the taxon names listed within in the matrix are compared against the other inputs for matching (completely, across all inputs that include taxon names). Thus, it is up to the user to ensure the same taxa are found in all inputs. However, some NEXUS files may not parse correctly (particularly if character data for taxa stretches across more than a single line in the matrix). This may necessitate setting this argument to FALSE, which will instead do a straight scan of the NEXUS matrix without parsing it, and without checking the taxon names against other outputs. Some options for whichAppearance will not be available, however.

createEmptyMorphMat

If origNexusFile is not specified (implying there is no pre-existing morphological character matrix for this dataset), then an 'empty' NEXUS-formatted
createMrBayesTipDatingNexus

matrix will be appended to the set of MrBayes commands if this command is TRUE (the default). This 'empty' matrix will have each taxon in tipTimes coded for a single missing character (i.e., '?'). This allows tip-dating analyses with hard topological constraints, and ages determined entirely by the fossilized birth-death prior, with no impact from a presupposed morphological clock (thus a 'clock-less analysis').

orderedChars Should be a vector of numbers, indicating which characters should have their character-type in MrBayes changed to 'ordered'. If NULL, the default, then all characters will be treated as essentially unordered. No character ID should be listed that is higher than the number of characters in the matrix provided in origNexusFile. If origNexusFile is not provided, while orderedChars is defined, then an error will be returned.

morphModel This argument can be used to switch between two end-member models of morphological evolution in MrBayes, here named 'strong' and 'relaxed', for the 'strong assumptions' and 'relaxed assumptions' models described by Bapst, Schreiber and Carlson (Systematic Biology). The default is a model which makes very 'strong' assumptions about the process of morphological evolution, while the 'relaxed' alternative allows for considerably more heterogeneity in the rate of morphological evolution across characters, and in the forward and reverse transition rates between states. Note that in both cases, the character data is assumed to be filtered to only parsimony-informative characters, without autapomorphies.

runName The name of the run, used for naming the log files. If not set, the name will be taken from the name given for outputting the NEXUS script (newFile). If newFile is not given, and runName is not set by the user, the default run name will be "new_run_paleotree".

ngen Number of generations to set the MCMCMC to run for. Default (ngen = 100000000) is very high.

donotrun If TRUE, the commands that cause a script to automatically begin running in MrBayes will be left out. Useful for troubleshooting initial runs of scripts for non-fatal errors and warnings (such as ignored constraints). Default for this argument is FALSE.

cleanNames If TRUE (the default), then special characters (currently, this only contains the forward-slashes: '/') are removed from taxon names before construction of the NEXUS file.

printExecute If TRUE (the default) and if output is directed to a newFile (i.e. a newFile is specified), a line for pasting into MrBayes for executing the newly created file will be messaged to the terminal.

Details

Users must supply a data set of tip ages (in various formats), which are used to construct age calibrations commands on the tip taxa (via paleotree function createMrBayesTipCalibrations). The user must also supply some topological constraint: either a set of taxa designated as the outgroup, which is then converted into a command constraining the monophyly on the ingroup taxa, which is presumed to be all taxa not listed in the outgroup. Alternatively, a user may supply a tree which is then converted into a series of hard topological constraints (via function createMrBayesConstraints).
Both types of topological constraints cannot be applied. Many of the options available with `createMrBayesTipCalibrations` are available with this function, allowing users to choose between fixed calibrations or uniform priors that approximate stratigraphic uncertainty. In addition, the user may also supply a path to a text file presumed to be a NEXUS file containing character data formatted for use with `MrBayes`.

The taxa listed in `tipTimes` must match the taxa in `treeConstraints`, if such is supplied. If supplied, the taxa in `outgroupTaxa` must be contained within this same set of taxa. These all must have matches in the set of taxa in `origNexusFile`, if provided and if `parseOriginalNexus` is `TRUE`.

Note that because the same set of taxa must be contained in all inputs, relationships are constrained as 'hard' constraints, rather than 'partial' constraints, which allows some taxa to float across a partially fixed topology. See the documentation for `createMrBayesConstraints`, for more details.

**Value**

If argument `newFile` is `NULL`, then the text of the generated NEXUS script is output to the console as a series of character strings.

**Note**

This function allows a user to take an undated phylogenetic tree in R, and a set of age estimates for the taxa on that tree, and produce a posterior sample of dated trees using the MCMC in `MrBayes`, while treating an 'empty' morphological matrix as an uninformative set of missing characters. This 'clock-less tip-dating' approach is essentially an alternative to the `cal3` method in paleotree, sharing the same fundamental theoretical model (a version of the fossilized birth-death model), but with a better algorithm that considers the whole tree simultaneously, rather than evaluating each node individually, from the root up to the tips (as `cal3` does it, and which may cause artifacts). That said, `cal3` still has a few advantages: tip-dating as of April 2017 still only treats OTUs as point observations, contained in a single time-point, while `cal3` can consider taxa as having durations with first and last occurrences. This means it may be more straightforward to assess the extent of budding cladogenesis patterns of ancestor-descendant relationships in `cal3`, than in tip-dating.

**Author(s)**

David W. Bapst. This code was produced as part of a project funded by National Science Foundation grant EAR-1147537 to S. J. Carlson.

The basic `MrBayes` commands utilized in the output script are a collection of best practices taken from studying NEXUS files supplied by April Wright, William Gearty, Graham Slater, Davey Wright, and guided by the recommendations of Matzke and Wright, 2016 in Biol. Lett.

**References**

The basic fundamentals of tip-dating, and tip-dating with the fossilized birth-death model are introduced in these two papers:


For recommended best practices in tip-dating analyses, please see:

The rationale behind the two alternative morphological models are described in more detail here:

**See Also**

This function wraps various aspects of the functions `createMrBayesConstraints` and `createMrBayesTipCalibrations`. In many ways, this functionality is a replacement for the probabilistic dating method `cal3`.

For putting the posterior estimated trees on an absolute time scale, see functions `obtainDatedPosteriorTreesMrB`, with `getFixedTimes = TRUE` if you used a taxon with a fixed age, and function `setRootAges`.

**Examples**

```r
# let's do some examples

# load retiolitid dataset
data(retiolitinae)

# let's try making a NEXUS file!

# Use a uniform prior, with a 10 million year offset for
# the expected tree age from the earliest first appearance
# set average tree age to be 10 Ma earlier than first FAD

outgroupRetio <- "Rotaretiolites" # sister to all other included taxa

# the following will create a NEXUS file with an 'empty' morph matrix
# with the only topological constraint on ingroup monophyly
# Probably shouldn't do this: leaves too much to the FBD prior

# with doNotRun set to TRUE for troubleshooting

cREATEMrBayesTipDatingNexus(tipTimes = retioRanges,
outgroupTaxa = outgroupRetio, treeConstraints = NULL,
ageCalibrationType = "uniformRange", whichAppearance = "first",
treeAgeOffset = 10, newFile = NULL, origNexusFile = NULL, createEmptyMorphMat = TRUE,
runName = "retio_dating", doNotRun = TRUE)

# let's try it with a tree for topological constraints
# this requires setting outgroupTaxa to NULL
# let's also set doNotRun to FALSE

cREATEMrBayesTipDatingNexus(tipTimes = retioRanges,
outgroupTaxa = NULL, treeConstraints = retioTree,
```
dateNodes

Absolute Dates for Nodes of a Time-Scaled Phylogeny

Description

This function returns the ages of nodes (both internal and terminal tips) for a given phylogeny of class ‘phylo’. Specialized for use with time-scaled trees from paleotree, see Details.
Usage

dateNodes(tree, rootAge = tree$root.time, labelDates = FALSE, tolerance = 0.001)

Arguments

tree A phylogeny object of class 'phylo'. Must have edge.lengths!

rootAge The root age of the tree, assumed by default to be equal to the element tree$root.time, which is a standard element for trees time-scaled by the paleotree package. If not given by the user and if the $root.time element does not exist, then the maximum depth of the tree will be taken as the root age, which implicitly assumes the latest most terminal tip is an extant taxon at the modern day (time = 0). If rootAge is so defined that some nodes may occur later than time = 0, this function may return negative dates.

labelDates If FALSE (the default), the dates returned are labeled with the tip/node numbers as in tree$edge. If TRUE, they are labeled with the tip labels of every descendant tip, which for terminal tips means a single taxon label, and for internal tips a label that might be very long, composed of multiple tip labels pasted together. Thus, by default, this argument is FALSE.

tolerance The tolerance within which a node date has to be removed from time = 0 (i.e. the modern) to issue a warning that there are ‘negative’ node dates.

Details

This function is specialized for phylo objects time-scaled or simulated with functions from paleotree, and thus have a $root.time element. This function will still work without such, but users should see the details for the rootAge argument.

Value

Returns a vector of length Ntip(tree) + Nnode(tree) which contains the dates for all terminal tip nodes and internal nodes for the tree, in that order, as numbered in the tree$edge matrix. These dates are always on a descending scale (i.e. time before present); see help for argument rootAge for how the present time is determined. If rootAge is so defined that some nodes may occur later than time = 0 units before present, this function may (confusingly) return negative dates and a warning message will be issued.

Author(s)

David W. Bapst, based on a function originally written by Graeme Lloyd.

See Also

compareTimescaling, nodeDates2branchLengths
Examples

```r
# let's simulate some example data
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1, nTotalTaxa = c(30, 40), nExtant = 0)
taxa <- fossilRecord2fossilTaxa(record)
# get the true time-scaled tree
tree1 <- taxa2phylo(taxa)

# now let's try dateNodes
dateNodes(tree1)

# let's ignore $root.time
dateNodes(tree1, rootAge = NULL)

# with the lengthy tip-label based labels
# some of these will be hideously long
dateNodes(tree1, labelDates = TRUE)
```

degradeTree  
Randomly Collapse a Portion of Nodes on a Phylogeny

Description
degradeTree removes a proportion of the total nodes in a tree, chosen randomly, collapsing the nodes to produce a less-resolved tree. The related function `collapseNodes` given a tree and a vector of nodes to collapse, removes those nodes from a tree, creating a polytomy.

Usage
degradeTree(tree, prop_collapse = NULL, nCollapse = NULL, node.depth = NA, leave.zlb = FALSE)

collapseNodes(tree, nodeID, collapseType, leave.zlb = FALSE)

Arguments

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>tree</td>
<td>A phylogeny of class phylo</td>
</tr>
<tr>
<td>prop_collapse</td>
<td>Proportion of nodes to collapse</td>
</tr>
<tr>
<td>nCollapse</td>
<td>Number of nodes to collapse, can be supplied as an alternative to prop_collapse</td>
</tr>
<tr>
<td>node.depth</td>
<td>A number between 0 to 1, which conditions the depth of nodes removed. If NA, no conditioning (this is the default).</td>
</tr>
<tr>
<td>leave.zlb</td>
<td>If FALSE, the default option, the original branch length distribution is destroyed and branches set to zero by this function will return polytomies. If TRUE, then the original edge lengths are kept for unmodified edges, and modified edges are changed to zero length, and are not collapsed into polytomies. The removed branch length is not shifted to other edges.</td>
</tr>
</tbody>
</table>
nodeID  The node ID number(s) to be collapsed into a polytomy, as identified in the $edge matrix of the phylo object. Must be a vector of one or more ID numbers.

collapseType  Whether to collapse the edge leading the listed node (if collapseType = "forward"), or to collapse the child edges leading away from the node (if collapseType = "backward"). Collapsing a node 'into' a polytomy conceptually could be either and users should heed this option carefully. A third option, if "collapseType = clade" is to collapse the entire clade that is descended from a node (i.e. forward).

Details

In the function degradeTree, the nodes are removed at random using the basic R function sample. degradeTree can be conditioned to remove nodes of a particular depth with greater probability/frequency by setting node.depth to a value between zero (favoring the removal of deep nodes close to the root) or one (shallow nodes far from the root). Depth is evaluated based on the number of descendant tips. If node.depth is not NA, the relative proportion of descendants from each node is calculated, summed to 1 and the node.depth value subtracted from this proportion. These values are then squared, normalized again to equal to 1 and then used as the probabilities for sampling nodes for removal.

By default, branch lengths are removed from the input tree prior to degradation and entirely absent from the output tree. This is changed if argument leave.zlb is TRUE.

Value

Returns the modified tree as an object of class phylo, with no edge lengths by default.

Author(s)

David W. Bapst

See Also

di2multi, timeLadderTree

Examples

set.seed(444)
tree <- rtree(100)
tree1 <- degradeTree(tree, prop-collapse = 0.5)
tree3 <- degradeTree(tree, nCollapse = 50)

# let's compare the input and output
layout(matrix(1:2,2))
plot(tree, show.tip.label = FALSE, use.edge.length = FALSE)
plot(tree1, show.tip.label = FALSE, use.edge.length = FALSE)

# now with collapseNodes
tree <- rtree(100)
# collapse nodes backwards
# let's collapse lucky node number 13!
depthRainbow

Paint Tree Branch Depth by Color

Description

Paints the edges of a phylogeny with colors relative to their depth.

Usage

depthRainbow(tree)

Arguments

tree A phylo object

Details

The only purpose of this function is to make an aesthetically-pleasing graphic of one’s tree, where branches are color-coded with a rainbow palette, relative to their depth. Depth is defined relative to the number of branching nodes between the basal node of a branch and the root, not the absolute distance (i.e. branch length) to the root or the distance from the tips.

Value

No value returned, just plots a colorful phylogeny.

Examples

set.seed(444)
tree <- rtree(500)
depthRainbow(tree)
divCurveFossilRecordSim

_Diversity-Curve Plotting for Simulations of Diversification and Sampling In the Fossil Record_

Description

An extremely simple plotting function, which plots the original taxonomic diversity versus the sampled taxonomic diversity, for use with output from the function `simfossilrecord`. If sampling processes were not included in the model, then it plots simply the single diversity curve.

Usage

divCurveFossilRecordSim(fossilRecord, merge.cryptic = TRUE,
plotLegend = TRUE, legendPosition = "topleft", curveColors = c("black", "red"), curveLineTypes = c(1, 2))

Arguments

- **fossilRecord**: A list object output by `simFossilRecord`, often composed of multiple elements, each of which is data for 'one taxon', with the first element being a distinctive six-element vector composed of numbers, corresponding to the six variable tables by `fossilRecord2fossilTaxa` after simulating with `simFossilRecord` (originally produced by deprecated function `simFossiltaxa`).
- **merge.cryptic**: If `TRUE`, cryptic taxon-units (i.e. those in the same cryptic complex) will be merged into single taxa for the sake of being counted in the diversity curves presented by this function.
- **plotLegend**: A logical. Should a legend be plotted? Only applies if sampling processes were modeled.
- **legendPosition**: Where should the legend be plotted? See help for `legend` for details. Only applies if sampling processes were modeled.
- **curveColors**: A vector of length two indicating what colors the original and sampled diversity curves should be displayed in. Only applies if sampling processes were modeled.
- **curveLineTypes**: A vector of length two indicating what colors the original and sampled diversity curves should be displayed in. Only applies if sampling processes were modeled.

Details

This function is essentially a wrapper for `paleotree` function `multiDiv`.

Value

This function returns nothing: it just creates a plot.
Author(s)
David W. Bapst

See Also
simFossilRecord

Examples

```r
set.seed(44)
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, nruns = 1, nTotalTaxa = c(20,30), nExtant = 0, plot = FALSE)

# now let's plot it
divCurveFossilRecordSim(record)
```

Description

Functions to plot diversity curves based on taxic range data, in both discrete and continuous time, and for phylogenies.

Usage

```r
taxicDivCont(timeData, int.length = 1, int.times = NULL, plot = TRUE, plotLogRich = FALSE, timelims = NULL, drop.cryptic = FALSE)
taxicDivDisc(timeList, int.times = NULL, drop.singletons = FALSE, plot = TRUE, plotLogRich = FALSE, timelims = NULL, extant.adjust = 0.001, split.int = TRUE)
phyloDiv(tree, int.length = 0.1, int.times = NULL, plot = TRUE, plotLogRich = FALSE, drop.ZLB = TRUE, timelims = NULL)
```

Arguments

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>timeData</td>
<td>Two-column matrix giving the per-taxon first and last appearances in absolute time. The simulated data tables output by fossilRecord2fossilTaxa following simulation with simFossilRecord can also be supplied to taxicDivCont.</td>
</tr>
<tr>
<td>int.length</td>
<td>The length of intervals used to make the diversity curve. Ignored if int.times is given.</td>
</tr>
<tr>
<td>int.times</td>
<td>An optional two-column matrix of the interval start and end times for calculating the diversity curve. If NULL, calculated internally. If given, the argument split.int and int.length are ignored.</td>
</tr>
</tbody>
</table>
DiversityCurves

plot
If TRUE, a diversity curve generated from the data is plotted.

plotLogRich
If TRUE, taxic diversity is plotted on log scale.

timelims
Limits for the x (time) axis for diversity curve plots. Only affects plotting. Given as either NULL (the default) or as a vector of length two as for xlim in the basic R function plot. Time axes will be plotted exactly to these values.

drop.cryptic
If TRUE, cryptic taxa are merged to form one taxon for estimating taxon curves. Only works for objects from simFossilRecord via fossilRecord2fossilTaxa.

timelist
A list composed of two matrices, giving interval start and end dates and taxon first and last occurrences within those intervals. See details.

drop.singletons
If TRUE, taxa confined to a single interval will be dropped prior to the diversity curve calculation. This is sometimes done if single intervals have overly high diversities due to the 'monograph' effect where more named taxa are known in certain intervals largely due to taxonomic expert effort and not real changes in historical biotic diversity.

extant.adjust
Amount of time to be added to extend start time for (0,0) bins for extant taxa, so that the that 'time interval' does not appear to have an infinitely small width.

split.int
For discrete time data, should calculated/input intervals be split at discrete time interval boundaries? If FALSE, can create apparent artifacts in calculating the diversity curve. See below.

tree
A time-scaled phylogeny of class phylo.

drop.ZLB
If true, zero-length terminal branches are dropped from the input tree for phylogenetic datasets, before calculating standing diversity.

Details

First, some background. Diversity curves are plots of species/taxon/lineage richness over time for a particular group of organisms. For paleontological studies, these are generally based on per-taxon range data while more recently in evolutionary biology, molecular phylogenies have been used to calculate lineage-through-time plots (LTTs). Neither of these approaches are without their particular weaknesses; reconstructing the true history of biodiversity is a difficult task no matter what data is available.

The diversity curves produced by these functions will always measure diversity within binned time intervals (and plot them as rectangular bins). For continuous-time data or phylogenies, one could decrease the int.length used to get what is essentially an 'instantaneous' estimate of diversity. This is warned against, however, as most historical diversity data will have some time-averaging or uncertain temporal resolution and thus is probably not finely-resolved enough to calculate instantaneous estimates of diversity.

As with many functions in the paleotree library, absolute time is always decreasing, i.e. the present day is zero.

As diversity is counted within binned intervals, the true standing diversity may be somewhat lower than the measured (observed) quantity, particularly if intervals are longer than the mean duration of taxa is used. This will be an issue with all diversity curve functions, but particularly the discrete-time variant. For diversity data in particularly large discrete time intervals, plotting this data in smaller bins which do not line up completely with the original intervals will create a 'spiky' diversity curve,
as these smaller intersecting bins will have a large number of taxa which may have been present in either of the neighboring intervals. This will give these small bins an apparently high estimated standing diversity. This artifact is avoided with the default setting split.int = TRUE, which will split any input or calculated intervals so that they start and end at the boundaries of the discrete-time range bins.

The timeList object should be a list composed of two matrices, the first matrix giving by-interval start and end times (in absolute time), the second matrix giving the by-taxon first and last appearances in the intervals defined in the first matrix, numbered as the rows. Absolute time should be decreasing, while the intervals should be numbered so that the number increases with time. Taxa alive in the modern should be listed as last occurring in a time interval that begins at time 0 and ends at time 0. See the documentation for the time-scaling function bin_timePaleoPhy and the simulation function binTimeData for more information on formatting.

Unlike some paleotree functions, such as perCapitaRates, the intervals can be overlapping or of unequal length. The diversity curve functions deal with such issues by assuming taxa occur from the base of the interval they are first found in until the end of the last interval they are occur in. Taxa in wide-ranging intervals that contain many others will be treated as occurring in all nested intervals.

phyloDiv will resolve polytomies to be dichotomous nodes separated by zero-length branches prior to calculating the diversity curve. There is no option to alter this behavior, but it should not affect the use of the function because the addition of the zero-length branches should produce an identical diversity history as a polytomy. phyloDiv will also drop zero-length terminal branches, as with the function dropZLB. This the default behavior for the function but can be turned off by setting the argument drop.zlb to FALSE.

Value

These functions will invisibly return a three-column matrix, where the first two columns are interval start and end times and the third column is the number of taxa/lineages counted in that interval.

Author(s)

David W. Bapst

See Also

multiDiv, timeSliceTree, binTimeData

There are several different functions for traditional LTT plots (phylogenetic diversity curves), such as the function ltt.plot in the package ape, the function ltt in the package phytools, the function plotLtt in the package laser and the function LTT.average.root in the package TreeSim.

Examples

#taxicDivDisc with the retiolinae dataset
data(retiolitinae)
taxicDivDisc(retioRanges)

#simulation examples
set.seed(444)
DiversityCurves

```r
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1, nTotalTaxa = c(30, 40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
# let's see what the 'true' diversity curve looks like in this case
# plot the FADs and LADs with taxicDivCont()
taxicDivCont(taxa)
# simulate a fossil record with imperfect sampling with sampleRanges
rangesCont <- sampleRanges(taxa, r = 0.5)
# plot the diversity curve based on the sampled ranges
layout(1:2)
taxicDivCont(rangesCont)
# Now let's use binTimeData to bin in intervals of 1 time unit
rangesDisc <- binTimeData(rangesCont, int.length = 1)
# plot with taxicDivDisc
taxicDivDisc(rangesDisc)
# compare to the continuous time diversity curve
layout(1)
# Now let's make a tree using taxa2phylo
tree <- taxa2phylo(taxa, obs_time = rangesCont[,2])
phyloDiv(tree)

# a simple example with phyloDiv
  # using a tree from rtree in ape
set.seed(444)
tree <- rtree(100)
phyloDiv(tree)

# a neat example of using phyloDiv with timeSliceTree
  # to simulate doing molecular-phylogeny studies
  # of diversification...in the past
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1, nTotalTaxa = c(30, 40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
taxicDivCont(taxa)
# that's the whole diversity curve
# with timeSliceTree we could look at the lineage accumulation curve
  # we'd get of species sampled at a point in time
tree <- taxa2phylo(taxa)
# use timeSliceTree to make tree of relationships up until time = 950
tree950 <- timeSliceTree(tree, sliceTime = 950, plot = TRUE, drop.extinct = FALSE)
# use drop.extinct = T to only get the tree of lineages extant at time = 950
tree950 <- timeSliceTree(tree, sliceTime = 950, plot = TRUE, drop.extinct = TRUE)
# now it's an ultrametric tree with many fewer tips...
# let's plot the lineage accumulation plot on a log scale
phyloDiv(tree950, plotLogRich = TRUE)

# an example of a 'spiky' diversity curve and why split.int is a good thing
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1, nTotalTaxa = c(30, 40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
```
durationFreq

Models of Sampling and Extinction for Taxonomic Duration Datasets

Description

These functions construct likelihood models of the observed frequency of taxon durations, given either in discrete (make_durationFreqDisc) or continuous time (make_durationFreqCont). These models can then be constrained using functions available in this package and/or analyzed with commonly used optimizing functions.

Usage

make_durationFreqCont(timeData, groups = NULL, drop.extant = TRUE, threshold = 0.01, tol = 1e-04)

make_durationFreqDisc(timeList, groups = NULL, drop.extant = TRUE)

Arguments

timeData

Two-column matrix of per-taxon first and last occurrence given in continuous time, relative to the modern (i.e. older dates are also the 'larger' dates). Unsampled taxa (e.g. from a simulation of sampling in the fossil record, listed as NAs in the supplied matrix) are automatically dropped from the matrix and from groups simultaneously.

groups

Either NULL (the default) or matrix with the number of rows equal to the number of taxa and the number of columns equal to the number of 'systems' of categories for taxa. Taxonomic membership in different groups is indicated by numeric values. For example, a dataset could have a 'groups' matrix composed of a column representing thin and thick shelled taxa, coded 1 and 2 respectively, while the second column indicates whether taxa live in coastal, outer continental shelf, or deep marine settings, coded 1-3 respectively. Different combinations of groups will be treated as having independent sampling and extinction parameters in the default analysis, for example, thinly-shelled deep marine species will have separate parameters from thinly-shelled coastal species. Grouping systems
durationFreq could also represent temporal heterogeneity, for example, categorizing Paleozoic versus Mesozoic taxa. If groups are NULL (the default), all taxa are assumed to be of the same group with the same parameters. Unsampled taxa (e.g. from a simulation of sampling in the fossil record, listed as NAs in timeData or timeList) are automatically dropped from groupings and the time dataset (either timeData or timeList) and from groups simultaneously.

**drop.extant** Drops all extant taxa from a dataset, w

**threshold** The smallest allowable duration (i.e. the measured difference in the first and last occurrence dates for a given taxon). Durations below this size will be treated as "one-hit" sampling events.

**tol** Tolerance level for determining whether a taxon from a continuous-time analysis is extant or not. Taxa which occur at a date less than tol are treated as occurring at the modern day (i.e. being functionally identical as occurring at 0 time).

**timeList** A 2 column matrix with the first and last occurrences of taxa given in relative time intervals (i.e. ordered from first to last). If a list of length two is given for timeData, such as would be expected if the output of binTimeData was directly input, the second element is used. See details. Unsampled taxa (e.g. from a simulation of sampling in the fossil record, listed as NAs in the second matrix) are automatically dropped from the timeList and from groups simultaneously. Living taxa observed in the modern day are expected to be listed as last observed in a special interval (0,0), i.e. begins and ends at 0 time. This interval is always automatically removed prior to the calculation intermediary data for fitting likelihood functions.

**Details**

These functions effectively replace two older functions in paleotree, now removed, getSampRateCont and getSampProbDisc. The functions here do not offer the floating time interval options of their older siblings, but do allow for greater flexibility in defining constrains on parameter values. Differences in time intervals, or any other conceivable discrete differences in parameters, can be modeled using the generic groups argument in these functions.

These functions use likelihood functions presented by Foote (1997). These analyses are ideally applied to data from single stratigraphic section but potentially are applicable to regional or global datasets, although the behavior of those datasets is less well understood.

As with many functions in the paleotree library, absolute time is always decreasing, i.e. the present day is zero and older dates are 'larger'. On the contrary, relative time is in intervals with non-zero integers that increase sequentially beginning with 1, from earliest to oldest.

For make_durationFreqDisc, the intervals in timeList should be non-overlapping sequential intervals of roughly equal length. These should be in relative time as described above, so the earliest interval should be 1 and the numbering should increase as the intervals go up with age. If both previous statements are true, then differences in interval numbers will represent the same rough difference in the absolute timing of those intervals. For example, a dataset where all taxa are listed from a set of sequential intervals of similar length, such as North American Mammal assemblage zones, microfossil faunal zones or graptolite biozones can be given as long as they are correctly numbered in sequential order in the input. As a counter example, a dataset which includes taxa resolved only to intervals as wide as the whole Jurassic and taxa resolved to biozones within the Jurassic should not be included in the same input. Drop taxa from less poorly resolved intervals...
from such datasets if you want to apply this function, as long as this retains a large enough sample of taxa listed from the sequential set of intervals.

Please check that the optimizer function you select actually converges. The likelihood surface can be very flat in some cases, particularly for small datasets (<100 taxa). If the optimizer does not converge, consider increasing iterations or changing the starting values.

**Value**

A function of class "paleotreeFunc", which takes a vector equal to the number of parameters and returns the *negative* log likelihood (for use with optim and similar optimizing functions, which attempt to minimize support values). See the functions listed at modelMethods for manipulating and examining such functions and constrainParPaleo for constraining parameters.

Parameters in the output functions are named 'q' for the instantaneous per-capita extinction rate, 'r' for the instantaneous per-capita sampling rate and 'R' for the per-interval taxonomic sampling probability. Groupings follow the parameter names, separated by periods; by default, the parameters will be placed in at least group 1 (of a grouping with a single group), such that make_durationFreqCont by default creates a function with parameters named 'q.1' and 'r.1', while make_durationFreqDisc creates a function with parameters named 'q.1' and 'R.1'.

Note that the 'q' parameters estimated by make_durationFreqDisc is scaled to per lineage intervals and not to per lineage time-units. If intervals are the same length, this can be easily corrected by multiplying 1 by the interval length. It is unclear how to treat uneven intervals and I urge workers to consider multiple strategies.

For translating these sampling probabilities and sampling rates, see SamplingConv.

**Author(s)**

David W. Bapst

**References**


**See Also**

See freqRat, sRate2sProb, qsRate2Comp sProb2sRate and qsProb2Comp.

**Examples**

```r
# let's simulate some taxon ranges from an imperfectly sampled fossil record
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(30, 40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
rangesCont <- sampleRanges(taxa, r = 0.5)
# bin the ranges into discrete time intervals
rangesDisc <- binTimeData(rangesCont, int.length = 1)
```
#note that we made interval lengths = 1:
# thus q (per int) = q (per time) for make_durationFreqDisc

## Not run:
# old ways of doing it (defunct as of paleotree version 2.6)
getSampRateCont(rangesCont)
getSampProbDisc(rangesDisc)

## End(Not run)

# new ways of doing it
  # we can constrain our functions
  # we can use parInit, parLower and parUpper to control parameter bounds

# as opposed to getSampRateCont, we can do:
likFun <- make_durationFreqCont(rangesCont)
optim(parInit(likFun),likFun,lower = parLower(likFun),upper = parUpper(likFun),
  method = "L-BFGS-B",control = list(maxit = 1000000))

# as opposed to getSampProbDisc, we can do:
likFun <- make_durationFreqDisc(rangesDisc)
optim(parInit(likFun),likFun,lower = parLower(likFun),upper = parUpper(likFun),
  method = "L-BFGS-B",control = list(maxit = 1000000))

# these give the same answers (as we'd expect them to!)

# with newer functions we can constrain our functions easily
  # what if we knew the extinction rate = 0.1 a priori?
likFun <- make_durationFreqCont(rangesCont)
likFun <- constrainParPaleo(likFun,q.1~0.1)
optim(parInit(likFun),likFun,lower = parLower(likFun),upper = parUpper(likFun),
  method = "L-BFGS-B",control = list(maxit = 1000000))

# actually decreases our sampling rate estimate
  # gets further away from true rate estimate, r = 0.5 (geesh!)
  # but this *is* a small dataset...

---

equation2function  

**Turn a Character String of the Right-Hand Side of an Equation into an R Function**

description

equation2function converts the right-hand side of an equation that can be written as a single line (like the right-hand side of an object of class formula) and creates an R function which calls the variables within as arguments and returns values consistent with the parameters of the input equation as written.

usage

equation2function(equation, envir = parent.frame(), notName = "XXXXXXXXXX")
Arguments

- **equation**: The right-hand-side (RHS) of an equation, given as a character string. If not of type character, `equation2function` attempts to coerce `equation` to type character and fails if it cannot.

- **envir**: The environment the resulting function will be evaluated in. See `as.function`.

- **notName**: A useless string used simply as a placeholder in turning `equation` into a function, which should not match any actual variable in `equation`. Only supplied as an argument in case any

Details

This simple little function is rather 'hacky' but seems to get the job done, for a functionality that does not seem to be otherwise exist elsewhere in R.

Value

A function, with named blank (i.e. no default value) arguments.

Author(s)

David W. Bapst

Examples

```r
# some simple examples
foo <- equation2function("x+y")
foo
foo(x = 4, y = 0.1)

foo <- equation2function("x+2*sqrt(2*y+3)\cdot 2")
foo
foo(x = 4, y = 0.1)

# what about weird long argument names and spaces
foo <- equation2function("stegosaur + 0.4 \times P")
foo
foo(stegosaur = 5, P = 0.3)
```

Description

The following functions are for measuring and fitting various distributions to the gradual exhaustion of unexpressed character states, as originally described by Wagner (2000, Evolution).
Usage

accioExhaustionCurve(phylotree, charData, charTypes = "unordered", 
                      outgroup = NULL, firstAppearances = NULL, missingCharValue = "?",
                      inapplicableValue = "-")

accioBestAcquisitionModel(exhaustion_info, changesType, 
                          models = c("exponential", "gamma", "lognormal", "zipf"))

charExhaustPlot(exhaustion_info, changesType, xlab = "Total Characters", 
                 ylab = NULL, main = NULL, xsize = 3)

Arguments

phyloTree A phylogenetic tree of class phylo as used by package ape.
charData A data.frame of morphological character codings (a morphological 'matrix'),
           with taxon units as rows and characters as columns.
charTypes A vector of length equal to the number of characters in charData, with elements
           indicating whether the corresponding character in charData is "unordered" or
           "ordered". However, if length(charTypes) = 1, then it is repeated for all
taxa. The default value for this argument is "unordered".
outgroup A string matching to one of the tip labels as given by tip.label,
firstAppearances A vector, with length equal to the same number of taxa (rows) as in charData,
in the same corresponding order.
missingCharValue The string value indicating a missing character coding value, by default "?".
inapplicableValue The string value indicating an inapplicable character coding value, by default 
                  "-".
exhaustion_info The list of results output from function accioExhaustionCurve.
changesType A single character value, indicating the character change data to be assessed
             from the result of the character exhaustion analysis, must be one of either 'total-
             lAcc' (to the total number of accumulated character changes, ideal for modeling
             the size and distribution of state space) or 'charAlt' (to plot the total number of
             character alterations, ideal for modeling the size and distribution of character
             space).
models A vector of type character naming models to be fit. Default is c("exponential", "gamma", "lognormal"

xlab Label for the X axis; "Total Characters" by default.
ylab Label for the Y axis. If not provided by the user, a label based on the changesType
       argument will be used.
main Main title label for the plot. If not provided by the user, a label based on the
       changesType argument will be used.
xsize Parameter controlling size of the axes, which are forced to be symmetric.
Details

acciEXhaustionCurve uses a Sankoff parsimony ancestral-reconstruction algorithm (written by PJ Wagner, not the one from phangorn used elsewhere in paleotree) to calculate character changes across each branch (internode edge) of a tree, and then reports the counts of character state.

acciBestAcquisitionModel takes output from accioExhaustionCurve, calculates one of two character change indices, and then fits a series of user-selected models to the dataset, returning details pertaining to the best-fit model.

charExhaustPlot is a wrapper for accioBestAcquisitionModel that produces a plot of the observed character change data against the expectation under the best-fit model.

The functions accioBestAcquisitionModel and charExhaustPlot offer users two different options for examining character change: totalAcc fits models to the total accumulated number of state changes over the phylogeny, thus using exhaustion to explore the size and distribution of character space. The other option charAlt fits models to the number of character that alter from primitive to derived over phylogeny, thus reflecting the size and distribution of state space.

acciExhaustionCurve can order its reconstruction of change by stratigraphic order of first appearances. It is unclear what this means.

Value

acciExhaustionCurve outputs a list containing two objects: first, a matrix named exhaustion consisting of three columns: "Steps", "Novel_States", and "Novel_Characters", respectively giving the counts of these respective values for each branch (internode edge). The second element of this list is named State_Derivations and is a count of how often each state, across all characters, was derived relative to the primitive position along each internode edge.

The output of accioBestAcquisitionModel is a list object containing information on the best-fit model, the parameters of that model, the calculated probabilition distribution for that model at the same intervals, for use in quantile plots.

charExhaustPlot produces a plot, and outputs no data.

Note

This family of functions presented here were originally written by Peter J. Wagner, and then modified and adapted by David W. Bapst for wider release in a CRAN-distributed package: paleotree. This makes the code presented here a very different beast than typical paleotree code, for example, there are fewer tests of correct input type, length, etc.

Author(s)

Initially written by Peter J. Wagner, with modification and documentation by David W. Bapst.

References

See Also

Also see paleotree functions `minCharChange` and `ancPropStateMat`, the latter of which is a wrapper for phangorn’s function `ancestral.pars`.

Examples

```r
# get data
data(SongZhangDicrano)
dicranoTree <- cladogramDicranoX13

# modify char data
charMat <- data.matrix(charMatDicrano)
charMat[is.na(charMatDicrano)] <- 0
charMat <- (charMat-1)
colnames(charMat) <- NULL
# replace missing values
charMat[is.na(charMatDicrano)] <- "??"

# the 'outgroup' is Exigraptus, first taxon listed in the matrix
exhaustionResults <- accioExhaustionCurve(phyloTree = dicranoTree,
                                           charData = charMat, charTypes = "unordered",
                                           outgroup = "Exigraptus_uniformis")

# fits models to exhaustion for total accumulation
accioBestAcquisitionModel(exhaustion_info = exhaustionResults,
                          changesType = "totalAcc",
                          models = c("exponential","gamma","lognormal","zipf"))

# plot of exhausation of total accumulation of character states
charExhaustPlot(exhaustion_info = exhaustionResults,
                changesType = "totalAcc")

# plot of exhausation of character alterations
charExhaustPlot(exhaustion_info = exhaustionResults,
                changesType = "charAlt")
```

Description

This function takes a tree composed of higher-level taxa and a vector of lower-level taxa belonging to the set of higher-level taxa included in the input tree and produces a tree composed of the lower-level taxa, by treating the higher-level taxa as unresolved monophyletic polytomies. A user can
also mark higher taxa as paraphyletic such that these are secondarily collapsed and do not form monophyletic clades in the output tree.

Usage

```
expandTaxonTree(taxonTree, taxaData, collapse = NULL, keepBrLen = FALSE,
    plot = FALSE)
```

Arguments

- **taxonTree**: A phylo object where tips represent higher taxa
- **taxaData**: Character vector of higher taxa, with elements names equal to the lower taxa. See below.
- **collapse**: Character vector of non-monophyletic higher taxa to be collapsed
- **keepBrLen**: Logical, decides if branch lengths should be kept or discarded. FALSE by default. See details below.
- **plot**: If true, plots a comparison between input and output trees

Details

The output tree will probably be a rough unresolved view of the relationships among the taxa, due to the treatment of higher-level taxa as polytomies. This is similar to the methods used in Webb and Donoghue (2005) and Friedman (2009). Any analyses should be done by resolving this tree with `multi2di` in the ape package or via the various time-scaling functions found in this package (paleotree).

The taxaData vector should have one element per lower-level taxon that is to be added to the tree. The name of each element in the vector should be the names of the lower-level taxa, which will be used as new tip labels of the output lower-taxon tree. There should be no empty elements! `expandTaxonTree` won’t know what to do with taxa that don’t go anywhere.

By default, all higher-level taxa are treated as monophyletic clades if not otherwise specified. The collapse vector can (and probably should) be used if there is doubt about the monophyly of any higher-level taxa included in the input taxon-tree, so that such a group would be treated as a paraphyletic group in the output tree.

Also by default, the output tree will lack branch lengths and thus will not be time-scaled. If keepBrLen is true, then the tree’s edge lengths are kept and new taxa are added as zero length branches attaching to a node that represents the previous higher-taxon. This tree is probably not useful for most applications, and may even strongly bias some analyses. USE WITH CAUTION! The ‘collapse’ vector will cause such edges to be replaced by zero-length branches rather than fully collapsing them, which could have odd effects. If ‘collapse’ is not null and keepBrLen is true, a warning is issued that the output probably won’t make much sense at all.

Value

Outputs the modified tree as an object of class phylo, with the higher-level taxa expanded into polytomies and the lower-level taxa as the tip labels.
footeValues

Author(s)

David W. Bapst

References


See Also

multi2di, bind.tree

Examples

```r
set.seed(444)
# lets make our hypothetical simulated tree of higher taxa
taxr <- rtree(10)
taxd <- sample(taxr$tip.label,30,replace = TRUE) # taxa to place within higher taxa
names(taxd) <- paste(taxd,"_x",1:30,sep = "")
coll <- sample(taxr$tip.label,3) # what to collapse?
expandTaxonTree(taxonTree = taxr,taxaData = taxd,collapse = coll,plot = TRUE)
```

footeValues

*Calculates Values for Foote’s Inverse Survivorship Analyses*

Description

This function calculates the intermediary values needed for fitting Foote’s inverse survivorship analyses, as listed in the table of equations in Foote (2003), with the analyses themselves described further in Foote (2001) and Foote (2005).

Usage

```r
footeValues(p, q, r, PA_n = 0, PB_1 = 0, p_cont = TRUE, q_cont = TRUE, Nb = 1)
```

Arguments

- `p`  
  Instantaneous origination/branching rate of taxa. Under a continuous model, assumed to be *per interval*, or equal to the product of interval lengths and the rates per lineage time units for each interval. Under a pulsed mode (p_cont = FALSE), p is a per-interval 'rate' which can exceed 1 (because diversity can more than double; Foote, 2003a). Given as a vector with length equal to the number of intervals, so a different value may be given for each separate interval. Must be the same length as q and r.
q

Instantaneous extinction rate of taxa. Under a continuous model, assumed to be per interval, or equal to the product of interval lengths and the rates per lineage time units for each interval. Under a pulsed mode (q_cont = FALSE), q is a per-interval ‘rate’ but which cannot be observed to exceed 1 (because you can’t have more taxa go extinct than exist). Given as a vector with length equal to the number of intervals, so a different value may be given for each separate interval. Must be the same length as p and r.

r

Instantaneous sampling rate of taxa, assumed to be per interval, or equal to the product of interval lengths and the rates per lineage time units for each interval. Given as a vector with length equal to the number of intervals, so a different value may be given for each separate interval. Must be the same length as p and q.

PA_n

The probability of sampling a taxon after the last interval included in a survivorship study. Usually zero for extinct groups, although more logically has the value of 1 when there are still extant taxa (i.e., if the last interval is the Holocene and the group is still alive, the probability of sampling them later is probably 1...). Should be a value of 0 to 1.

PB_1

The probability of sampling a taxon before the first interval included in a survivorship study. Should be a value of 0 to 1.

p_cont

If TRUE (the default), then origination is assumed to be a continuous time process with an instantaneous rate. If FALSE, the origination is treated as a pulsed discrete-time process with a probability.

q_cont

If TRUE (the default), then extinction is assumed to be a continuous time process with an instantaneous rate. If FALSE, the extinction is treated as a pulsed discrete-time process with a probability.

Nb

The number of taxa that enter an interval (b is for ‘bottom’). This is an arbitrary constant used to scale other values in these calculations and can be safely set to 1.

Details

Although most calculations in this function agree with the errata for Foote’s 2003 table (see references), there were some additional corrections for Prob(D|FL) made as part of a personal communication in 2013 between the package author and Michael Foote.

Value

Returns a matrix with number of rows equal to the number of intervals (i.e. the length of p, q and r) and named columns representing the different values calculated by the function: ”Nb”, ”Nbt”, ”NbL”, ”NFL”, ”PD_bt”, ”PD_bL”, ”PD_Ft”, ”PD_FL”, ”PA”, ”PB”, ”Xbt”, ”XbL”, ”XFl” and ”XFL”.

Author(s)

David W. Bapst, with advice from Michael Foote.
References


Examples

```r
#very simple example with three intervals, same value for all parameters

data <- matrix(c(0, 100, 200, 300, 400), ncol = 2, byrow = TRUE)

#example rates (for the most part)
rate <- rep(0.1, 3)
#all continuous
footeValues(rate, rate, rate)
#origination pulsed
footeValues(rate, rate, rate, p_cont = FALSE)
#extinction pulsed
footeValues(rate, rate, rate, q_cont = FALSE)
#all pulsed
footeValues(rate, rate, rate, p_cont = FALSE, q_cont = FALSE)
```

freqRat

*Frequency Ratio Method for Estimating Sampling Probability*

Description

Estimate per-interval sampling probability in the fossil record from a set of discrete-interval taxon ranges using the frequency-ratio method described by Foote and Raup (1996). Can also calculate extinction rate per interval from the same data distribution.

Usage

```
freqRat(timeData, calcExtinction = FALSE, plot = FALSE)
```

Arguments

- `timeData`: A 2 column matrix with the first and last occurrences of taxa given in relative time intervals. If a list of length two is given for `timeData`, such as would be expected if the output of `binTimeData` was directly input, the second element is used.
- `calcExtinction`: If TRUE, the per-interval, per-lineage extinction rate is estimated as the negative slope of the log frequencies, ignoring single hits (as described in Foote and Raup, 1996.)
- `plot`: If true, the histogram of observed taxon ranges is plotted, with frequencies on a linear scale
Details

This function uses the frequency ratio ("freqRat") method of Foote and Raup (1996) to estimate the per-interval sampling rate for a set of taxa. This method assumes that intervals are of fairly similar length and that taxonomic extinction and sampling works similar to homogenous Poisson processes. These analyses are ideally applied to data from single stratigraphic section but potentially are applicable to regional or global datasets, although the behavior of those datasets is less well understood.

The frequency ratio is a simple relationship between the number of taxa observed only in a single time interval (also known as singletons), the number of taxa observed only in two time intervals and the number of taxa observed in three time intervals. These respective frequencies, respectively f1, f2 and f3 can then be related to the per-interval sampling probability with the following expression.

\[ \text{Sampling Probability} = \frac{f_2^2}{f_1 \times f_3} \]

This frequency ratio is generally referred to as the 'freqRat' in paleobiological literature. It is relatively easy to visually test if range data fits expectation that true taxon durations are exponentially distributed by plotting the frequencies of the observed ranges on a log scale: data beyond the 'singletons' category should have a linear slope, implying that durations were originally exponentially distributed. (Note, a linear scale is used for the plotting diagram made by this function when 'plot' is TRUE, so that plot cannot be used for this purpose.)

The accuracy of this method tends to be poor at small interval length and even relatively large sample sizes. A portion at the bottom of the examples in the help file examine this issue in greater detail with simulations. This package author recommends using the ML method developed in Foote (1997) instead, which is usable via the function \texttt{make\_durationFreqDisc}.

As extant taxa should not be included in a freqRat calculation, any taxa listed as being in a bin with start time 0 and end time 0 (and thus being extant without question) are dropped before the model fitting it performed.

Value

This function returns the per-interval sampling probability as the "freqRat", and estimates

Author(s)

David W. Bapst

References


See Also

Model fitting methods in \texttt{make\_durationFreqDisc} and \texttt{make\_durationFreqCont}. Also see conversion methods in \texttt{sProb2sRate}, \texttt{qsProb2Comp}
Examples

```r
# Simulate some fossil ranges with simFossilRecord
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(30,40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
# Simulate a fossil record with imperfect sampling with sampleRanges
rangesCont <- sampleRanges(taxa, r = 0.1)
# Now let's use binTimeData to bin in intervals of 5 time units
rangesDisc <- binTimeData(rangesCont,int.length = 5)

# Now, get an estimate of the sampling rate (we set it to 0.5 above)
# For discrete data we can estimate the sampling probability per interval (R)
# i.e. this is not the same thing as the instantaneous sampling rate (r)
# can use sRate2sProb to see what we would expect
sRate2sProb(r = 0.1,int.length = 5)
# Expect R = -0.39

# Now we can apply freqRat to get sampling probability
SampProb <- freqRat(rangesDisc,plot = TRUE)
SampProb

# Est. R = -0.25
# Not wildly accurate, is it?

# can also calculate extinction rate per interval of time
freqRat(rangesDisc,calcExinction = TRUE)

# Est. ext rate = -0.44 per interval
# 5 time-unit intervals, so -0.44 / 5 = -0.08 per time-unit
# That's pretty close to the generating value of 0.01, used in sampleRanges

## Not run:

# The following example code (which is not run by default) examines how
# the freqRat estimates vary with sample size, interval length
# and compare it to using make_durationFreqDisc

# how good is the freqRat at 20 sampled taxa on avg?
set.seed(444)
r <- runif(100)
int.length = 1
R <- sapply(r,sRate2sProb,int.length = 1) # estimate R from r, assuming stuff like p = q
ntaxa <- freqRats <- numeric()
for(i in 1:length(r)){
  # assuming budding model
  record <- simFossilRecord(p = 0.1, q = 0.1, r = r[i], nruns = 1,
nSamp = c(15,25), nExtant = 0, plot = TRUE)
ranges <- fossilRecord2FossilRanges(record)
timeList <- binTimeData(ranges,int.length = int.length)
ntaxa[i] <- nrow(timeList[[2]])
}
```

```
freqRats[i] <- freqRat(timeList)
}
plot(R,freqRats):abline(0,1)
#without the gigantic artifacts bigger than 1...
plot(R,freqRats,ylim = c(0,1));abline(0,1)
#very worrisome lookin'!

#how good is it at 100 sampled taxa on average?
set.seed(444)
r <- runif(100)
int.length = 1
R <- sapply(r,sRate2sProb,int.length = 1)
ntaxa <- freqRats <- numeric()
for(i in 1:length(r)){
  #assuming budding model
  record <- simFossilRecord(p = 0.1, q = 0.1, r = r[i], nruns = 1,
  nSamp = c(80,150), nExtant = 0, plot = TRUE)
  ranges <- fossilRecord2FossilRanges(record)
  timeList <- binTimeData(ranges,int.length = int.length)
  ntaxa[i] <- nrow(timeList[[2]])
  freqRats[i] <- freqRat(timeList)
}
plot(R,freqRats,ylim = c(0,1));abline(0,1)
#not so hot, eh?

###################
#LETS CHANGE THE TIME BIN LENGTH!

#how good is it at 100 sampled taxa on average, with longer time bins?
set.seed(444)
r <- runif(100)
int.length <- 10
R <- sapply(r,sRate2sProb,int.length = int.length)
ntaxa <- freqRats <- numeric()
for(i in 1:length(r)){
  #assuming budding model
  record <- simFossilRecord(p = 0.1, q = 0.1, r = r[i], nruns = 1,
  nSamp = c(80,150), nExtant = 0, plot = TRUE)
  ranges <- fossilRecord2FossilRanges(record)
  timeList <- binTimeData(ranges,int.length = int.length)
  ntaxa[i] <- nrow(timeList[[2]])
  freqRats[i] <- freqRat(timeList)
}
plot(R,freqRats,ylim = c(0,1));abline(0,1)
#things get more accurate as interval length increases... odd, eh?

#how good is it at 20 sampled taxa on average, with longer time bins?
set.seed(444)
r <- runif(100)
int.length <- 10
R <- sapply(r,sRate2sProb,int.length = int.length)
ntaxa <- freqRats <- numeric()
for(i in 1:length(r)){
#assuming budding model
record <- simFossilRecord(p = 0.1, q = 0.1, r = r[i], nruns = 1, nSamp = c(15, 25), nExtant = 0, plot = TRUE)
ranges <- fossilRecord2fossilRanges(record)
timelist <- binTimeData(ranges,int.length = int.length)
nextant[i] <- nrow(timelist[[2]])
plot(R,freqRats[i],ylim = c(0,1));abline(0,1)
#still not so hot at low sample sizes, even with longer bins

#how good is the ML method at 20 taxa, 1 time-unit bins?
set.seed(444)
r <- runif(100)
int.length <- 1
R <- sapply(r,sRate2sProb,int.length = int.length)
ntaxa <- ML_sampprob <- numeric()
for(i in 1:length(r)){
  #assuming budding model
  record <- simFossilRecord(p = 0.1, q = 0.1, r = r[i], nruns = 1, nSamp = c(15,25), nExtant = 0, plot = TRUE)
ranges <- fossilRecord2fossilRanges(record)
timelist <- binTimeData(ranges,int.length = int.length)
nextant[i] <- nrow(timelist[[2]])
likFun <- make_durationFreqDisc(timelist)
ML_sampprob[i] <- optim(parInit(likFun),likFun,
  lower = parLower(likFun),upper = parUpper(likFun),
  method = "L-BFGS-B",control = list(maxit = 1000000))[[1]][2]
}
plot(R,ML_sampprob);abline(0,1)
# Not so great due to likelihood surface ridges
# but it returns values between 0-1

#how good is the ML method at 100 taxa, 1 time-unit bins?
set.seed(444)
r <- runif(100)
int.length <- 1
R <- sapply(r,sRate2sProb,int.length = int.length)
ntaxa <- ML_sampprob <- numeric()
for(i in 1:length(r)){
  #assuming budding model
  record <- simFossilRecord(p = 0.1, q = 0.1, r = r[i], nruns = 1, nSamp = c(80,150), nExtant = 0, plot = TRUE)
ranges <- fossilRecord2fossilRanges(record)
timelist <- binTimeData(ranges,int.length = int.length)
nextant[i] <- nrow(timelist[[2]])
likFun <- make_durationFreqDisc(timelist)
ML_sampprob[i] <- optim(parInit(likFun),likFun,
  lower = parLower(likFun),upper = parUpper(likFun),
  method = "L-BFGS-B",control = list(maxit = 1000000))[[1]][2]
graptDisparity

*Morphological Character and Range Data for late Ordovician and Early Silurian Graptoloidea*

**Description**

This dataset contains a morphological character matrix (containing both a set of 45 discrete characters, and 4 continuous characters coded as minimum and maximum range values), along with biostratigraphic range data for 183 graptoloid species-level taxa from Bapst et al. (2012, PNAS). Also includes a pre-calculated distance matrix based on the character matrix, using the algorithm applied by Bapst et al (2012). Interval dates for biostratigraphic zones is taken from Sadler et al. 2011.

**Format**

This dataset is composed of three objects:

- **graptCharMatrix** A matrix composed of mixed character data and a group code for 183 graptoloid taxa, with rows named with species names and columns named with character names.

- **graptRanges** A list containing two matrices: the first matrix describes the first and last interval times for 20 graptolite biozones and the second matrix contains the first and last appearances of 183 graptolite species in those same biozones. (In other words, graptRanges has the ‘timeList’ format called by some paleotree functions).

- **graptDistMat** A 183x183 matrix of pair-wise distances (dissimilarities) for the 183 graptolite species, using the algorithm for discrete characters and min-max range values described in Bapst et al.

**Details**

The character matrix contains characters of two differing types with a (very) small but non-negligible amount of missing character data for some taxa. This required the use of an unconventional ad hoc distance metric for the published analysis, resulting in a (very slightly) non-Euclidean distance matrix. This breaks some assumptions of some statistical analyses or requires special corrections, such as with PCO.

Note that taxonomic data were collected only for species present within an interval defined by the base of the Uncinatus biozone (~448.57 Ma) to the end of the cyphus biozone (~439.37 Ma). Many taxa have first and last appearance dates listed in graptRanges which are outside of this interval (see examples).
Source

Source for stratigraphic ranges and character data:

Source for interval dates for graptolite zones:

See Also

For more example graptolite datasets, see *retiolitinae*

This data was added mainly to provide an example dataset for *nearestNeighborDist*

Examples

```r
# load data
data(graptDisparity)

# separate out two components of character matrix

# 45 discrete characters
discChar <- graptCharMatrix[,1:45]

# min ranges for 4 continuous characters
cMinChar <- graptCharMatrix[,c(46,48,50,52)]

# max ranges for 4 continuous characters
cMaxChar <- graptCharMatrix[,c(47,49,51,53)]

# group (clade/parACLade) coding
groupID <- graptCharMatrix[,54]

# number of species
nspec <- nrow(graptCharMatrix)

# some plotting information from Bapst et al.'s plotting scripts
grpLabel <- c("Normalo.", "Monogr.", "Climaco.", "Dicrano.", "Lasiogr.", "Diplogr.", "Reticol.")
grpColor <- c("red", "purple", colors()[257], colors()[614],
              colors()[124], "blue", colors()[556])

###########

# plot diversity curve of taxa
taxicDivDisc(graptRanges)

# but the actual study interval for the data is much smaller
abline(v = 448.57, lwd = 3) # start of study interval
abline(v = 439.37, lwd = 3) # end of study interval
```
#plot diversity curve just for study interval
taxicDivDisc(graptRanges, timelims = c(448.57, 439.37))

##############

#distance matrix is given as graptDistMat
#to calculate yourself, see code below in DoNotRun section

#now, is the diagonal zero? (it should be)
all(diag(graptDistMat) == 0)

#now, is the matrix symmetric? (it should be)
isSymmetric(graptDistMat)

#can apply cluster analysis
clustRes <- hclust(as.dist(graptDistMat))
plot(clustRes, labels = FALSE)

#use ape to plot with colors at the tips
dev.new(width = 15)  # for a prettier plot
plot.phylo(as.phylo(clustRes), show.tip.label = FALSE,
no.margin = TRUE, direction = "upwards")
tiplabels(pch = 16, col = grpColor[groupID+1])
legend("bottomright", legend = grpLabel, col = grpColor, pch = 16)
dev.set(2)

#can apply PCO (use lingoes correction to account for negative values
#resulting from non-euclidean matrix
pco_res <- pcoa(graptDistMat, correction = "lingoes")

#relative corrected eigenvalues
rel_corr_eig <- pco_res$values$Rel_corr_eig
layout(1:2)
plot(rel_corr_eig)
#cumulative
plot(cumsum(rel_corr_eig))

#first few axes account for very little variance!!

#well let’s look at those PCO axes anyway
layout(1)
pco_axes <- pco_res$ vectors
plot(pco_axes[,1], pco_axes[,2], pch = 16, col = grpColor[groupID+1],
xlab = paste("PCO Axis 1, Rel. Corr. Eigenvalue = ", round(rel_corr_eig[1],3)),
ylab = paste("PCO Axis 2, Rel. Corr. Eigenvalue = ", round(rel_corr_eig[2],3)))
legend("bottomright", legend = grpLabel, col = grpColor, pch = 16, ncol = 2, cex = 0.8)
# Not run:

calculate a distance matrix (very slow!)
Bapst et al. calculated as # char diffs / total # of chars
but both calculated for only non-missing characters for both taxa
non-identical discrete states = difference for discrete traits
non-overlapping ranges for continuous characters = difference for cont traits

```
distMat <- matrix(,n-spec,n-spec)
rownames(distMat) <- colnames(distMat) <- rownames(graptCharMatrix)
for(i in 1:n-spec){ for(j in 1:n-spec){
    #calculate for each pair of species
    #discrete characters
    di <- discChar[i,]   #discrete character vector for species i
dj <- discChar[j,]   #discrete character vector for species j
    #now calculate pair-wise differences for non-missing characters
discDiff <- (di != dj)[!is.na(di)&!is.na(dj)] #logical vector
    #continuous characters: need another for() loop
    contDiff <- numeric()
    for(ct in 1:4){
        #if they do not overlap, a min must be greater than a max value
    }
    #remove NAs
    contDiff <- contDiff[!is.na(contDiff)]
    #combine
    totalDiff <- c(discDiff,contDiff)
    #divide total difference
    distMat[i,j] <- sum(totalDiff)/length(totalDiff)
}}

#but is it identical to the distance matrix already provided?
identical(distMat,graptDistMat)
#ehh, numerical rounding issues...

#A somewhat speeder alternative to calculate a distance matrix
distMat <- matrix(,n-spec,n-spec)
rownames(distMat) <- colnames(distMat) <- rownames(graptCharMatrix)
for(i in 1:(n-spec-1)){ for(j in (i+1):n-spec){
    #calculate for each pair of species
    #now calculate pair-wise differences for non-missing characters
    discDiff <- (discChar[i,] != discChar[j,])[!is.na(discChar[i,])&!is.na(discChar[j,])] #logical vector
    #continuous characters: if they do not overlap, a min must be greater than a max value
    contDiff <- sapply(1:4,function(ct)
        cMinChar[i,ct]>cMaxChar[j,ct] | cMinChar[j,ct]>cMaxChar[i,ct])
    #remove NAs, combine, divide total difference
    distMat[i,j] <- distMat[j,i] <- sum(c(discDiff,contDiff[!is.na(contDiff)]))/length(c(discDiff,contDiff[!is.na(contDiff)]))
}}
diag(distMat) <- 0

#but is it identical to the distance matrix already provided?
identical(distMat,graptDistMat)
Example datasets consisting of (a) occurrence data and (b) taxonomic data downloaded from the Paleobiology Database API for the Graptolithina.

The example occurrence dataset (grapt0ccPBDB) is a data.frame consisting of 5900 occurrences (rows) and 35 variables (columns). The example taxonomy dataset (graptTaxaPBDB) is a data.frame consisting of 364 formal taxa (rows) and 53 variables (columns). Variables are coded in the ‘pbdb’ vocabulary of the PBDB API v1.2.

This example PBDB data is included here for testing functions involving occurrence data and taxonomy in paleotree.

Source

See examples for the full R code used to obtain the data from the API. You can find the Paleobiology Database at [http://paleobiodb.org](http://paleobiodb.org)

The occurrence data was entered by (in order of relative portion) P. Novack-Gottshall, M. Krause, M. Foote, A. Hendy, T. Hanson, M. Sommers and others. This same data was authorized mainly by A. Miller, W. Kiessling, M. Foote, A. Hendy, S. Holland, J. Sepkoski and others.

See Also

-taxonSortPBDBocc, occData2timeList, makePBDBtaxonTree, plotOccData

Examples

```r
# original code used to obtain this dataset on March 21st, 2015
# using version 1.2 of the Paleobiology Database API

# (sorry, URLs removed as they lead to the PBDB test server...)
```
horizonSampRate

Estimate Sampling Rate from Sampling Horizon Data (Solow and Smith, 1997)

Description

This function implements the exact maximum likelihood estimator for the instantaneous sampling rate from Solow and Smith (1997, Paleobiology), which is based on the relationship between the number of collections for a set of taxa and their durations (known precisely in continuous time).

Usage

horizonSampRate(sampOcc = NULL, durations = NULL, nCollections = NULL)

Arguments

sampOcc  A list with the number of elements equal to the number of taxa, and each element of the list being a numerical vector with the length equal to the number of collections for each taxon, and each value equal to the precise date of that fossil's time of collection. These dates do not need to be ordered. If not supplied, the elements durations and nCollections must be supplied.

durations  A vector of precise durations in continuous time, with the length equal to the number of taxa. If not supplied, this is calculated from SampOcc, which must be supplied.
nCollections  A vector of integers representing the number of collections for each taxon in the input durations. If not supplied this is calculated from SampOcc, which must be supplied.

Details

Given a dataset of taxa with a vector \( N \), representing the number of collections for each taxon, and a vector \( D \), giving the precise duration for each taxon, we can use the following maximum likelihood estimator from Solow and Smith (1997) to obtain the instantaneous sampling rate:

\[
samplingRate = \frac{\text{sum}(N - 1)^2}{\text{sum}(D) \times \text{sum}(N)}
\]

This method is exclusively for datasets with very precisely dated horizons, such as microfossils from deep sea cores with very precise age models. The first and last appearance must be known very precisely to provide an equally precise estimate of the duration. Most datasets are not precise enough for this method, due to chronostratigraphic uncertainty. However, note that the age of individual collections other than the first and last appearance dates do not need to be known: its only the number of collections that matters.

Value

Returns the instantaneous sampling (in per lineage*time-units) as a single numerical value. Note that this is the instantaneous sampling rate and not the probability of sampling a taxon per interval.

References


See Also

Duration frequency methods (Foote and Raup, 1996; Foote, 1997) use ranges alone to estimate sampling parameters, implemented in *durationFreq*.

Also see the conversion functions for sampling parameters at *SamplingConv*.

Examples

```r
# can simulate this type of data with sampleRanges
# just set ranges.only = FALSE
# let's try a simulation example:
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nrums = 1, nTotalTaxa = c(30, 40), nExtant = 0)
taxa <- fossilRecord2fossilTaxa(record)
sampledOccurrences <- sampleRanges(taxa, r = 0.5, ranges.only = FALSE)

# now try with horizonSampRate
horizonSampRate(sampOcc = sampledOccurrences)

# but we could also try with the *other* inputs
# useful because some datasets we may only have durations
# and number of sampling events for
```
inverseSurv <- sampledOccurrences(!is.na(sampledOccurrences)]
dur <- sapply(filtered, max) - sapply(filtered, min)
nCol <- sapply(filtered, length)
# supply as durations and nCollections
horizonSampRate(durations = dur, nCollections = nCol)

inverseSurv  Inverse Survivorship Models in the Fossil Record

Description

This function replicates the model-fitting procedure for forward and reverse survivorship curve data, described by Michael Foote in a series of papers (2001, 2003a, 2003b, 2005). These methods are discrete interval taxon ranges, as are used in many other functions in paleotree (see function arguments). This function can implement the continuous time, pulsed interval and mixed models described in Foote (2003a and 2005).

Usage

make_inverseSurv(timeList, groups = NULL, p_cont = TRUE, q_cont = TRUE,
                 PA_n = "fixed", PB_1 = 0, Nb = 1, drop.extant = TRUE)

Arguments

timelist A 2 column matrix with the first and last occurrences of taxa given in relative time intervals (i.e. ordered from first to last). If a list of length two is given for timeData, such as would be expected if the output of binTimeData was directly input, the second element is used. See details. Unsampled taxa (e.g. from a simulation of sampling in the fossil record, listed as NAs in the second matrix) are automatically dropped from the timeList and from groups simultaneously. Living taxa observed in the modern day are expected to be listed as last observed in a special interval (0,0), i.e. begins and ends at 0 time. This interval is always automatically removed prior to the calculation intermediary data for fitting likelihood functions.

groups Either NULL (the default) or matrix with the number of rows equal to the number of taxa and the number of columns equal to the number of 'systems' of categories for taxa. Taxonomic membership in different groups is indicated by numeric values. For example, a dataset could have a 'groups' matrix composed of a column representing thin and thick shelled taxa, coded 1 and 2 respectively, while the second column indicates whether taxa live in coastal, outer continental shelf, or deep marine settings, coded 1-3 respectively. Different combinations of groups will be treated as having independent sampling and extinction parameters in the default analysis, for example, thinly-shelled deep marine species will have separate parameters from thinly-shelled coastal species. Grouping systems could also represent temporal heterogeneity, for example, categorizing Paleozoic versus Mesozoic taxa. If groups are NULL (the default), all taxa are assumed to be of the same group with the same parameters. Unsampled taxa (e.g.
from a simulation of sampling in the fossil record, listed as NAs in timeData or timeList) are automatically dropped from groupings and the time dataset (either timeData or timeList) and from groups simultaneously.

**p_cont**

If TRUE (the default), then origination is assumed to be a continuous time process with an instantaneous rate. If FALSE, the origination is treated as a pulsed discrete-time process with a probability.

**q_cont**

If TRUE (the default), then extinction is assumed to be a continuous time process with an instantaneous rate. If FALSE, the extinction is treated as a pulsed discrete-time process with a probability.

**PA_n**

The probability of sampling a taxon after the last interval included in a survivorship study. Usually zero for extinct groups, although more logically has the value of 1 when there are still extant taxa (i.e., if the last interval is the Holocene and the group is still alive, the probability of sampling them later is probably 1...). Should be a value of 0 to 1, NULL, or can be simply "fixed", the default option. This default "fixed" option allows make_inverseSurv to decide the value based on whether there is a modern interval (i.e., an interval that is cHPLPI) or not: if there is one, then PA_n = 1, if not, then PA_n = 0. If NULL, PA_n is treated as an additional free parameter in the output model.

**PB_1**

The probability of sampling a taxon before the first interval included in a survivorship study. Should be a value of 0 to 1, or NULL. If NULL, PB_1 is treated as an additional free parameter in the output model.

**Nb**

The number of taxa that enter an interval (b is for 'bottom'). This is an arbitrary constant used to scale other values in these calculations and can be safely set to 1.

**drop.extant**

Drops all extant taxa from a dataset, w

**Details**

The design of this function to handle mixed continuous and discrete time models means that parameters can mean very different things, dependent on how a model is defined. Users should carefully evaluate their function arguments and the discussion of parameter values as described in the Value section.

**Value**

A function of class "paleotreeFunc", which takes a vector equal to the number of parameters and returns the *negative* log likelihood (for use with optim and similar optimizing functions, which attempt to minimize support values). See the functions listed at modelMethods for manipulating and examining such functions and constrainParPaleo for constraining parameters.

The function output will take the largest number of parameters possible with respect to groupings and time-intervals, which means the number of parameters may number in the hundreds. Constraining the function for optimization is recommended except when datasets are very large.

Parameters in the output functions are named 'p', 'q' and 'r', which are respectively the origination, extinction and sampling parameters. If the respective arguments 'p_cont' and 'q_cont' are TRUE, then 'p' and 'q' will represent the instantaneous per-capita origination and extinction rates (in units of per lineage time-units). When one of these arguments is given as FALSE, the respective parameter (p or q) will represent per-lineage-interval rates. For p, this will be the per lineage-interval
rate of a lineage producing another lineage (which can exceed 1 because diversity can more than double) and for q, this will be the per lineage-interval 'rate' of a lineage going extinct, which cannot be observed to exceed 1 (because the proportion of diversity that goes extinct cannot exceed 1). To obtain the per lineage-interval rates from a set of per lineage-time unit rates, simply multiply the per lineage-time-unit rate by the duration of an interval (or divide, to do the reverse; see Foote, 2003 and 2005). 'r' is always the instantaneous per-capita sampling rate, in units per lineage-time units.

If PA_n or PB_1 were given as NULL in the arguments, two additional parameters will be added, named respectively 'PA_n' and 'PB_1', and listed separately for every additional grouping. These are the probability of a taxon occurring before the first interval in the dataset (PB_1) and the probability of a taxon occurring after the last interval in a dataset (PA_n). These will be listed as 'PA_n.0' and 'PB_1.0' to indicate that they are not related to any particular time-interval included in the analysis, unlike the p, q, and r parameters (see below).

Groupings follow the parameter names, separated by periods; by default, the parameters will be placed in groups corresponding to the discrete intervals in the input timeList, such that make_inverseSurv will create a function with parameters 'p.1', 'q.1' and 'r.1' for interval 1; 'p.2', 'q.2' and 'r.2' for interval 2 and so on. Additional groupings given by the user are listed after this first set (e.g. 'p.1.2.2').

Calculating The Results of an Inverse Survivorship Model: Because of the complicated grouping and time interval scheme, combined with the probable preference of users to use constrained models rather than the full models, it may be difficult to infer what the rates for particular intervals and groups actually are in the final model, given the parameters that were found in the final optimization.

To account for this, the function output by inverseSurv also contains an alternative mode which takes input rates and returns the final values along with a rudimentary plotting function. This allows users to output per-interval and per-group parameter estimates. To select these feature, the argument altMode must be TRUE. This function will invisibly return the rate values for each group and interval as a list of matrices, with each matrix composed of the p, q and r rates for each interval, for a specific grouping.

This plotting is extremely primitive, and most users will probably find the invisibly returned rates to be of more interest. The function layout is used to play plots for different groupings in sequence, and this may lead to plots which are either hard to read or even cause errors (because of too many groupings, producing impossible plots). To repress this, the argument plotPar can be set to FALSE.

This capability means the function has more arguments that just the usual par argument that accepts the vector of parameters for running an optimization. The first of these additional arguments, altMode enables this alternative mode, instead of trying to estimate the negative log-likelihood from the given parameters. The other arguments augment the calculation and plotting of rates.

To summarize, a function output by inverseSurv has the following arguments:

**par** A vector of parameters, the same length as the number of parameters needed. For plotting, can be obtained with optimization

**altMode** If FALSE (the default) the function will work like ordinary model-fitting functions, returning a negative log-likelihood value for the input parameter values in par. If TRUE, however, the input parameters will instead be translated into the by-interval, by-group rates used for calculating the log-likelihoods, plotted (if plotPar is TRUE) and these final interval-specific rates will be returned invisibly as described above.
plotPar  If TRUE (the default) the calculated rates will be plotted, with each grouping given a separate plot. This can be repressed by setting plotPar to FALSE. As the only conceivable purpose for setting plotPar to FALSE is to get the calculated rates, these will not be returned invisibly if plotPar is FALSE.

ratesPerInt  If FALSE, the default option, the rates plotted and returned will be in units per lineage-time units, if those rates were being treated as rates for a continuous-time process (i.e. \( p_{\text{cont}} = \text{TRUE} \) and \( q_{\text{cont}} = \text{TRUE} \) for \( p \) and \( q \), respectively, while \( r \) is always per lineage-time units). Otherwise, the respective rate will be in units per lineage-interval. If ratesPerInt is TRUE instead, then all rates, even rates modeled as continuous-time process, will be returned as per lineage-interval rates, even the sampling rate \( r \).

logRates  If FALSE (the default) rates are plotted on a linear scale. If TRUE, rates are plotted on a vertical log axis.

jitter  If TRUE (default) the sampling rate and extinction rate will be plotted slightly ahead of the origination rate on the time axis, so the three rates can be easily differentiated. If false, this is repressed.

legendPosition  The position of a legend indicating which line is which of the three rates on the resulting plot. This is given as the possible positions for argument 'x' of the function `legend`, and by default is "topleft", which will be generally useful if origination and extinction rates are initially low. If legendPosition is NA, then a legend will not be plotted.

Note

This function is an entirely new rewrite of the methodology derived and presented by Foote in his studies. Thus, whether it would give identical results cannot be assumed nor is it easy to test given differences in the way data is handled between our coded functions. Furthermore, there may be differences in the math due to mistakes in the derivations caught while this function was programmed. I have tested the function by applying it to the same Sepkoski genus-level dataset that Foote used in his 2003 and 2005 papers. Users can feel free to contact me for detailed figures from this analysis. Overall, it seems my function captured the overall pattern of origination and sampling rates, at least under a model where both origination and extinction are modeled as continuous-time processes. Extinction showed considerably more variability relative to the published figures in Foote (2005). Additional analyses are being run to identify the sources of this discrepancy, and the function is being released here in paleotree on a trial basis, so that it can be more easily loaded onto remote servers. Users should be thus forewarned of this essentially 'beta' status of this function.

Author(s)

David W. Bapst, with some advice from Michael Foote.

References


See Also

This function extensively relies on `footeValues`. A similar format for likelihood models can be seen in `durationFreq`. Also see `freqRat`, `sRate2sProb`, `qsRate2Comp`, `sProb2sRate` and `qsProb2Comp`. For translating between sampling probabilities and sampling rates, see `SamplingConv`.

Examples

```r
# let's simulate some taxon ranges from an imperfectly sampled fossil record
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1, nTotalTaxa = c(30, 40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
rangesCont <- sampleRanges(taxa, r = 0.5)
# bin the ranges into discrete time intervals
rangesDisc <- binTimeData(rangesCont, int.length = 5)

# apply `make_inverseSurv`
likFun <- make_inverseSurv(rangesDisc)
# use `constrainParPaleo` to make the model time-homogenous

# `match.all=match.all` will match parameters so only 2 pars: p = q and r
constrFun <- constrainParPaleo(likFun, match.all=match.all)
results <- optim(parInit(constrFun), constrFun,
                 lower = parLower(constrFun), upper = parUpper(constrFun),
                 method = "L-BFGS-B", control = list(maxit = 1000000))
results

# plot the results
constrFun(results$par, altMode = TRUE)

# unconstrained function with ALL of 225 parameters!!!
# this will take forever to converge, so it isn't run
optim(parInit(likFun),likFun,lower = parLower(likFun),upper = parUpper(likFun),
       method = "L-BFGS-B",control = list(maxit = 1000000))
```

---

**Example Species Abundances Tables**

**Description**

A totally fictional example of species abundance data, for testing functions that require a site-by-taxon table of community ecology data.

**Format**

A table of type integer, representing terrestrial fauna and flora abundance counts.
Details

A classic dataset of ecological data collected by Satoshi and Okido, consisting of individual counts for 54 terrestrial faunal and floral species, from 23 sites across the mainland Kanto region.

Different ontogenetic stages were compounded and recorded by the common name for the first ontogenetic stage, with some inconsistency for species whose earliest stage have only been recently recognized. When separate names are commonly applied to sexual dimorphic forms, these were also combined and a single common name was used.

*Note: This data is a totally made-up, satirical homage to a well-known video game series (thus constituting fair-use).*

Source

Pokemon And All Respective Names are Trademark and Copyright of Nintendo 1996-2015.

See Also

twoWayEcologyCluster, communityEcology

Examples

data(kanto)

#visualize site abundances as barplots
barplotAbund <- function(x){
x <- x[,colSums(x)>0]
layout(1:(nrow(x)*1))
xpar <- par(mar = c(0,7,2,0))
for(i in 1:(nrow(x)-1)){
  barplot(x[i,],ylab = rownames(x)[i],
         names.arg = "")
}
barplot(x[nrow(x),],
        ylab = rownames(x)[nrow(x)],las = 3)
par(xpar)
layout(1)
mtext("Abundances",side = 2,line = 3,adj = 0.8)
}

#first five sites
kanto5 <- kanto[1:5,]
barplotAbund(kanto5)

#get pairwise Spearman rho coefficients
rhoCoeff <- pairwiseSpearmanRho(kanto,dropAbsent = "bothAbsent")

#what are the nearest-neighbor rhos (largest rho correlations)?
diag(rhoCoeff) <- NA
rhoNearest <- apply(rhoCoeff,1,max,na.rm = TRUE)
rhoNearest
# We can see the power plant sample is extremely different from the rest
# measure evenness: Hurlbert's PIE

kantoPIE <- HurlbertPIE(kanto)

# compare to dominance (relative abundance of most abundant taxon)
dominance <- apply(kanto,1,function(x) max(x)/sum(x) )

plot(kantoPIE,dominance)

# relatively strong relationship!

# Not run:
# Some Cool Ecology Stuff With Other Packages

# basically all the analyses & visualizations
# for ecology in R that I think are awesome

Ordination (PCO, DCA)

# get bray-curtis distances
library(vegan)
bcDist <- vegdist(kanto,method = "bray")

# do a PCO on the bray-curtis distances
pcoRes <- pcoa(bcDist,correction = "lingoes")
scores <- pcoRes$vectors

# plot the PCO
plot(scores,type = "n")
text(labels = rownames(kanto),scores[,1],scores[,2],cex = 0.5)

# the way the power plant and the pokémon tower converge
# is very suspicious: may be distortion due to a long gradient

# do a DCA instead with vegan's decorana
dcaRes <- decorana(kanto)

# plot using native vegan functions
# will show species scores in red
plot(dcaRes,cex = 0.5)
# kind of messy

# show just the sites scores
plot(dcaRes,cex = 0.5,display = "sites")

# show just the species scores
plot(dcaRes,cex = 0.5,display = "species")
#well, that's pretty cool

# get the nearest neighbor for each site
# based on pair-wise rho coefficients
rhoNeighbor <- apply(rhoCoef,1,function(x)
  rownames(kanto)[tail(order(x,na.last = NA),1)]
)

# let's plot the nearest neighbor connections with igraph
NNTable <- cbind(rownames(kanto),rhoNeighbor)

# now plot with igraph
library(igraph)
NList <- graph.data.frame(NNTable)
plot(NList)

# arrows point at the nearest neighbor of each sample
# based on maximum Spearman rho correlation

# Two Way Cluster With Heatmap

# This example based on code provided by Max Christie

# load pheatmap library for this example
library(pheatmap)

# get distance matrices for sites and taxa
# based on bray-curtis dist
# standardized to total abundance

# standardize site matrix to relative abundance
siteStand <- decostand(kanto, method = "total")
# site distance matrix (Bray-Curtis)
siteDist <- vegdist(siteStand, "bray", diag = TRUE)

# standardize taxa matrix to relative abundance
taxaStand <- decostand(t(kanto), method = "total")
# taxa distance matrix (Bray-Curtis)
taxaDist <- vegdist(taxaStand, "bray", diag = TRUE)

### Need to set graphic parameters for table

# Check out range of values for relative abundance
# hist(myStand) # none get very high...

# number of breaks: number of colors for heatmap
nBreaks <- 15

# set underValue
# anything below this counts as not appearing
# at that site for visualization purposes
```r
underValue <- min(siteStand[siteStand>0])-min(siteStand[siteStand>0])/10
# set overValue (max relative abundance)
overValue <- max(siteStand)

# you can set your breaks to any sequence you want
# and they don't have to be the same length.
# You can do this manually too.
# here we added a 0 to 'underValue' bin to
# the heatmap, making this bin essentially 0.
colorBreaks <- c(0, seq(underValue, max(siteStand), by = max(siteStand)/(nBreaks-1)))

# here we used the function rainbow to create a vector of colors.
# You can set these colors yourself too.
# It is important that this vector is one element
# less than the myBreaks vector
rainColors <- rainbow(nBreaks)

# now we can add "white" onto the vector,
# this will be the first color bin,
# which we're going to set to be (essentially) 0.
rainColors <- c("white", rainColors)
# If you don't add white, taxa at 0 abundance get colored in

### Plot the 2-Way Cluster

# heatmap, with user-set colors
# feed the function a distance matrix we wanted to use.
# siteDist and taxaDist made above by vegdist (bray-curtis)
# scale is the relative abundance, let's label it as such

dev.new(width = 10)
# for some reason, mtext() doesn't recognize pheatmap as plot.new
plot.new(width = 7)
pheatmap(siteStand, color = rainColors, breaks = colorBreaks,
        clustering_method = "ward.D",
        clustering_distance_rows = siteDist,
        clustering_distance_cols = taxaDist)
mtext("Relative Abundance", side = 4, line = -1.4, adj = 0.95)

# pretty cool looking!

#########################################
## Testing for differences between groups of sites

# is there a difference between routes and non-routes
groups <- rep(0, nrow(kanto))
groups[grep(rownames(kanto), pattern = "Route")]

# anosim (in vegan)
# are distances within groups smaller than distances between?
# we could also use adonis from vegan instead
library(vegan)

anosim(dat = kanto, grouping = groups)
```
macroperforateForam

Ancestor-Descendant Relationships for Macroperforate Foraminifera, from Aze et al. (2011)

Description

An example dataset of ancestor-descendant relationships and first and last appearance dates for a set of macroperforate Foraminifera, taken from the supplemental materials of Aze et al. (2011). This dataset is included here primarily for testing functions parentChild2taxonTree and taxa2phylo.

Format

The ‘foramAM’ and ‘foramAL’ tables include budding taxon units for morphospecies and lineages respective, with four columns: taxon name, ancestral taxon’s name, first appearance date and last appearance date (note that column headings vary). The ‘foramAMb’ and ‘foramALb’ tables are composed of data for the same taxon units as the previous branching events are split so that the relationships are fully ‘bifurcating’, rather than ‘budding’. As this obscures taxonomic identity, taxon identification labels are included in an additional, fifth column in these tables. See the examples section for more details.

Details

This example dataset is composed of four tables, each containing information on the ancestor-descendant relationships and first and last appearances of species of macroperforate foraminifera species from the fossil record. Each of the four tables are for the same set of taxa, but divide and concatenate the included foram species in four different ways, relating to the use of morphospecies versus combined anagenetic lineages (see Ezard et al., 2012), and whether taxa are retained as units related by budding-cladogenesis or the splitting of taxa at branching points to create a fully ‘bifurcating’ set of relationships, independent of ancestral morphotaxon persistence through branching events. See the examples section for more details.
macroperforateForam

Source

This dataset is obtained from the supplementary materials of, specifically ‘Appendix S5’:

References

This dataset has been used or referenced in a number of works, including:

Examples

# Following Text Reproduced from Aze et al. 2011's Supplemental Material
# Appendix S5
#
# 'Data required to produce all of the phylogenies included in the manuscript
# using paleoPhylo (Ezard & Purvis, 2009) a free software package to draw
# paleobiological phylogenies in R.'
#
# 'The four tabs hold different versions of our phylogeny:
# aMb: fully bifurcating morphospecies phylogeny
# aM: budding/bifurcating morphospecies phylogeny
# aLb: fully bifurcating lineage phylogeny
# aL: budding/bifurcating lineage phylogeny
#
# 'Start Date gives the first occurrence of the species according
# to the particular phylogeny; End Date gives the last occurrence
# according to the particular phylogeny.'

## Not run:

#load the data
# given in supplemental as XLS sheets
# converted to separate tab-delimited text files

# aM: budding/bifurcating morphospecies phylogeny
foramAM <- read.table(file.choose(),stringsAsFactors = FALSE,header = TRUE)

# aL: budding/bifurcating lineage phylogeny
foramAL <- read.table(file.choose(),stringsAsFactors = FALSE,header = TRUE)

# aMb: fully bifurcating morphospecies phylogeny
foramAMb <- read.table(file.choose(),stringsAsFactors = FALSE,header = TRUE)

# aLb: fully bifurcating lineage phylogeny
foramALb <- read.table(file.choose(),stringsAsFactors = FALSE,header = TRUE)

save.image("macroperforateForam.rdata")

## End(Not run)

# instead, we'll just load the data directly
data(macroperforateForam)

# Two distinctions among the four datasets:
# (1): morphospecies vs morphospecies combined into sequences of anagenetic
# morphospecies referred to as 'lineages'. Thus far more morphospecies
# than lineages. The names of lineages are given as the sequence of
# their respective component morphospecies.
# (2): Datasets where taxon units (morphospecies or lineages) are broken up
# at 'budding' branching events (where the ancestral taxon persists)
# so that final dataset is 'fully bifurcating', presumably
# to make comparison easier to extant-taxon only datasets.
# (This isn't a limitation for paleotree, though!).
# This division of taxon units requires abstracting the taxon IDs,
# requiring another column for Species Name.

dim(foramAM)
dim(foramAL)
dim(foramAMb)
dim(foramALb)

# Need to convert these to same format as fossilRecord2fossilTaxa output.
# those 'taxa' tables has 6 columns:
# taxon.id ancestor.id orig.time ext.time still.alive looks.like

# for the purposes of this, we'll make taxon.id = looks.like
# (That's only for simulating cryptic speciation anyway)
# still.alive should be TRUE (1) if ext.time = 0

# a function to convert Aze et al's suppmat to paleotree-readable format

createTaxaData <- function(table){
  # reorder table by first appearance time
  table <- table[order(as.numeric(table[,3])),]
  ID <- 1:nrow(table)
  anc <- sapply(table[,2],function(x)
if(!is.na(x)){
  which(x == table[,1])
}else{ NA })
stillAlive <- as.numeric(table[,4] == 0)
ages <- cbind(as.numeric(table[,3]),as.numeric(table[,4]))
res <- cbind(ID,anc,ages,stillAlive,ID)
colnames(res) <- c('taxon.id','ancestor.id','orig.time','ext.time','still.alive','looks.like')
rownames(res) <- table[,1]
return(res)
}
taxaAM <- createTaxaData(foramAM)
taxaAMb <- createTaxaData(foramAMb)
taxaAL <- createTaxaData(foramAL)
taxaALb <- createTaxaData(foramALb)

#########################################################################
#Checking Ancestor-Descendant Relationships for Irregularities
#For each of these, there should only be a single taxon
#without a parent listed (essentially, the root ancestor)

countParentsWithoutMatch <- function(table){
  parentMatch <- match(unique(table[,2]),table[,1])
  sum(is.na(parentMatch))
}
#test this on the provided ancestor-descendant relationships
countParentsWithoutMatch(foramAM)
countParentsWithoutMatch(foramAL)
countParentsWithoutMatch(foramAMb)
countParentsWithoutMatch(foramALb)

#and on the converted datasets
countParentsWithoutMatch(taxaAM)
countParentsWithoutMatch(taxaAL)
countParentsWithoutMatch(taxaAMb)
countParentsWithoutMatch(taxaALb)

#can construct the parentChild2taxonTree
#using the ancestor-descendant relationships
#can be very slow...

treeAM <- parentChild2taxonTree(foramAM[,2:1])
treeAL <- parentChild2taxonTree(foramAL[,2:1])
treeAMb <- parentChild2taxonTree(foramAMb[,2:1])
treeALb <- parentChild2taxonTree(foramALb[,2:1])
macroperforateForam

```
layout(matrix(1:4,2,2))
plot(treeAM,main = 'treeAM',show.tip.label = FALSE)
plot(treeAL,main = 'treeAL',show.tip.label = FALSE)
plot(treeAMB,main = 'treeAMB',show.tip.label = FALSE)
plot(treeALB,main = 'treeALB',show.tip.label = FALSE)

# FYI
# in case you were wondering
# you would *not* time-scale these Frankenstein monsters

########################################################################

# Checking stratigraphic ranges

# do all first occurrence dates occur before last occurrence dates?
# we'll check the original datasets here

ccheckFolO <- function(data){
diffDate <- data[,3]-data[,4] #subtract LO from FO
isGood <- all(diffDate >= 0) #is it good
return(isGood)
}

ccheckFolO(foramAM)
ccheckFolO(foramAL)
ccheckFolO(foramAMB)
ccheckFolO(foramALB)

#cool, but do all ancestors appear before their descendants?
# easier to check unified fossilRecord2fossilTaxa format here

ccheckAncOrdE <- function(taxa){
#get ancestor's first occurrence
ancFO <- taxa[taxa[,2],3]
#get descendant's first occurrence
descFO <- taxa[,3]
diffDate <- ancFO-descFO #subtract descFO from ancFO
#remove NAs due to root taxon
diffDate <- diffDate[!is.na(diffDate)]
isGood <- all(diffDate >= 0) #is it all good
return(isGood)
}

ccheckAncOrdE(taxaAM)
ccheckAncOrdE(taxaAL)
ccheckAncOrdE(taxaAMB)
ccheckAncOrdE(taxaALB)

#now, are there gaps between the last occurrence of ancestors
# and the first occurrence of descendants?
# (shall we call these 'stratophenetic ghost branches'?)
```
# These shouldn't be problematic, but do they occur in this data?
# After all, fossilRecord2FossilTaxa output tables are designed for
# fully observed simulated fossil records with no gaps.

sumAncDescGap <- function(taxa){
  #get ancestor's last occurrence
  ancLO <- taxa[,2,4]
  #get descendant's first occurrence
  descFO <- taxa[,3]
  diffDate <- ancLO-descFO #subtract descFO from ancFO
  #remove NAs due to root taxon
  diffDate <- diffDate[!is.na(diffDate)]
  #should be negative or zero, positive values are gaps
  gaps <- c(0,diffDate[diffDate>0])
  sumGap <- sum(gaps)
  return(sumGap)
}

#get the total gap between ancestor LO and child FO
sumAncDescGap(taxaAM)
sumAncDescGap(taxaAL)
sumAncDescGap(taxaAMb)
sumAncDescGap(taxaALb)

#It appears there is *no* gaps between ancestors and their descendants
#in the Aze et al. foram dataset... wow!

Creating time-scaled phylogenies from the Aze et al. data

# Aze et al. (2011) defines anagenesis such that taxa may overlap
# in time during a transitional period (see Ezard et al. 2012
# for discussion of this definition). Thus, we would expect that
# paleotree obtains very different trees for morphospecies versus
# lineages, but very similar phylogenies for datasets where budding
# taxa are retained or arbitrarily broken into bifurcating units.

# We can use the function taxa2phylo to directly create
# time-scaled phylogenies from the Aze et al. stratophenetic data

timetreeAM <- taxa2phylo(taxaAM)
timetreeAL <- taxa2phylo(taxaAL)
timetreeAMb <- taxa2phylo(taxaAMb)
timetreeALb <- taxa2phylo(taxaALb)

layout(matrix(1:4,2,2))
plot(timetreeAM,main = 'timetreeAM',show.tip.label = FALSE)
axisPhylo()
plot(timetreeAL,main = 'timetreeAL',show.tip.label = FALSE)
axisPhylo()}
plot(timetreeAMB, main = 'timetreemAMB', show.tip.label = FALSE)
axisPhylo()
plot(timetreeALb, main = 'timetreemALb', show.tip.label = FALSE)
axisPhylo()

# visually compare the two pairs we expect to be close to identical

# morphospecies
layout(1:2)
plot(timetreeAM, main = 'timetreemAM', show.tip.label = FALSE)
axisPhylo()
plot(timetreeAMB, main = 'timetreemAMB', show.tip.label = FALSE)
axisPhylo()

# lineages
layout(1:2)
plot(timetreeAL, main = 'timetreemAL', show.tip.label = FALSE)
axisPhylo()
plot(timetreeALb, main = 'timetreemALb', show.tip.label = FALSE)
axisPhylo()

layout(1)

# compare the summary statistics of the trees
Ntip(timetreeAM)
Ntip(timetreeAMB)
Ntip(timetreeAL)
Ntip(timetreeALb)

# very different!

# after dropping anagenetic zero-length-terminal-edge ancestors
# we would expect morphospecies and lineage phylogenies to be very similar

# morphospecies
Ntip(dropZLB(timetreeAM))
Ntip(dropZLB(timetreeAMB))

# identical!

# lineages
Ntip(dropZLB(timetreeAL))
Ntip(dropZLB(timetreeALb))

# ah, very close, off by a single tip
# ... probably a very short ZLB outside tolerance

# we can create some diversity plots to compare

multiDiv(data = list(timetreeAM, timetreemAMB),
plotMultCurves = TRUE)

multiDiv(data = list(timetreeAL, timetreemALb),
plotMultCurves = TRUE)

# we can see that the morphospecies datasets are identical
# that's why we can only see one line
# some very slight disagreement between the lineage datasets
# around ~38-20 Ma

# can also compare morphospecies and lineages diversity curves

multiDiv(data = list(timetreeAM, timetreeAL),
plotMultCurves = TRUE)

# they are similar, but some peaks are missing from lineages
# particularly around ~20-10 Ma

---

**makePBDBtaxonTree**

*Creating a Taxon-Tree from Taxonomic Data Downloaded from the Paleobiology Database*

**Description**

This function creates phylogeny-like object of type phylo from the taxonomic information recorded in a taxonomy download from the PBDB for a given group. Two different algorithms are provided, the default being based on parent-child taxon relationships, the other based on the nested Linnean hierarchy.

**Usage**

```r
makePBDBtaxonTree(data, rank, method = "parentChild", solveMissing = NULL,
tipSet = "nonParents", cleanTree = TRUE, APIversion = "1.1")
```

**Arguments**

- **data**
  A table of taxonomic data collected from the Paleobiology Database, using the taxa list option with show = phylo. Should work with versions 1.1-1.2 of the API, with either the 'pdbb' or 'com' vocab. However, as 'accepted_name' is not available in API v1.1, the resulting tree will have a taxon’s *original* name and not any formally updated name.

- **rank**
  The selected taxon rank; must be one of 'species', 'genus', 'family', 'order', 'class' or 'phylum'.

- **method**
  Controls which algorithm is used for calculating the taxon-tree: either method = "parentChild" (the default option) which converts the listed binary parent-child taxon relationships from the input PBDB data, or method = "Linnean", which converts a taxon-tree by creating a table of the Linnean taxonomic assignments (family, order, etc), which are provided when option 'show = phylo' is used in PBDB API calls.
makePBDBtaxonTree

solveMissing  Under method = "parentChild", what should makePBDBtaxonTree do about multiple 'floating' parent taxa, listed without their own parent taxon information in the input dataset under data? Each of these is essentially a separate root taxon, for a different set of parent-child relationships, and thus poses a problem as far as returning a single phylogeny is concerned. If solveMissing = NULL (the default), nothing is done and the operation halts with an error, reporting the identity of these taxa. Two alternative solutions are offered: first, solveMissing will combine these disparate potential roots and link them to an artificially-constructed pseudo-root, which at least allows for visualization of the taxonomic structure in a limited dataset. Secondly, solveMissing = "mergeroots" queries the Paleobiology Database repeatedly via the API for information on parent taxa of the 'floating' parents, and continues within a while() loop until only one such unassigned parent taxon remains. This latter option may talk a long time or never finish, depending on the linearity and taxonomic structures encountered in the PBDB taxonomic data; i.e. if someone a taxon was ultimately its own indirect child in some grand loop by mistake, then under this option makePBDBtaxonTree might never finish. In cases where taxonomy is bad due to weird and erroneous taxonomic assignments reported by the PBDB, this routine may search all the way back to a very ancient and deep taxon, such as the Eukaryota taxon. Users should thus use solveMissing = "queryPBDB" only with caution.

tipSet  This argument only impacts analyses where the argument method = "parentChild" is also used. This tipSet controls which taxa are selected as tip taxa for the output tree. The default tipSet = "nonParents" selects all child taxa which are not listed as parents in parentChild. Alternatively, tipSet = "all" will add a tip to every internal node with the parent-taxon name encapsulated in parentheses.

cleanTree By default, the tree is run through a series of post-processing, including having singles collapsed, nodes reordered and being written out as a Newick string and read back in, to ensure functionality with ape functions and ape-derived functions. If FALSE, none of this post-processing is done and users should beware, as such trees can lead to hard-crashes of R.

APIversion Version of the Paleobiology Database API used by makePBDBtaxonTree when solveMissing = "queryPBDB". The current default is "1.1", which is the only option available as of 05/05/2015. In the future, the improved API version "1.2" will be released on the public PBDB server, which will become the new default for this function, but the option to return to "1.1" behavior will be retained for.

Details

This function should not be taken too seriously. Many groups in the Paleobiology Database have out-of-date or very incomplete taxonomic information. This function is meant to help visualize what information is present, and by use of time-scaling functions, allow us to visualize the intersection of temporal and phylogenetic, mainly to look for incongruence due to either incorrect taxonomic placements, erroneous occurrence data or both.

Note however that, contrary to common opinion among some paleontologists, taxon-trees may be just as useful for macroevolutionary studies as reconstructed phylogenies (Soul and Friedman, in press.).
Value

A phylogeny of class 'phylo', where each tip is a taxon of the given 'rank'. See additional details regarding branch lengths can be found in the sub-algorithms used to create the taxon-tree by this function: `parentChild2taxonTree` and `taxonTable2taxonTree`.

Depending on the method used, either the element `parentChild` or `taxonTable` is added to the list structure of the output phylogeny object, which was used as input for one of the two algorithms mentioned above.

Please note that when applied to output from the taxa option of the API version 1.1, the taxon names returned are the original taxon names as 'accepted_name' is not available in API v1.1, while under API v1.2, the returned taxon names should be the most up-to-date formal names for those taxa. Similar issues also affect the identification of parent taxa, as the accepted name of the parent ID number is only provided in version 1.2 of the API.

Author(s)

David W. Bapst

References


See Also

Two other functions in paleotree are used as sub-algorithms by `makePBDBtaxonTree` to create the taxon-tree within this function, and users should consult their manual pages for additional details: `parentChild2taxonTree` and `taxonTable2taxonTree`.

Other functions for manipulating PBDB data can be found at `taxonSortPBDBocc.occData2timeList`, and the example data at `graptPBDB`.

Examples

```r
## Not run:

easyGetPBDBtaxa <- function(taxon,show = c("phylo","img","app")){
  #let's get some taxonomic data
taxaData <- read.csv(paste0("http://paleobiodb.org/",
  "data1.1/taxa/list.txt?base_name = ",taxon,
  "&rel = all_children&show = ",
paste0(show,collapse = ","),"&status = senior"),
  stringsAsFactors = FALSE)
  return(taxaData)
}

#graptolites
graptData <- easyGetPBDBtaxa("Graptolithina")
graptTree <- makePBDBtaxonTree(graptData,"genus",
  method = "parentChild", solveMissing = "queryPBDB")
#try Linnean
```
graptTree <- makePBDBtaxonTree(graptData, "genus", method = "Linnean")
plot(graptTree, show.tip.label = FALSE, no.margin = TRUE, edge.width = 0.35)
nodelabels(graptTree$node.label, adj = c(0,1/2))

# conodonts
conoData <- easyGetPDBBtaxa("Conodonta")
conoTree <- makePBDBtaxonTree(conoData, "genus", method = "parentChild", solveMissing = "queryPDBB")
plot(conoTree, show.tip.label = FALSE, no.margin = TRUE, edge.width = 0.35)
nodelabels(conoTree$node.label, adj = c(0,1/2))

# asaphid trilobites
asaData <- easyGetPDBBtaxa("Asaphida")
asaTree <- makePBDBtaxonTree(asaData, "genus", method = "parentChild", solveMissing = "queryPDBB")
plot(asaTree, show.tip.label = FALSE, no.margin = TRUE, edge.width = 0.35)
nodelabels(asaTree$node.label, adj = c(0,1/2))

# Ornithischia
ornithData <- easyGetPDBBtaxa("Ornithischia")
ornithTree <- makePBDBtaxonTree(ornithData, "genus", method = "parentChild", solveMissing = "queryPDBB")
# try Linnean
# need to drop repeated taxon first: Hylaeeosaurus
ornithData <- ornithData[!(which(ornithData[, "taxon_name"] == "Hylaeeosaurus") == 1),]
ornithTree <- makePBDBtaxonTree(ornithData, "genus", method = "Linnean")
plot(ornithTree, show.tip.label = FALSE, no.margin = TRUE, edge.width = 0.35)
nodelabels(ornithTree$node.label, adj = c(0,1/2))

# Rhynchonellida
rynchData <- easyGetPDBBtaxa("Rhynchonellida")
rynchTree <- makePBDBtaxonTree(rynchData, "genus", method = "parentChild", solveMissing = "queryPDBB")
plot(rynchTree, show.tip.label = FALSE, no.margin = TRUE, edge.width = 0.35)
nodelabels(rynchTree$node.label, adj = c(0,1/2))

# some of these look pretty messy!

## End(Not run)

########################################################################

# let's try time-scaling the graptolite tree

# get some example occurrence and taxonomic data
data(graptPDBB)

# get the taxon tree: Linnean method
graptTree <- makePBDBtaxonTree(graptTaxaPDBB, "genus", method = "Linnean")
plot(graptTree$node.label,cex = 0.5)

#get the taxon tree: parentChild method
graptTree <- makePDBtaxonTree(graptTaxaPDB, "genus", method = "parentChild")
plot(graptTree, cex = 0.4)
modelabels(graptTree$node.label,cex = 0.5)

# get time data from occurrences
graptOccGenus <- taxonSortPDBOcc(graptOccPDB, rank = "genus", onlyFormal = FALSE)
graptTimeGenus <- occData2timeList(occList = graptOccGenus)

# let's time-scale the parentChild tree with paleotree
# use minimum branch length for visualization
# and nonstoch.bin so we plot maximal ranges
timeTree <- bin_timePaleoPhy(graptTree, timeList = graptTimeGenus, nonstoch.bin = TRUE, type = "mbl", vartime = 3)

# drops a lot of taxa; some of this is due to mispellings, etc

## Not run:

## make pretty plot with library strap
library(strap)
geoscalePhylo(timeTree, ages = timeTree$ranges.used)
modelabels(timeTree$node.label,cex = 0.5)

## End(Not run)

---

### minBranchLength

**Scales Edge Lengths of a Phylogeny to a Minimum Branch Length**

**Description**

Rescales a tree with edge lengths so that all edge lengths are at least some minimum branch length (mbl), without changing the relative distance of the tips from the root node. Edge lengths are transformed so they are greater than or equal to the input minimum branch length, by subtracting edge length from more rootward edges and added to later branches.

**Usage**

```r
minBranchLength(tree, mbl)
```

**Arguments**

- **tree**: A phylogeny with edge lengths of class `phylo`.
- **mbl**: The minimum branch length
Details

This function was formally an internal segment in `timePaleoPhy`, and now is called by `timePaleoPhy` instead, allowing users to apply `minBranchLength` to trees that already have edge lengths.

Value

A phylogeny with edge lengths of class `phylo`.

Author(s)

David W. Bapst

See Also

This function was originally an internal piece of `timePaleoPhy`, which implements the minimum branch length time-scaling method along with others, which may be what you’re looking for (instead of this miscellaneous function).

Examples

```r
# simulation with an example non-ultrametric tree

tree <- rtree(20)
# randomly replace edges with ZLBs, similar to multi2di output
tree <- degradeTree(tree, 0.3, leave.zlb = TRUE)

tree2 <- minBranchLength(tree, 0.1)

layout(1:2)
plot(tree); axisPhylo()
plot(tree2); axisPhylo()

layout(1)

# now let's try it with an ultrametric case

# get a random tree

tree <- rtree(30)
# randomly replace edges with ZLBs, similar to multi2di output

tree <- degradeTree(tree, 0.5, leave.zlb = TRUE)
# now randomly resolve
tree <- di2multi(tree)
# give branch lengths so its ultrametric

plot(tree) # and we have an ultrametric tree with polytomies, yay!

# now randomly resolve, get new branch lengths as would with real data


tree2 <- multi2di(tree)
tree2 <- minBranchLength(tree2, 0.1)
```
**minCharChange**

Estimating the Minimum Number of Character Transitions Using Maximum Parsimony

**Description**

`minCharChange` is a function which takes a cladogram and a discrete trait and finds the solutions of inferred character states for ancestral nodes that minimizes the number of character state transitions (either gains or losses/reversals) for a given topology and a set of discrete character data. `minCharChange` relies on `ancPropStateMat`, which is a wrapper for `phangorn`'s function `ancestralPars`.

**Usage**

```r
minCharChange(trait, tree, randomMax = 10000, maxParsimony = TRUE, orderedChar = FALSE, type = "MPR", cost = NULL, printMinResult = TRUE, ambiguity = c(NA, "?"), dropAmbiguity = FALSE, polySymbol = "&", contrast = NULL)
```

```r
ancPropStateMat(trait, tree, orderedChar = FALSE, type = "MPR", cost = NULL, ambiguity = c(NA, "?"), dropAmbiguity = FALSE, polySymbol = "&", contrast = NULL, returnContrast = FALSE)
```

**Arguments**

- **trait**
  A vector of trait values for a discrete character, preferably named with taxon names identical to the tip labels on the input tree.

- **tree**
  A cladogram of type 'phylo'. Any branch lengths are ignored.

- **randomMax**
  The maximum number of cladograms examined when searching a large number of solutions consistent with the reconstructed ancestral states from `ancestralPars` with the minimum number of character state transitions. If the number of potential solutions is less than `randomMax`, then solutions are exhaustively searched.

- **maxParsimony**
  If `maxParsimony` is TRUE (the default) then only solutions with the smallest number of total transitions examined will be returned. Note that since solutions are stochastically 'guessed' at, and the number of possible solutions may not be exhaustively searched, there may have been solutions not examined with a lower number of transitions even if `maxParsimony` = TRUE. Regardless, one may want to do `maxParsimony` = FALSE if one is interested in whether there are solutions with a smaller number of gains or losses and thus wants to return all solutions.
orderedChar If TRUE (not the default), then the character will be reconstructed with a cost (step) matrix of a linear, ordered character. This is not applicable if type = "ACCTRAN", as cost matrices cannot be used with ACCTRAN in ancestral.pars, and an error will be returned if orderedChar = TRUE but a cost matrix is given, as the only reason to use orderedChar is to produce a cost matrix automatically.

type The parsimony algorithm applied by ancestral.pars, which can apply one of two: "MPR" (the default) is a relatively fast algorithm developed by Hanazawa et al. (1995) and Narushima and Hanazawa (1997), which relies on reconstructing the states at each internal node by re-rooting at that node. "ACCTRAN", the 'accelerated transitions' algorithm (Swofford and Maddison, 1987), favors character reversal over independent gains when there is ambiguity. The "ACCTRAN" option in ancestral.pars avoids repeated rerooting of the tree to search for a smaller set of maximum-parsimony solutions that satisfy the ACCTRAN algorithm, but does so by assigning edge weights. As of phangorn v1.99-12, "MPR" is calculated using the Sankoff parsimony algorithm, which allows multifurcations (polytomies). This is not true for "ACCTRAN", which uses the Fitch algorithm, which does not allow for multifurcations. An error is returned if a tree with multifurcations is passed and type = "ACCTRAN".

cost A matrix of the cost (i.e. number of steps) necessary to change between states of the input character trait. If NULL (the default), the character is assumed to be unordered with equal cost to change from any state to another. Cost matrices only impact the "MPR" algorithm; if a cost matrix is given but type = "ACCTRAN", an error is issued.

printMinResult If TRUE (the default), a summary of the results is printed to the terminal. The information in this summary may be more detailed if the results of the analysis are simpler (i.e. fewer unique solutions).

ambiguity A vector of values which indicate ambiguous (i.e. missing or unknown) character state codings in supplied trait data. Taxa coded ambiguously as treated as being equally likely to be any state coding. By default, NA values and "?" symbols are treated as ambiguous character codings, in agreement with behavior of functions in packages phangorn and claddis. This argument is designed to mirror an hidden argument with an identical name in function phydat in package phangorn.

dropAmbiguity A logical. If TRUE (which is not the default), all taxa with ambiguous codings as defined by argument ambiguity will be dropped prior to ancestral nodes being inferred. This may result in too few taxa.

polySymbol A single symbol which separates alternative states for polymorphic codings; the default symbol is "&", following the output by claddis's ReadMorphNexus function, where polymorphic taxa are indicated by default with a string with state labels separated by an "&" symbol. For example, a taxon coded as polymorphic for states 1 or 2, would be indicated by the string "1&2". polySymbol is used to break up these strings and automatically construct a fitting contrast table for use with this data, including for ambiguous character state codings.

contrast A matrix of type integer with cells of 0 and 1, where each row is labeled with a string value used for indicating character states in trait, and each column is labeled with the formal state label to be used for assign taxa to particular character states. A value of 1 indicates that the respective coding string for that row should
be interpreted as reflecting the character state listed for that column. A coding could reflect multiple states (such as might occur when taxa are polymorphic for some morphological character), so the sums of rows and columns can sum to more than 1. If contrast is not NULL (the default), the arguments will nullify. This argument is designed to mirror an hidden argument with an identical name in function phyDat in package phangorn. This structure is based on the contrasts tables used for statistical evaluation of factors. See the phangorn vignette "Special features of phangorn" for more details on its implementation within phangorn including an example. See examples below for the construction of an example contrast matrix for character data with polymorphisms, coded as character data output by Cladist's ReadMorphNexus function, where polymorphic taxa are indicated with a string with state labels separated by an "&" symbol.

returnContrast
If TRUE, the contrast table used by ancestral.pars will be output instead for user evaluation that polymorphic symbols and ambiguous states are being parsed correctly.

Details
The wrapper function ancPropStateMat simply automates the application of functions ancestral.pars and phyDat from phangorn, along with several additional checks and code to present the result as a matrix, rather than a specialized list.

Note that although the default orderedChar argument assumes that multistate characters are unordered, the results of character change will always be reported as gains and losses relative to the numbering of the states in the output transitionSumChanges, exactly as if they had been ordered. In the case where the character is actually ordered, this may be considered a conservative approach, as using a parsimony algorithm for unordered character states allows fewer gains or losses to be counted on branches where multiple gains and losses are reported. If the character is presumably unordered and multistate, however, then the gains and losses division is arbitrary nonsense and should be combined to to obtain the total number of character changes.

Value
By default, ancPropStateMat returns a matrix, with rows corresponding to the ID numbers of tips and nodes in $edge, and columns corresponding to character states, with the value representing the proportional weight of that node being that state under the algorithm used (known tip values are always 1). If argument returnContrast is TRUE then ancPropStateMat will instead return the final contrast table used by phyDat for interpreting character state strings.

minCharChange invisibly returns a list containing the following elements, several of which are printed by default to the console, as controlled by argument printMinResult:

message Describes the performance of minCharChange at searching for a minimum solution.
sumTransitions A vector recording the total number of necessary transitions (sum total of gains and losses/reversal) for each solution; effectively the parsimony cost of each solution.
minTransitions A symmetrical matrix with number of rows and columns equal to the number of character states, with values in each cell indicating the minimum number of transitions from one ancestral state (i.e. the rows) to a descendant state (i.e. the columns), taken across the set of kept solutions (dependent on which are kept as decided by argument maxParsimony).
Generally guaranteed not to add up to the number of edges contained within the input tree, and thus may not represent any realistic evolutionary scenario but does represent a conservative approach for asking 'what is the smallest possible number of transitions from 0 to 1' or 'smallest possible number of transitions from 1 to 0', independently of each other.

**solutionArray** A three-dimensional array, where for each solution, we have a matrix with edges for rows and two columns indicating the ancestral and child nodes of that edge, with values indicating the states inferred for those nodes in a particular solution.

**transitionArray** A labeled three-dimensional array where for each solution we have a symmetrical matrix with number of rows and columns equal to the number of character states, with values in each cell indicating the total number of transitions from one ancestral state (i.e. the rows) to a descendant state (i.e. the columns).

**transitionSumChanges** Which is a three column matrix with a row for every solution, with the values in the three columns measuring the number of edges (branches) inferred to respectively have gains, no change or losses (i.e. reversals), as calculated relative to the order of character states.

**Author(s)**

David W. Bapst

**References**


**See Also**

The functions described here are effectively wrappers of phangorn's function `ancestral.pars`.

**Examples**

```r
# let's write a quick & dirty ancestral trait plotting function

quickAncPlotter <- function(tree, ancData, cex){
    ancCol <- (1:ncol(ancData)) + 1
    plot(tree, show.tip.label = FALSE, no.margin = TRUE, direction = "upwards")
    tiplabels(pch = 16, pie = ancData[(1:Ntip(tree)),], cex = cex, piecol = ancCol, col = 0)
    nodelabels(pie = ancData[-(1:Ntip(tree))], cex = cex, piecol = ancCol)
}
```
# example with retiolitid graptolite data
data(retiolitiniae)

#unordered, MPR
ancMPR <- ancPropStateMat(retioTree, trait = retioChar[,2], type = "MPR")
quickAncPlotter(retioTree,ancMPR,cex = 0.5)
text(x = 4,y = 5,"type = 'MPR',cex = 1.5)

minCharChange(retioTree,trait = retioChar[,2],type = "MPR")

# with simulated data

set.seed(444)
tree <- rtree(50)
#simulate under a likelihood model
char <- rTraitDisc(tree,k = 3,rate = 0.7)
tree$edge.length <- NULL
tree <- ladderize(tree)

#unordered, MPR
ancMPR <- ancPropStateMat(tree, trait = char, type = "MPR")
#ordered, MPR
ancMPRord <- ancPropStateMat(tree, trait = char, orderedChar = TRUE, type = "MPR")

#let's compare unordered versus ordered under MPR
layout(1:2)
quickAncPlotter(tree,ancMPR,cex = 0.3)
text(x = 8,y = 15,"unordered char\nMPR",cex = 1.5)
quickAncPlotter(tree,ancMPRord,cex = 0.3)
text(x = 9,y = 15,"ordered char\nMPR",cex = 1.5)
layout(1)

## Not run: # what ancPropStateMat automates (with lots of checks):

require(phangorn)
char1 <- matrix(char,1)
rownames(char1) <- names(char)
#translate into something for phangorn to read
char1 <- phyDat(char1,type = "USER",levels = sort(unique(char1)))
x <- ancestral.pars(tree,char1,type = "MPR")
y <- ancestral.pars(tree,char1,type = "ACCTRAN")

## End(Not run)

#estimating minimum number of transitions with MPR
minCharChange(tree,trait = char,type = "MPR")

#POLYMORPHISM IN CHARACTER DATA
# example trait data with a polymorphic taxon
# separated with `&` symbol
# similar to polymorphic data output by readMorphNexus from package Claddis
charPoly <- as.character(c(1,2,NA,0,0,1,"1&2",2,0,NA,0,2,1,1,"1&2"))

# simulate a tree with 16 taxa
set.seed(444)
tree <- rtree(15)
tree$edge.length <- NULL
tree <- ladderize(tree)
names(charPoly) <- tree$tip.label
ccharPoly

# need a contrast matrix that takes this into account
# can build row by row, by hand

# first, build contrast matrix for basic states
contrast012 <- rbind(c(1,0,0),c(0,1,0),c(0,0,1))
colnames(contrast012) <- rownames(contrast012) <- 0:2
ccontrast012

# add polymorphic state and NA ambiguity as new rows
contrastPoly <- c(0,1,1)
contrastNA <- c(1,1,1)
contrastNew <- rbind(contrast012,'1&2' = contrastPoly,contrastNA)
rownames(contrastNew)[5] <- NA

# let's look at contrast
contrastNew

# now try this contrast table we've assembled
# default: unordered, MPR
ancPoly <- ancPropStateMat(tree, trait = charPoly, contrast = contrastNew)

# but...!
# we can also do it automatically,
# by default, states with `&` are automatically treated
# as polymorphic character codings by ancPropStateMat
ancPolyAuto <- ancPropStateMat(tree, trait = charPoly, polySymbol = "&")

# but does this match what the table we constructed?
ancPropStateMat(tree, trait = charPoly,
polySymbol = ",", returnContrast = TRUE)

# compare to contrastNew above!
# only difference should be the default ambiguous
# character `?' is added to the table

# compare reconstructions
layout(1:2)
quickAncPlotter(tree,ancPoly, cex = 0.5)
text(x = 3.5,y = 1.2,"manually-constructed\ncontrast",cex = 1.3)
quickAncPlotter(tree,ancPolyAuto, cex = 0.5)
### Model Function Methods: Parameter Names, Bounds and Initial Values

**Description**

A large number of functions for obtaining and modifying the parameters of likelihood models made in paleotree. These functions allow users to obtain or set parameter names, or obtain and set parameter bounds, both of which are treated as an attribute of the function class used by paleotree. In practice, this allows users to quickly obtain parameter names and upper and lower values for use in bounded optimizers, including reasonable starting values.

**Usage**

```r
parnames(x, ...)  
# S3 method for class 'paleotreeFunc'
parnames(x, ...)

# S3 method for class 'constrained'
parnames(x, ...)
parnames(x) <- value

# S3 replacement method for class 'constrained'
parnames(x) <- value

# S3 replacement method for class 'paleotreeFunc'
parnames(x) <- value
parbounds(x, ...)
```

```r
# S3 method for class 'paleotreeFunc'
parbounds(x, ...)
```
## S3 method for class 'constrained'
parbounds(x, ...)

parbounds(x) <- value

## S3 replacement method for class 'constrained'
parbounds(x) <- value

## S3 replacement method for class 'paleotreeFunc'
parbounds(x) <- value

parLower(x, ...)

## S3 method for class 'constrained'
parLower(x, ...)

## S3 method for class 'paleotreeFunc'
parLower(x, ...)

parLower(x) <- value

## S3 replacement method for class 'constrained'
parLower(x) <- value

## S3 replacement method for class 'paleotreeFunc'
parLower(x) <- value

parUpper(x, ...)

## S3 method for class 'constrained'
parUpper(x, ...)

## S3 method for class 'paleotreeFunc'
parUpper(x, ...)

parUpper(x) <- value

## S3 replacement method for class 'constrained'
parUpper(x) <- value

## S3 replacement method for class 'paleotreeFunc'
parUpper(x) <- value

parInit(x, ...)

## S3 method for class 'constrained'
parInit(x, ...)

## S3 method for class 'paleotreeFunc'
parInit(x, ...)
## S3 method for class 'paleotreeFunc'
parInit(x, ...)

### Arguments

- **x**: A function of S3 class `paleotreeFunc` with all necessary attributes expected of that class, which include parameter names and upper and lower bounds. As I have deliberately not exported the function which creates this class, it should be impossible for regular users to obtain such objects easily without using one of the `make` functions, which automatically output a function of the appropriate class and attributes.

- **...**: 'Ignored arguments to future methods' (i.e. for diversitree). Kept here only so `constrainParPaleo` is kept as close to the parent method in diversitree as possible.

- **value**: The new value with which to replace the parameter names or bounds. Must be a vector of the same length as the number of parameters. For `parbounds`, must be a list composed of two vectors.

### Details

Parameter names cannot be changed for a constrained function.

The `parInit` function calls the bounds for each parameter and gives a randomly selected value selected from a uniform distribution, using the parameter bounds for each parameter as the bounds on the uniform distribution. This uses a shorthand to quickly generate initial parameter values which are within the set bounds, for use in functions such as `optim`. The random sampling of initial values allows a user to quickly assess if initial parameter values affect the optimization by simply rerunning the function on new values. Infinite initial parameter values (resulting from infinite bounds) are discarded, and replaced with the lower bound value (assuming only upper bounds are infinite...). Some randomly selected initial parameter values may be too high (due to the liberal upper bounds I set for parameters in many of the likelihood functions) and thus users should always try slightly different values to see if the resulting maximum likelihood parameter values change.

As `parInit` depends on the upper and lower bounds attribute, no function is offered to allow it to be replaced (as there is nothing to replace!).

### Value

Returns the sought parameter names, bounds or initial values or (for the replacement methods) returns a modified function with the respective attributes altered.

### Author(s)

These functions are strongly based on or inspired by the `argnames` functions provided for handling models in Rich Fitzjohn’s library `diversitree`, but the functions presented here are derivations written by David Bapst.

### See Also

These model methods were introduced to interact with the new model framework introduced in `paleotree` v1.9, in particular to interface with `constrainParPaleo`. 
modifyTerminalBranches

Modify, Drop or Bind Terminal Branches of Various Types (Mainly for Paleontological Phylogenies)

Description

These functions modify terminal branches or drop certain terminal branches based on various criteria. DropZLB drops tip-taxa that are attached to the tree via zero-length terminal branches ("ZLBs").

Examples

#example with make_durationFreqCont
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1, nTotalTaxa = c(30,40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
rangesCont <- sampleRanges(taxa, r = 0.5)
likFun <- make_durationFreqCont(rangesCont)

#get parameter names
parnames(likFun)

#get the bounds for those parameters
parbounds(likFun)

#can also get these seperately
parLower(likFun)
parUpper(likFun)

#initial parameter values
parInit(likFun)  #arbitrary midway value between par bounds

#can then use these in optimizers, such as optim with L-BFGS-B
#see the example for make_durationFreqCont

#renaming parameter names
likFun2 <- likFun
parnames(likFun2) <- c("extRate","sampRate")
parnames(likFun2)
#test if reset correctly
parnames(likFun2) == c("extRate","sampRate")
#also works for constrained functions
constrainFun <- constrainParPaleo(likFun,q.1-r.1)
parnames(constrainFun)
#also modified the parameter bounds, see!
parbounds(constrainFun)
parInit(constrainFun)
#but cannot rename parameter for constrained function!
This is sometimes useful for paleo-trees, as various time-scaling methods often produce these ZLBs, taxa whose early appearance causes them to be functionally interpreted as ancestors in some time-scaling methods. Removing ZLBs is advised for analyses of diversification/diversity, as these will appear as simultaneous speciation/extinction events. Note this function only drops tips attached to a terminal zero-length branch; if you want to collapse internal zero-length branches, see the ape function `di2multi`.

**Usage**

```r
modifyTerminalBranches
dropZLB(tree)
dropExtinct(tree, tol = 0.01, ignore.root.time = FALSE)
dropExtant(tree, tol = 0.01)
addTermBranchLength(tree, addtime = 0.001)
fixRootTime(treeOrig, treeNew, consistentDepth = TRUE,
nodeAgeTransfer = TRUE)
dropPaleoTip(tree, ...)
bindPaleoTip(tree, tipLabel, nodeAttach = NULL, tipAge = NULL,
edgeLength = NULL, positionBelow = 0, noNegativeEdgeLength = TRUE)
```

**Arguments**

- `tree` A phylogeny as a phylo object. `dropPaleoTip` requires this input object to also have a `tree$root.time` element. If not provided for `bindPaleoTip`, then the `root.time` will be presumed to be such that the furthest tip from the root is at time = 0.
- `tol` Tolerance for determining modern age; used for distinguishing extinct from extant taxa. Tips which end within `tol` of the furthest distance from the root will be treated as 'extant' taxa for the purpose of keeping or dropping.
- `ignore.root.time` Ignore `tree$root.time` in calculating which tips are extinct? `tree$root.time` will still be adjusted, if the operation alters the `root.time`.
- `addtime` Extra amount of time to add to all terminal branch lengths.
- `treeOrig` A phylo object of a time-scaled phylogeny with a $root.time element.
- `treeNew` A phylo object containing a modified form of `treeOrig` (with no extra tip taxa added, but possibly with some tip taxa removed).
- `consistentDepth` A logical, either TRUE or FALSE. If TRUE (the default) the tree’s root-to-furthest-tip depth is tested to make sure this depth is not greater than the new $root.time appended to the output tree.
- `nodeAgeTransfer` A logical. If TRUE, the $root.time of the new tree is determined by comparing the clades of taxa between the two input trees. The new root age assigned is
the age of (1) the treeOrig clade that contains all taxa present in treeNew and, if the set of (1) contains multiple clades, (2) the clade in the first set that contains the fewest taxa not in treeNew. If FALSE, the root.time assigned to treeNew is the $root.time of treeOrig, adjusted based on the change in total tree depth between treeOrig and treeNew, as measured between the root and the first matching taxon in both trees. The later is how fixRootTime functioned by default prior to paleotree v2.3.

... additional arguments passed to dropPaleoTip are passed to drop.tip.

tipLabel A character string of length = 1 containing the name of the new tip to be added to tree.

nodeAttach Node or tip ID number (as given in tree$edge) at which to attach the new tip. See documentation of bindNtip for more details.

tipAge The age of the tip taxon added to the tree, in time before present (i.e. where present is 0), given in the same units as the edges of the tree are already scaled. Cannot be given if edgeLength is given.

edgeLength The new edgeNlength of the terminal branch this tip is connected to. Cannot be given if tipAge is given.

positionBelow The distance along the edge below the node to be attached to (given in codenodeAttach to add the new tip. Cannot be negative or greater than the length of the edge below nodeAttach.

noNegativeEdgeLength Return an error if a negative terminal edge length is calculated for the new tip.

Details

DropExtinct drops all terminal branches which end before the modern (i.e. extinct taxa). DropExtant drops all terminal branches which end at the modern (i.e. extant/still-living taxa). In both cases, the modern is defined based on tree$root.time if available, or the modern is inferred to be the point in time when the tip furthest from the root (the latest tip) terminates.

If the input tree has a $root.time element, as expected for most paleo-tree objects handled by this library, that $root.time is adjusted if the relative time of the root divergence changes when terminal branches are dropped. Adjusted root.times are only given if the input tree has root.times.

addTermBranchLength adds an amount equal to the argument addtime to the terminal branch lengths of the tree. If there is a $root.time element, this is increased by an amount equal to addtime. A negative amount can be input to reduce the length of terminal branches. However, if negative branch lengths are produced, the function fails and a warning is produced. This function does not call fixRootTime, so the root.time elements in the result tree may be nonsensical, particularly if negative amounts are input.

When a tree is modified, such as having tips dropped or branches extended, fixRootTime can be used to find the new $root.time. It is mainly used as a utility function called by a majority of the other functions discussed in this help file.

dropPaleoTip is a wrapper for ape's drop.tip which also modifies the $root.time element if necessary, using fixRootTime. Similarly, bindPaleoTip is a wrapper for phytool's bind.tip which allows tip age as input and modifies the $root.time element if necessary (i.e. if a tip is added to edge leading up to the root).
Note that for bindPaleoTip, tips added below the root are subtracted from any existing $root.edge element, as per behavior of link(bind.tip) and bind.tree. However, bindPaleoTip will append a $root.edge of the appropriate value (i.e., root edge length) if one does not exist (or is not long enough) to avoid an error. After binding is finished, any $root.edge equal to 0 is removed before the resulting tree is output.

Value

Gives back a modified phylogeny as a phylo object, generally with a modified $root.time element.

Author(s)

David W. Bapst. The functions dropTipPaleo and bindTipPaleo are modified imports of drop.tip and bind.tip from packages ape and phytools.

See Also

compareTermBranches, phylodiv, drop.tip, bind.tip

Examples

```r
set.seed(444)
#Simulate some fossil ranges with simFossilRecord
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(30,40), nExtant = 0)
taxa <- fossilRecord2fossilTaxa(record)
#simulate a fossil record with imperfect sampling with sampleRanges
rangesCont <- sampleRanges(taxa,r = 0.5)
#Now let's make a tree using taxa2phylo
tree <- taxa2phylo(taxa,obs_time = rangesCont[,2])
#compare the two trees
layout(1:2)
plot(Ladderize(tree))
plot(Ladderize(dropZLB(tree)))
layout(1)

#example using dropExtinct and dropExtant
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(30,40), nExtant = c(10,20))
taxa <- fossilRecord2fossilTaxa(record)
tree <- taxa2phylo(taxa)
phylodiv(tree)
tree1 <- dropExtinct(tree)
phylodiv(tree1)
tree2 <- dropExtant(tree)
phylodiv(tree2)

#graphics.off()

#example using addTermBranchLength
```
```r
set.seed(444)
treeA <- rtree(10)
treeB <- addTermBranchLength(treeA, 1)
compareTermBranches(treeA, treeB)

# test dropPaleoTip
# (and fixRootTime by extension...)

tree <- read.tree(text = "(A:3,(B:2,(C:5,D:3):2):3);")
tree$root.time <- 10
plot(tree, no.margin = FALSE)
axisPhylo()

# now a series of tests, dropping various tips
(test <- dropPaleoTip(tree, "A")$root.time) # = 7
(test[2] <- dropPaleoTip(tree, "B")$root.time) # = 10
(test[3] <- dropPaleoTip(tree, "C")$root.time) # = 10
(test[4] <- dropPaleoTip(tree, "D")$root.time) # = 10
(test[5] <- dropPaleoTip(tree, c("A", "B"))$root.time) # = 5
(test[6] <- dropPaleoTip(tree, c("B", "C"))$root.time) # = 10
(test[7] <- dropPaleoTip(tree, c("A", "C"))$root.time) # = 7
(test[8] <- dropPaleoTip(tree, c("A", "D"))$root.time) # = 7

# is it all good? if not, fail so paleotree fails...
if(!identical(test, c(7,10,10,10,5,10,7,7))) {stop("fixRootTime fails!")}

# simple example
require(phytools)

# bindPaleoTip effectively wraps bind.tip from phytools
# using a conversion like below

tipAge <- 5
node <- 6

# new length = the root time - tipAge - nodeheight(tree, node)
newLength <- tree$root.time - tipAge - nodeheight(tree, node)
tree1 <- bind.tip(tree, "tip.label", where = node, edge.length = newLength)
```
modifyTerminalBranches

layout(1:2)
plot(tree); axisPhylo()
plot(tree1); axisPhylo()

## End (Not run)

# now with bindPaleoTip

tree1 <- bindPaleoTip(tree, "new", nodeAttach = 6, tipAge = 5)

layout(1:2)
plot(tree); axisPhylo()
plot(tree1); axisPhylo()
layout(1)

# then the tip age of "new" should be 5
test <- dateNodes(tree1)[which(tree1$tip.label == "new")][1] == 5
if(!test){stop("bindPaleoTip fails!")}

# with positionBelow

tree1 <- bindPaleoTip(tree, "new", nodeAttach = 6, tipAge = 5, positionBelow = 1)

layout(1:2)
plot(tree); axisPhylo()
plot(tree1); axisPhylo()
layout(1)

# at the root

tree1 <- bindPaleoTip(tree, "new", nodeAttach = 5, tipAge = 5)

layout(1:2)
plot(tree); axisPhylo()
plot(tree1); axisPhylo()
layout(1)

# then the tip age of "new" should be 5
test <- dateNodes(tree1)[which(tree1$tip.label == "new")][1] == 5
if(!test){stop("bindPaleoTip fails!")}

# at the root with positionBelow

tree1 <- bindPaleoTip(tree, "new", nodeAttach = 5, tipAge = 5, positionBelow = 3)

layout(1:2)
plot(tree); axisPhylo()
plot(tree1); axisPhylo()
layout(1)
multiDiv

Calculating Diversity Curves Across Multiple Datasets

Description

Calculates multiple diversity curves from a list of datasets of taxic ranges and/or phylogenetic trees, for the same intervals, for all the individual datasets. A median curve with 95 percent quantile bounds is also calculated and plotted for each interval.

Usage

multiDiv(data, int.length = 1, plot = TRUE, split.int = TRUE, drop.ZLB = TRUE, drop.cryptic = FALSE, extant.adjust = 0.01, plotLogRich = FALSE, yAxisLims = NULL, timelims = NULL, int.times = NULL, plotMultCurves = FALSE, multRainbow = TRUE, divPalette = NULL, divLineType = 1, main = NULL)

plotMultiDiv(results, plotLogRich = FALSE, timelims = NULL, yAxisLims = NULL, plotMultCurves = FALSE, multRainbow = TRUE, divPalette = NULL, divLineType = 1, main = NULL)

Arguments

data A list where each element is a dataset, formatted to be input in one of the diversity curve functions listed in DiversityCurves.

int.length The length of intervals used to make the diversity curve. Ignored if int.times is given.

plot If TRUE, the median diversity curve is plotted.

split.int For discrete time data, should calculated/input intervals be split at discrete time interval boundaries? If FALSE, can create apparent artifacts in calculating the diversity curve. See below.

drop.ZLB If true, zero-length terminal branches are dropped from the input tree for phylogenetic datasets, before calculating standing diversity.

drop.cryptic If TRUE, cryptic taxa are merged to form one taxon for estimating taxon curves. Only works for objects from simFossilRecord via fossilRecord2fossilTaxa.

extant.adjust Amount of time to be added to extend start time for (0,0) bins for extant taxa, so that the that 'time interval' does not appear to have an infinitely small width.

plotLogRich If TRUE, taxic diversity is plotted on log scale.
multiDiv

yAxisLims  Limits for the y (i.e. richness) axis on the plotted diversity curves. Only affects plotting. Given as either NULL (the default) or as a vector of length two as for xlim in the basic R function plot. Time axes will be plotted exactly to these values. The minimum value must be more than 1 if plotLogRich = TRUE.

timelims  Limits for the x (time) axis for diversity curve plots. Only affects plotting. Given as either NULL (the default) or as a vector of length two as for xlim in the basic R function plot. Time axes will be plotted exactly to these values.

int.times  An optional two-column matrix of the interval start and end times for calculating the diversity curve. If NULL, calculated internally. If given, the argument split.int and int.length are ignored.

plotMultCurves  If TRUE, each individual diversity curve is plotted rather than the median diversity curve and 95 percent quantiles. FALSE by default.

multRainbow  If TRUE and plotMultCurves are both TRUE, each line is plotted as a different, randomized color using the function 'rainbow'. If FALSE, each line is plotted as a black line. This argument is ignored if divPalette is supplied.

divPalette  Can be used so users can pass a vector of chosen color identifiers for each diversity curve in 'data' which will take precedence over multRainbow. Must be the same length as the number of diversity curves supplied.

divLineType  Used to determine line type (lty) of the diversity curves plotted when plotMultCurves = TRUE. Default is lty = 1 for all curves. Must be either length of 1 or exact length as number of diversity curves.

main  The main label for the figure.

results  The output of a previous run of multiDiv for reploting.

Details

This function is essentially a wrapper for the individual diversity curve functions included in paleotree. multiDiv will intuitively decide whether input datasets are continuous-time taxic ranges, discrete-time (binned interval) taxic ranges or phylogenetic trees, as long as they are formatted as required by the respective diversity curve functions. A list that contains a mix of data types is entirely acceptable. A list of matrices output from fossilRecord2fossilTaxa, via simulation with simFossilRecord is allowable, and treated as input for taxicDivCont. Data of an unknown type gives back an error.

The argument split.int splits intervals, if and only if discrete interval time data is included among the datasets. See the help file for taxicDivDisc to see an explanation of why split.int = TRUE by default is probably a good thing.

As with many functions in the paleotree library, absolute time is always decreasing, i.e. the present day is zero.

The 'averaged' curve is actually the median rather than the mean as diversity counts are often highly skewed (in this author's experience).

The shaded certainty region around the median curve is the two-tailed 95 percent lower and upper quantiles, calculated from the observed data. It is not a true probabilisitic confidence interval, as it has no relationship to the standard error.
Value

A list composed of three elements will be invisibly returned:

- `int.times` A two column matrix giving interval start and end times
- `div` A matrix of measured diversities in particular intervals by rows, with each column representing a different dataset included in the input
- `median.curve` A three column matrix, where the first column is the calculated median curve and the second and third columns are the 95 percent quantile upper and lower bounds

See Also

The diversity curve functions used include: `phyloDiv`, `taxicDivCont` and `taxicDivDisc`. Also see the function LTT.average.root in the package TreeSim, which calculates an average LTT curve for multiple phylogenies, the functions `codemltt.plot` in ape and `codeltt` in phytools.

Examples

```r
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(30,40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
rangesCont <- sampleRanges(taxa, r = 0.5)
rangesDisc <- binTimeData(rangesCont, int.length = 1)
cladogram <- taxa2Cladogram(taxa, plot = TRUE)
#using multiDiv with very different data types
ttree <- timePaleoPhy(cladogram, rangesCont, type = "basic", add.term = TRUE, plot = FALSE)
input <- list(rangesCont, rangesDisc, ttree)
multiDiv(input, plot = TRUE)

#using fixed interval times
multiDiv(input, int.times = rangesDisc[[1]], plot = TRUE)

#using multiDiv with samples of trees
ttrees <- timePaleoPhy(cladogram, rangesCont, type = "basic", randres = TRUE, ntrees = 10, add.term = TRUE)
multiDiv(ttrees)

#uncertainty in diversity history is solely due to
#the random resolution of polytomies

#multiDiv can also take output from simFossilRecord, via fossilRecord2FossilTaxa
#what do many simulations run under some set of conditions 'look' like on average?
set.seed(444)
records <- simFossilRecord(p = 0.1, q = 0.1, nruns = 10,
totalTime = 30, plot = TRUE)
taxa <- sapply(records, fossilRecord2FossilTaxa)
multiDiv(taxa)

#increasing cone of diversity!
#Even better on a log scale:
multiDiv(taxa, plotLogRich = TRUE)
```
nearestNeighborDist  Nearest Neighbor Distances for Morphological Disparity Studies

Description

This is a simple function for obtaining nearest neighbor distance from a symmetric pair-wises distance matrix, assumed here to be dissimilarities between pairs of taxa. Per-species NND is returned rather than a mean or other summary value.

Usage

nearestNeighborDist(distMat)

Arguments

distMat  A symmetric, square pair-wise distance matrix, assumed to be a dissimilarity matrix with a zero diagonal. Can be a 'dist' object rather than a numerical matrix. Taxon labels can be applied to the rows and columns (or as labels if a 'dist' object) and will be used to name the resulting output.

Details

This function is mainly included here for pedagogical (teaching) purposes. NND is so simple to calculate, users are urged to write their own functions for primary research purposes.

Typically, the mean NND for a group is reported and used to compare different groupings of taxa (such as different time intervals, or different clades). Bootstrapping should be used to generate confidence intervals.

Value

Returns a vector of the nearest neighbor distance for each unit (taxon) in the pair-wise distance matrix, named with the labels from the input distance matrix.
**nearestNeighborDist**

**Author(s)**

David W. Bapst

**References**


**See Also**

For the example dataset used in examples, see *graptDisparity*

**Examples**

```r
# example using graptolite disparity data from Bapst et al. 2012

# load data
data(graptDisparity)

# calculate mean NND
NND <- nearestNeighborDist(graptDistMat)
mean(NND)

# calculate NND for different groups

groupID <- graptCharMatrix[,54]+1

groupNND <- numeric(7)
names(groupNND) <- c("Normalo.","Monogr.","Climaco.",
                     "Dicrano.","Lasiogr.","Diplogr.","Retiol.")
for(i in unique(groupID)){
groupNND[i] <- mean(nearestNeighborDist(
graptDistMat[groupID == i,groupID == i]))
}

# the paraphyletic Normalograptids that survived the HME are most clustered
# but this looks at all the species at once
# and doesn't look for the nearest *co-extant* neighbor!
# need to bring in temporal info to test that
```
nodeDates2branchLengths

Obtaining Edge Lengths for Undated Phylogenies Using Known Branching Node and Tip Ages

Description

This function takes some undated phylogenetic topology, a set of ages in absolute time, for the internal nodes and (by default) the terminal tips of that phylogeny, and returns a dated phylogeny consistent with those input ages.

Usage

nodeDates2branchLengths(nodeDates, tree, allTipsModern = FALSE)

Arguments

- **nodeDates**: Under default `allTipsModern = FALSE` conditions, `nodeDates` should be a vector of length `Ntip(tree) + Nnode(tree)` which contains the dates for all terminal tip nodes and internal nodes for the tree, in that order, as numbered in the `tree$edge` matrix. Such a vector is produced as output by `dateNodes`. If `allTipsModern = TRUE`, then the vector should only be as long as the number of nodes, and contain the dates only for those same internal nodes in `tree`. These dates should always on a descending scale (i.e. time before present), with respect to an absolute time-scale. It is possible for the time 0 date to represent a date far in the future from the latest tip.

- **tree**: An undated phylogeny object, of class `phylo`, lacking edge lengths. If the tree appears to be dated (i.e. has edge lengths), the function will issue a warning.

- **allTipsModern**: A logical, default is `FALSE`. If `FALSE`, then the function expects `nodeDates` to contain ages for all 'nodes' - both internal branching nodes and terminal tips. If `TRUE`, then the function will expect `nodeDates` to contain ages only for internal branching nodes, and all tips will be assumed to be at time 0. (Thus, if your tree is ultrametric but tips aren’t all at the modern, do not use `allTipsModern = TRUE`).

Details

The function `compute.brtimes` in package `ape` does a very similar functionality, but is limited in its application for only ultrametric trees, as it does not allow for tips to have incongruent ages. It also only accepts node ages as on the relative scale where the latest tips are at zero, as assumed in general elsewhere in package `ape`.

Value

A dated tree as a list of class `phylo`, with a `$root.time` element for referencing the tree against absolute time.
Author(s)

David W. Bapst

See Also

This function will likely often be used in conjunction with dateNodes, such as for summarizing node and tip age estimates from a sample of trees, to produce a single dated tree to act as a point estimate. Beware however that point estimates of tree samples may have little resemblance to any individual tree in that sample.

This function should perform identically for ultrametric trees as package ape’s function compute.btime.

Examples

```r
set.seed(444)

# we'll do a number of tests, let's check at the end that all are TRUE
tests <- logical()

# with a non-ultrametric tree
chrono <- rtree(10)
# make an undated tree
notChrono <- chrono
notChrono$edge.length <- NULL

# now lets try with dateNodes in paleoTree
nodeTimes <- dateNodes(chrono)
# need to use allTipsModern = FALSE because tip ages are included
chronoRedux <- nodeDates2branchLengths(tree = notChrono,
                                        nodeDates = nodeTimes,
                                        allTipsModern = FALSE)
# test that its the same
(tests <- c(tests, all.equal.numeric(chrono$edge.length, chronoRedux$edge.length)))

# modern ultrametric tree
chrono <- rcoal(10)
# make an undated tree
notChrono <- chrono
notChrono$edge.length <- NULL

# with ultrametric trees, you could just use ape's compute.btime

# getting branching times with ape
branchingTimes <- branching.times(chrono)
# setting those branching times with ape
chronoRedux <- compute.btime(notChrono, branchingTimes)
# test that its the same
(tests <- c(tests, all.equal.numeric(chrono$edge.length, chronoRedux$edge.length)))

# lets do the same thing but with nodeDates2branchLengths

# can use branching.times from ape
```
# (but only for ultrametric trees!)
chronoRedux <- nodeDates2branchLengths(tree = notChrono,
    nodeDates = branchingTimes, allTipsModern = TRUE)
# test that its the same
(tests <- c(tests, all.equal.numeric(chrono$edge.length, chronoRedux$edge.length)))

# now lets try with dateNodes in paleotree
nodeTimes <- dateNodes(chrono)
# need to use allTipsModern = FALSE because tip ages are included
chronoRedux <- nodeDates2branchLengths(tree = notChrono,
    nodeDates = nodeTimes, allTipsModern = FALSE)
# test that its the same
(tests <- c(tests, all.equal.numeric(chrono$edge.length, chronoRedux$edge.length)))

# get just the node times (remove tip dates)
nodeOnlyTimes <- nodeTimes[-(1:Ntip(chrono))]
# let's use the allTipsModern = TRUE setting
chronoRedux <- nodeDates2branchLengths(tree = notChrono,
    nodeDates = nodeOnlyTimes, allTipsModern = TRUE)
# test that its the same
(tests <- c(tests, all.equal.numeric(chrono$edge.length, chronoRedux$edge.length)))

# did all tests come out as TRUE?
if(!all(tests))(stop("nodeDates2branchLengths isn't functioning correctly"))

---

**obtainDatedPosteriorTreesMrB**

*Get the Sample of Posterior Trees from a Dated Phylogenetic Analysis with MrBayes (Or a Summary Tree, such as the MCCT)*

**Description**

MrBayes is not great for getting samples of dated posterior phylogenies, or for obtaining certain summary trees from the posterior (specifically the MCCT and MAP, which are specific trees in the posterior). This is because the tree samples as returned are scaled relative to rate parameters in a separate file. This function attempts to automate the handling of multiple files (both .t tree files and .p parameter files), as well as multiple files associated with separate runs, to obtain samples of posterior trees, or summary trees such as the MCCT or MAP. These resulting trees are now scaled to units of time, but not be placed correctly on an absolute time-scale if all tips are extinct. See details of output below.

**Usage**

obtainDatedPosteriorTreesMrB(runFile, nRuns = 2, burnin = 0.5, outputTrees,
    labelPostProb = FALSE, getFixedTimes = FALSE, getRootAges = FALSE,
    originalNexusFile = NULL, file = NULL)
Arguments

runFile  A filename in the current directory, or a path to a file that is either a .p or .t file from a MrBayes analysis. This filename and path will be used for finding additional .t and .p files, via the nRuns settings and assuming that files are in the same directory and these files are named under typical MrBayes file naming conventions. (In other words, if you have renamed your .p or .t files, this function probably won’t be able to find them.)

nRuns The number of runs in your analysis. This variable is used for figuring out what filenames will be searched for: if you specify that you have less runs than you actually ran in reality, then some runs won’t be examined in this function. Conversely, specify too many, and this function will throw an error when it cannot find files it expects but do not exist. The default for this argument (two runs) is based on the default number of runs in MrBayes.

burnin The fraction of trees sampled in the posterior discarded and not returned by this function directly, nor included in calculation of summary trees. Must be a numeric value greater than 0 and less than 1.

outputTrees Determines the output trees produced; for format of output, see section on returned Value below. Must be of length one, and either “all”, which means all trees from the post-burnin posterior will returned, a number greater than zero, which will be the number of trees randomly sampled from across the post-burning posterior and returned, or ”MCCT” and ”MAP”, which stand for ’maximum clade compatibility tree’ and ’maximum a posteri tree’ respectively. The MAP is the single tree from the post-burnin posterior with the highest marginal likelihood. The MCCT is the single tree from the post-burnin posterior which contains clades with the highest product of posterior probabilities for its component clades. Thus, the MAP is the best overall tree, while the MCCT may be the best tree for summarizing topological support.

labelPostProb Logical. If TRUE, then nodes of the output tree will be labeled with their respective posterior probabilities, as calculated based on the frequency of a clade occurring across the post-burnin posterior tree sample. If FALSE, this is skipped.

getFixedTimes If TRUE, this function will also look for, scan, and parse an associated NEXUS file. Ignoring any commented lines (i.e. anything between ”[]”), commands for fixing taxa will be identified, parsed and returned to the user, either as a message printed to the R console if output is read to a file, or as a attribute named ’fixed ages’ if output as an R object (formatted as a two-column table of OTU names and their respective fixed ages). If the output is an R object, these objects with Please note: the code for getFixedTimes = TRUE contains a while() loop in it for removing nested series of square brackets (i.e. treated as comments in NEXUS files). Thus files with ridiculously nested series of brackets may cause this code to take a while to complete, or may even cause it to hang.

getRootAges FALSE by default. If TRUE, and getFixedTimes = TRUE as well as file = NULL (such that trees will be assigned within the R memory rather than saved to an external file), the functions setRootAge and its wrapper function setRootAges will be applied to the output so that all output trees have root.time elements for use with other functions in paleotree as well as other packages.
Details

This function is most useful for dealing with dating analyses in MrBayes, particularly when tip-dating a tree with fossil taxa, as the half-compatibility and all-compatibility summary trees offered by the 'sumt' command in MrBayes can have issues properly portraying summary trees from such datasets.

Value

Depending on argument file, the output tree or trees is either returned directly, or instead written out in NEXUS format via ape's write.NEXUS function to an external file. The output will consist either of multiple trees sampled from the post-burnin posterior, or will consist of a single phylogeny (a summary tree, either the MCCT or the MAP).

If the argument setRootAges = TRUE is not used, users are warned that the resulting dated trees will not have $root.time elements necessary for comparison against an absolute time-scale. While the trees may be scaled to units of absolute time now, rather than with branch lengths expressed in the rate of character change, the dates estimated by some phylogenetics functions in R may give inaccurate estimates of when events occur on the absolute time-scale if all tips are extinct. This is because most functions for phylogenetics in R (and elsewhere) will instead presume that the latest tip will be at time 0 (the modern), which may be wrong if you are using paleotree for analyzing paleontological datasets consisting of entirely extinct taxa. This can be solved by using argument getFixedTimes = TRUE to obtain fixed tip ages, and then scaling the resulting output to absolute time using the argument setRootAges = TRUE, which obtains a $root.time element for each tree using the functions setRootAge and setRootAges (for single and multiple phylogenies).

Author(s)

David Bapst, with rescaling of raw output trees via code originally written by Nicholas Crouch.

See Also

When the arguments getFixedTimes = TRUE and setRootAges = TRUE are used, the resulting output will be scaled to absolute time with the available fixed ages using functions setRootAge and setRootAges (for single and multiple phylogenies). This is only done if fixed ages are available and if the tree is not being saved to an external file.

Maximum Clade Credibility trees are estimated using the function maxCladeCred in package phangorn.
Examples

```r
## Not run:

MCCT <- obtainDatedPosteriorTreesMrB(
  runfile = "C:\myTipDatingAnalysis\MrB_run_fossil_05-10-17.nex.run1.t",
  nRuns = 2, burnin = 0.5,
  outputTrees = "MCCT", file = NULL)

MAP <- obtainDatedPosteriorTreesMrB(
  runfile = "C:\myTipDatingAnalysis\MrB_run_fossil_05-10-17.nex.run1.t",
  nRuns = 2, burnin = 0.5, getFixedTimes = TRUE,
  outputTrees = "MAP", file = NULL)

# get a root age from the fixed ages for tips
getRootAge(tree = MAP)

# pull a hundred trees randomly from the posterior
hundredRandomlySelectedTrees <- obtainDatedPosteriorTreesMrB(
  runfile = "C:\myTipDatingAnalysis\MrB_run_fossil_05-10-17.nex.run1.t",
  nRuns = 2, burnin = 0.5, getFixedTimes = TRUE,
  getRootAges = TRUE,
  outputTrees = 100, file = NULL)

## End(Not run)
```

### occData2timeList

**Converting Occurrences Data to a timeList Data Object**

**Description**

This function converts occurrence data, given as a list where each element is a different taxon’s occurrence table (containing minimum and maximum ages for each occurrence), to the 'timeList' format, consisting of a list composed of a matrix of lower and upper age bounds for intervals, and a second matrix recording the interval in which taxa first and last occur in the given dataset.

**Usage**

```r
occData2timeList(occList, intervalType = "dateRange")
```

**Arguments**

- **occList**
  A list where every element is a table of occurrence data for a different taxon, such as that returned by `taxonSortPBDBocc`. The occurrence data can be either a two-column matrix composed of the lower and upper age bounds on each taxon occurrence, or has two named variables which match any of the field names given by the PBDB API under either the 'pbdb' vocab or 'com' (compact) vocab for early and late age bounds.
intervalType Must be either "dateRange" (the default), "occRange" or "zoneOverlap". Please see details below.

Details

This function should translate taxon-sorted occurrence data, which could be Paleobiology Database datasets sorted by taxonSortPDBocc or any data object where occurrence data (i.e. age bounds for each occurrence) for different taxa is separated into different elements of a named list.

The argument intervalType:

The argument intervalType controls the algorithm used for obtain first and last interval bounds for each taxon, of which there are several to select from:

"dateRange" The default option. The bounds on the first appearances are the span between the oldest upper and lower bounds of the occurrences, and the bounds on the last appearances are the span between the youngest upper and lower bounds across all occurrences. This is guaranteed to provide the smallest bounds on the first and last appearances, and was originally suggested to the author by J. Marcot.

"occRange" This option returns the smallest bounds among (a) the oldest occurrences for the first appearance (i.e. all occurrences with their lowest bound at the oldest lower age bound), and (b) the youngest occurrences for the last appearance (i.e. all occurrences with their uppermost bound at the youngest upper age bound).

"zoneOverlap" This option is an attempt to mimic the stratigraphic range algorithm used by PBDB Classic which "finds the oldest base that is older than at least part of all the intervals and the youngest that is younger than at least part of all the intervals" (personal communication, J. Alroy). This is a somewhat more complex case as we are trying to obtain a timeList object. So, for calculating the bounds of the first interval a taxon occurs in, the zoneOverlap algorithm looks for all occurrences that overlap with the age range of the earliest-most occurrence and (1) obtains their earliest boundary ages and returns the latest-most earliest age boundary among these overlapping occurrences and (2) obtains their latest boundary ages and returns the earliest-most latest age boundary among these overlapping occurrences. Similarly, for calculating the bound of the last interval a taxon occurs in, the zoneOverlap algorithm looks for all occurrences that overlap with the age range of the latest-most occurrence and (1) obtains their earliest boundary ages and returns the latest-most earliest age boundary among these overlapping occurrences and (2) obtains their latest boundary ages and returns the earliest-most latest age boundary among these overlapping occurrences.

On theoretical grounds, one could probably describe the zone-of-overlap algorithm as minimizing taxonomic age ranges by assuming that all overlapping occurrences at the start and end of a taxon’s range probably describe a very similar first and last appearance (FADs and LADs), and thus picks the occurrence with bounds that extends the taxonomic range the least. However, this does come with a downside that if these occurrences are not essentially repeated attempts to capture the same FAD or LAD, then the zone-of-overlap algorithm is not an accurate depiction of the uncertainty in the ages. The true biological range of a taxon might be well outside the bounds obtained using the zone-of-overlap algorithm. A more conservative approach is the "dateRange" algorithm which finds the smallest possible bounds on the endpoints of a taxon’s range without ignoring uncertainty from any particular set of occurrences.
Value

Returns a standard timeList data object, as used by many other paleotree functions, like `bin_timePaleoPhy`, `bin_cal3TimePaleoPhy` and `taxicDivDisc`

Author(s)

David W. Bapst, with the 'dateRange' algorithm suggested by Jon Marcot.

See Also

taxonSortPBDBocc, plotOccData and the example graptolite dataset at graptPBDB

Examples

data(graptPBDB)

graptOccSpecies <- taxonSortPBDBocc(graptOccPBDB, rank = "species", onlyFormal = FALSE)
graptTimeSpecies <- occData2timeList(occList = graptOccSpecies)

head(graptTimeSpecies[[1]])
head(graptTimeSpecies[[2]])

graptOccGenus <- taxonSortPBDBocc(graptOccPBDB, rank = "genus", onlyFormal = FALSE)
graptTimeGenus <- occData2timeList(occList = graptOccGenus)

layout(1:2)
taxicDivDisc(graptTimeSpecies)
taxicDivDisc(graptTimeGenus)

# the default interval calculation is "dateRange"
# let's compare to the other option, "occRange"
# for species

graptOccRange <- occData2timeList(occList = graptOccSpecies, intervalType = "occRange")

# we would expect no change in the diversity curve
# because there are only changes in th
# earlist bound for the FAD
# latest bound for the LAD
# so if we are depicting ranges within maximal bounds
# dateRanges has no effect
layout(1:2)
taxicDivDisc(graptTimeSpecies)
taxicDivDisc(graptOccRange)
# yep, identical

# so how much uncertainty was gained by using dateRange?

# write a simple function for getting uncertainty in first and last
# appearance dates from a timeList object
sumAgeUncert <- function(timeList){
  fourDate <- timeList$FourDate(timeList)
perOcc <- (fourDate[,1]-fourDate[,2])+(fourDate[,3]-fourDate[,4])
sum(perOcc)
}

# total amount of uncertainty in occRange dataset
sumAgeUncert(graptOccRange)
# total amount of uncertainty in dateRange dataset
sumAgeUncert(graptTimeSpecies)
# the difference
sumAgeUncert(graptOccRange)-sumAgeUncert(graptTimeSpecies)
# as a proportion
1-(sumAgeUncert(graptTimeSpecies)/sumAgeUncert(graptOccRange))

# a different way of doing it
dateChange <- timeList2fourDate(graptTimeSpecies)-timeList2fourDate(graptOccRange)
apply(dateChange,2,sum)
# total amount of uncertainty removed by dateRange algorithm
sum(abs(dateChange))

layout(1)

optimPaleo

Simplified Optimizer for paleoTree Likelihood Functions

Description

This function is a deliberately simplistic automation wrapper for the function `optim` and the use of the "L-BFGS-B" optimizing method, with initial parameter values and bounds provided with `parInit`, `parLower` and `parUpper`. It is mainly provided here as a shorthand to be used in educational demonstrations where model-fitting is not the primary focus, and use in actual analyses should be avoided.

Usage

optimPaleo(modelFun)

Arguments

modelFun A likelihood function for a model, of class 'paleoTreeFunc'.

Details

This is mainly provided in this publicly released package for pedagogical reasons. Users seeking an optimizer for their own analytical purposes should write their own optim function.

Value

Returns the results from using optim.
See Also

constrainParPaleo and modelMethods

Examples

# This function simply replicates optim() as shown below
# where modelFun is the likelihood function

#optim(parInit(modelFun),modelFun,
# lower = parLower(modelFun),upper = parUpper(modelFun),
# method = "L-BFGS-B",control = list(maxit = 1000000))

parentChild2taxonTree  Create a Taxonomy-Based Phylogeny ('Taxon Tree') from a Table of Parent-Child Taxon Relationships

Description

This function takes a two-column matrix of taxon names, indicating a set of binary parent-taxon:child-taxon paired relationships with a common root, and returns a 'taxonomy-tree' phylogeny object of class 'phylo'.

Usage

parentChild2taxonTree(parentChild, tipSet = "nonParents", cleanTree = TRUE)

Arguments

parentChild  A two-column matrix of type character where each element is a taxon name. Each row represents a parent-child relationship with first the parent (column 1) taxon name and then the child (column 2).

tipSet  This argument controls which taxa are selected as tip taxa for the output tree. The default tipSet = "nonParents" selects all child taxa which are not listed as parents in parentChild. Alternatively, tipSet = "all" will add a tip to every internal node with the parent-taxon name encapsulated in parentheses.

cleanTree  By default, the tree is run through a series of post-processing, including having singles collapsed, nodes reordered and being written out as a Newick string and read back in, to ensure functionality with ape functions and ape-derived functions. If FALSE, none of this post-processing is done and users should beware, as such trees can lead to hard-crashes of R.

Details

All taxa listed must be traceable via their parent-child relationships to a single, common ancestor which will act as the root node for output phylogeny. Additionally, the root used will be the parent taxon to all tip taxa closest in terms of parent-child relationships to the tip taxa: i.e., the most recent common ancestor. Ancestral taxa which are singular internal nodes that trace to this root are removed, and a message is printed.
Value
A phylogeny of class ‘phylo’, with tip taxa as controlled by argument tipSet. The output tree is returned with no edge lengths.
The names of higher taxa than the tips should be appended as the element $node.label for the internal nodes.

Author(s)
David W. Bapst

See Also
makePBDBtaxontree, taxontable2taxontree

Examples

# let's create a small, really cheesy example
pokexample <- rbind(cbind("Squirtadae",c("Squirtle","Blastoise","Wartortle")),
c("Shelloidea","Lapras"),c("Shelloidea","Squirtadae"),
c("Pokezooa","Shelloidea"),c("Pokezooa","Parasect"),
c("Rodentapokemonmorpha","Linoone"),c("Rodentapokemonmorpha","Sandshrew"),
c("Rodentapokemonmorpha","Pikachu"),c("Hirsutamona","Ursaring"),
c("Hirsutamona","Rodentapokemonmorpha"),c("Pokezooa","Hirsutamona"))

# Default: tipSet = ‘nonParents’
pokeTree <- parentChild2taxonTree(pokexample, tipSet = "nonParents")
plot(pokeTree); nodelabels(pokeTree$node.label)

# Get ALL taxa as tips with tipSet = ‘all’
pokeTree <- parentChild2taxonTree(pokexample, tipSet = "all")
plot(pokeTree); nodelabels(pokeTree$node.label)

## Not run:

# let's try a dataset where not all the taxon relationships lead to a common root
pokexample_bad <- rbind(cbind("Squirtadae",c("Squirtle","Blastoise","Wartortle")),
c("Shelloidea","Lapras"),c("Shelloidea","Squirtadae"),
c("Pokezooa","Shelloidea"),c("Pokezooa","Parasect"),
c("Rodentapokemonmorpha","Linoone"),c("Rodentapokemonmorpha","Sandshrew"),
c("Rodentapokemonmorpha","Pikachu"),c("Hirsutamona","Ursaring"),
c("Hirsutamona","Rodentapokemonmorpha"),c("Pokezooa","Hirsutamona"),
c("Umbrarcheota","Gengar"))

# this should return an error, as Gengar doesn't share common root
pokeTree <- parentChild2taxonTree(pokexample_bad)

# another example, where a taxon is listed as both parent and child
perCapitaRates

Description

Calculates and plots per-capita origination and extinction rates from sequential discrete-time taxon ranges, following Foote (2000).

Usage

perCapitaRates(timeList, plot = TRUE, logRates = FALSE,
               drop.extant = FALSE, isExtant = NULL, jitter = TRUE,
               legendPosition = "topleft")

Arguments

timeList     A list composed of two matrices, giving interval start and end dates and taxon first and last occurrences within those intervals. See details.
plot         If TRUE, the per-capita origination and extinctions rates are plotted for each interval. Rates which cannot be calculated for an interval will not be plotted, thus appearing as a gap in the plotted graph. The author takes no responsibility for the aesthetics of this plot.
logRates  If TRUE, rates plotted on log scale.
drop.extant  Drops all extant taxa from a dataset before calculating per-capita origination and extinction rates.
isExtant  A vector of TRUE and FALSE values, same length as the number of taxa in the second matrix of timeList, where TRUE values indicate taxa that are alive in the modern day (and thus are boundary crossers which leave the most recent interval). By default, this argument is NULL and instead which taxa are extant is inferred based on which taxa occur in an interval with start and end times both equal to zero. See details.
jitter  If TRUE (default) the extinction rate will be plotted slightly ahead of the origination rate on the time axis, so the two can be differentiated.
legendPosition  The position of a legend indicating which line is origination rate and which is extinction rate on the resulting plot. This is given as the possible positions for argument 'x' of the function legend, and by default is "topleft", which will be generally useful if origination and extinction rates are initially low. If legendPosition = NA, then a legend will not be plotted.

Details
This function calculates the per-capita rates of taxonomic origination and extinction from paleontological range data, as described by Foote (2000). These values are the instantaneous rate of either type of event occurring per lineage time-units. Although Foote (2001) also presents a number of alternative rates collected from the prior literature such as the 'Van Valen' rate metrics, these are not implemented here, but could be estimated using the matrix invisibly output by this function (See Foote, 2000, for the relevant equations for calculating these).

The timeList object should be a list composed of two matrices, the first matrix giving by-interval start and end times (in absolute time), the second matrix giving the by-taxon first and last appearances in the intervals defined in the first matrix, numbered as the rows. Absolute time should be decreasing, while the intervals should be numbered so that the number increases with time. Taxa alive in the modern should be either (a) listed in isExtant or (b) listed as last occurring in a time interval that begins at time 0 and ends at time 0. See the documentation for the time-scaling function bin_timePaleoPhy and the simulation function binTimeData for more information on formatting.

Unlike some functions in paleotree, such as the diversity curve functions, intervals must be both sequential and non-overlapping. The diversity curve functions deal with such issues by assuming taxa occur from the base of the interval they are first found in until the end of the last interval they are occur in. This inflation of boundary crossers could badly bias estimates of per-capita diversification rates.

Value
This function will invisibly return a ten column matrix, where the number of rows is equal to the number of intervals. The first two columns are interval start and end times and the third column is interval length. The fourth through eighth column is the four fundamental classes of taxa from Foote (2001): Nbt, NbL, NFt, NFL and their sum, N. The final two columns are the per-capita rates estimated for each interval in units per lineage time-units; the ninth column is the origination rate (’pRate’) and the tenth column is the extinction rate (’qRate’).
References


See Also

DiversityCurves, SamplingConv

Examples

# with the retiolinae dataset
data(retiolitinae)
perCapitaRates(retioRanges)

# Simulate some fossil ranges with simFossilRecord
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(0,100), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
# simulate a fossil record with imperfect sampling with sampleRanges()
rangesCont <- sampleRanges(taxa,r = 0.5)
# Now let's use binTimeData() to bin in intervals of 5 time units
rangesDisc <- binTimeData(rangesCont,int.length = 5)
# and get the per-capita rates
perCapitaRates(rangesDisc)
# on a log scale
perCapitaRates(rangesDisc,logRates = TRUE)

# get mean and median per-capita rates
res <- perCapitaRates(rangesDisc,plot = FALSE)
apply(res[,c("pRate","qRate")],2,mean,na.rm = TRUE)
apply(res[,c("pRate","qRate")],2,median,na.rm = TRUE)

# with modern taxa
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nExtant = c(10,50))
taxa <- fossilRecord2FossilTaxa(record)
# simulate a fossil record with imperfect sampling with sampleRanges()
rangesCont <- sampleRanges(taxa,r = 0.5,modern.samp.prob = 1)
# Now let's use binTimeData() to bin in intervals of 5 time units
rangesDisc <- binTimeData(rangesCont,int.length = 5)
# and now get per-capita rates
perCapitaRates(rangesDisc)
perfectParsCharTree

Simulate a Set of Parsimony-Informative Characters for a Phylogeny

Description

Creates a simulated set of parsimony-informative characters for a given rooted phylogeny, with characters shared out equally across nodes in the phylogeny, with any remaining characters assigned randomly to nodes.

Usage

perfectParsCharTree(tree, nchar)

Arguments

- **tree**: A phylogeny of class 'phylo'
- **nchar**: Number of parsimonious binary characters to simulate on the phylogeny.

Details

This function takes some a tree and places a number of binary characters on the tree, with character states arranged as if the derived condition was gained once, at a single node, and never lost. This ensures that the resulting simulated character matrices have no character conflict, supporting a single solution under maximum parsimony.

If `nchar` is greater than the number of nodes on the input phylogeny (ignoring the root), then characters are first placed to evenly cover all nodes, with as many full passes of tree as possible. Any characters in excess are placed at random nodes, without replacement. In other words, if a tree has 10 nodes (plus the root) and 25 characters are simulated, 20 of those characters will consist of two 10-character 'full passes' of the tree. The remaining five will be randomly dropped on the tree.

If few characters are simulated than the number of nodes, these are randomly placed on the given topology without replacement, just as described above.

This function assumes, like almost every function in paleotree, that the tree given is rooted, even if the most basal node is a polytomy.

Value

A matrix of `nchar` parsimonious binary characters for each taxon on `tree`, with states 0 and 1.

Author(s)

David W. Bapst
Examples

data(retiolitinae)

# fewer characters than nodes
perfectParsCharTree(retioTree, nchar = 10)

# same as number of nodes (minus root)
perfectParsCharTree(retioTree, nchar = 12)

# more characters than the number of nodes
perfectParsCharTree(retioTree, nchar = 20)

---

plotOccData

**Plotting Occurrence Data Across Taxa**

Description

plotOccData takes occurrence data which has been sorted into a by-taxon list, such as that output by taxonSortBPDBocc or may be output by simulations using sampleRanges and produces a plot showing the age uncertainty associated with individual occurrences, with occurrences of the same taxon grouped by color.

Usage

plotOccData(occlist, groupLabel = NULL, occColors = NULL, 
            lineWidth = NULL, xlims = NULL)

Arguments

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>occList</td>
<td>A list where every element is a table of occurrence data for a different taxon, such as that returned by taxonSortBPDBocc. The occurrence data can be either a two-column matrix composed of the lower and upper age bounds on each taxon occurrence, or has two named variables which match any of the field names given by the PBDB API under either the 'pdbb' vocab or 'com' (compact) vocab for early and late age bounds.</td>
</tr>
<tr>
<td>groupLabel</td>
<td>A character vector with a single string giving the name for the occurrence dataset used, such as the taxonomic name of the group examined. If not given (the default) a generic plot title is appended.</td>
</tr>
<tr>
<td>occColors</td>
<td>A vector of numbers or characters indicating colors on a color palette for use with basic plot. Must be the same length as occList. If empty, the default, the colors used are sampled randomly from the rainbow() function.</td>
</tr>
<tr>
<td>lineWidth</td>
<td>A numeric value giving the length to be used for the width of lines plotted in plotOccData. If not given (the default), this is calculated using an algorithm that selects an optimal line width for plotting.</td>
</tr>
<tr>
<td>xlims</td>
<td>A two element vector controlling the width of the horizontal time-scale the occurrence bars are plotted against. By default, this is not given and calculated internally.</td>
</tr>
</tbody>
</table>
Details

This function was originally conceived of in the following blog post: Link

Value

This function will invisibly return a list, with each per-taxon element containing the two-column matrix of age bounds for occurrences.

Author(s)

David W. Bapst

See Also

taxonSortPBDBocc, occData2timeList and the example graptolite dataset at graptPBDB

Examples

#load example graptolite PBDB occ dataset
data(graptPBDB)

#get formal genera
occSpecies <- taxonSortPBDBocc(graptOccPBDB, rank = "species")

#plot it!
plotOccData(occSpecies)

#this isn't too many occurrences, because there are so few
  #formal grapt species in the PBDB

#genera is messier...

#get formal genera
occGenus <- taxonSortPBDBocc(graptOccPBDB, rank = "genus")

#plot it!
plotOccData(occGenus)

#some of those genera have occurrences with very large
  #age uncertainties on them!
plotTraitgram plots a traitgram showing the evolution of a continuous trait. If node values are not given (i.e. the data is empirical data collected from tips, rather than simulated data), maximum-likelihood ancestral trait estimation is used to calculate node values. (Ackerly, 2009) given a tree and a set of continuous trait values.

Usage

```
plotTraitgram(trait, tree, main = "", conf.int = TRUE, lwd = 1.5)
```

Arguments

- `trait` A vector of continuous trait values. If the length of `trait` is equal to the number of tips and number of nodes, than it is presumed internal node values are given. If the length of `trait` is equal to the number of tips of `tree` then these will be treated as tip values and ancestral trait reconstruction will be used to reconstruct the missing ancestral values. If `trait` is not named, or if internal node values are given, then values will be presumed to be in the order of tip/node numbering in `tree$edge`, which for tips is the same as the ordering in `tree$tip.label`.

- `tree` A `phylo` object.

- `main` Main title of traitgram plot.

- `conf.int` If `TRUE` (the default), confidence intervals are plotted.

- `lwd` The line width used for branches in the figure.

Details

By default, this function will use `ace` from the library ape to reconstruct ancestral traits and confidence intervals using the PIC method, if internal node values (i.e. ancestral node values) are not given.

As with many functions in the paleotree library, absolute time is always decreasing, i.e. the present day is zero.

Value

Return no value, just plot the traitgram.

Note

One should probably never do ancestral trait estimation without looking at the confidence intervals, as these reconstructed estimates tend to be very uncertain.

Author(s)

David W. Bapst
References

See Also
ace

Also see the functions traitgram in the library picante and phenogram in the library phytools.

Examples

```r
set.seed(444)
tree <- rtree(10)
trait <- rTraitCont(tree)

# first, traitgram without conf intervals
plotTraitgram(trait, tree, conf.int = FALSE)

# now, with
plotTraitgram(trait, tree)
# not much confidence, eh?

# plotting simulated data
  # with values for ancestral nodes as input
trait <- rTraitCont(tree, ancestor = TRUE)
plotTraitgram(tree = tree, trait = trait)
```

---

**pqr2Ps**

*Joint Probability of A Clade Surviving Infinity or Being Sampled Once*

**Description**

Given the rates of branching, extinction and sampling, calculates the joint probability of a random clade (of unknown size, from 1 to infinite) either (a) never going extinct on an infinite time-scale or (b) being sampled at least once, if it does ever go extinct. As we often assume perfect or close to perfect sampling at the modern (and thus we can blanket assume that living groups are sampled), we refer to this value as the Probability of Being Sampled, or simply P(s). This quantity is useful for calculating the probability distributions of waiting times that depend on a clade being sampled (or not).

**Usage**

```r
pqr2Ps(p, q, r, useExact = TRUE)
```
Arguments

- **p**: Instantaneous rate of speciation (lambda). If the underlying model assumed is anagenetic (e.g., taxonomic change within a single lineage, 'phyletic evolution') with no branching of lineages, then p will be used as the rate of anagenetic differentiation.

- **q**: Instantaneous rate of extinction (mu)

- **r**: Instantaneous rate of sampling

- **useExact**: If TRUE, an exact solution developed by Emily King is used; if FALSE, an iterative, inexact solution is used, which is somewhat slower (in addition to being inexact...).

Details

Note that the use of the word 'clade' here can mean a monophyletic group of any size, including a single 'species' (i.e., a single phylogenetic branch) that goes extinct before producing any descendants. Many scientists I have met reserve the word 'clade' for only groups that contain at least one branching event, and thus contain two 'species'. I personally prefer to use the generic term 'lineage' to refer to monophyletic groups of one to infinity members, but others reserve this term for a set of morphospecies that reflect an unbroken anagenetic chain.

Obviously the equation used makes assumptions about prior knowledge of the time-scales associated with clades being extant or not: if we’re talking about clades that originated a short time before the recent, the clades that will go extinct on an infinite time-scale probably haven’t had enough time to actually go extinct. On reasonably long time-scales, however, this infinite assumption should be reasonable approximation, as clades that survive 'forever' in a homogenous birth-death scenario are those that get very large immediately (similarly, most clades that go extinct also go extinct very shortly after originating... yes, life is tough).

Both an exact and inexact (iterative) solution is offered; the exact solution was derived in an entirely different fashion but seems to faithfully reproduce the results of the inexact solution and is much faster. Thus, the exact solution is the default. As it would be very simple for any user to look this up in the code anyway, here’s the unpublished equation for the exact solution:

\[
Ps = 1 - \frac{((p + q + r) - \sqrt{((p + q + r)^2 - (4 * p * q)))}}{(2 * p)}
\]

Value

Returns a single numerical value, representing the joint probability of a clade generated under these rates either never going extinct or being sampled before it goes extinct.

Author(s)

This function is entirely the product of a joint effort between the package author (David W. Bapst), Emily A. King and Matthew W. Pennell. In particular, Emily King solved a nasty bit of calculus to get the inexact solution and later re-derived the function with a quadratic methodology to get the exact solution. Some elements of the underlying random walk model were provided by S. Nalayanan (a user on the website stackexchange.com) who assisted with a handy bit of math involving Catalan numbers.
References


See Also

SamplingConv

Examples

```r
#with exact solution
pqr2Ps(p = 0.1,q = 0.1,r = 0.1,useExact = TRUE)
#with inexact solution
pqr2Ps(p = 0.1,q = 0.1,r = 0.1,useExact = TRUE)
```

---

**probAnc**  
*Probability of being a sampled ancestor of another sampled taxon*

Description

Uses models from Foote (1996) to calculate the probability

Usage

```r
probAnc(p, q, R, mode = "budding", analysis = "directDesc", Mmax = 85, nrep = 10000)
```

Arguments

- **p**  
  Instantaneous rate of speciation (lambda). If the underlying model assumed is anagenetic (e.g. taxonomic change within a single lineage, 'phyletic evolution') with no branching of lineages, then p will be used as the rate of anagenetic differentiation.

- **q**  
  Instantaneous rate of extinction (mu)

- **R**  
  Per-interval probability of sampling a taxon at least once

- **mode**  
  Mode of morphotaxon differentiation, based on definitions in Foote, 1996. Can be pure cladogenetic budding ("budding"), pure cladogenetic bifurcating ("bifurcating") or pure anagenetic within-lineage change ("anagenesis"; i.e. Foote's 'phyletic change'). Default mode is "budding".

- **analysis**  
  The type of analysis to be performed, either the probability of sampling direct descendants ("directDesc") or of sampling indirect descendants ("indirectDesc").
probAnc

Mmax

The maximum number of direct descendants (M) to sum over in the function, which is ideally meant to be a sum from zero to infinity, like nrep. Unfortunately, \((2+M)\) is used in a factorial, which means we are limited to a relatively small upper bound on M.

nrep

Number of repetitions to run in functions which are meant to sum over infinity. Default is arbitrarily high.

Details

probAnc obtains the probability of sampling a descendant of a morphotaxon in the fossil record, given the sampling probability and estimates of origination and extinction rates. These values are always calculated assuming infinite time for the potential ancestor to produce daughter taxa (assuming it lives that long) and under homogenous birth, death and sampling rates/probabilities, which is a situation that may be overly ideal relative to many real fossil records.

This can be calculated for either direct descendants, i.e. the probability of sampling any morphotaxa that arise immediately from the particular morphotaxon that could be an ancestor, or indirect descendants, i.e. the probability for any morphotaxon that has the morphotaxon of question as an ancestor, no matter how distant. See the argument analysis for details. Mode of differentiation can also be varied for three different models, see the argument mode.

This probability is calculated including the probability that extinction might occur before any descendants are produced. Thus, if \(p = q\), the probability of a taxon going extinct before it produces any descendants will be 0.5, which means that even when sampling is perfect (\(R = 1\), meaning completeness of 100 can be no higher than 0.5. See Foote (1996) for a graphic depiction of this non-intuitive ceiling. For reasons (probably?) having to do with finite approximations of infinite summations, values close to perfect sampling may have values slightly higher than this ceiling, which is also apparent visually in the figures in Foote (1996). Thus, values higher than 0.5 when \(p = q\) should be discounted, and in general when sampling rate is high, results should be treated cautiously as overestimates.

References


See Also

SamplingConv

Examples

#examples, run at very low nrep for sake of speed (examples need to be fast)

#default: probability of sampling a direct descendant
probAnc(p = 0.1, q = 0.1, R = 0.5,  
    mode = "budding", analysis = "directDesc", nrep = 100)

#other modes
probAnc(p = 0.1, q = 0.1, R = 0.5,  
    mode = "bifurcating", analysis = "directDesc", nrep = 100)  
probAnc(p = 0.1, q = 0.1, R = 0.5,  


mode = "anagenesis", analysis = "directDesc",nrep = 100)

#probability of having sampled indirect descendants of a taxon
probAnc(p = 0.1, q = 0.1, R = 0.5,
mode = "budding", analysis = "indirectDesc",nrep = 100) #default
probAnc(p = 0.1, q = 0.1, R = 0.5,
mode = "bifurcating", analysis = "indirectDesc",nrep = 100)
probAnc(p = 0.1, q = 0.1, R = 0.5,
mode = "anagenesis", analysis = "indirectDesc",nrep = 100)

---

RaiaCopesRule

*Dated Trees and Trait Data for Ammonites, Ceratopsians and Cervids from Raia et al. 2015*

**Description**

Dated phylogenetic trees for fossil ammonite genera, fossil ceratopsian species and (both extinct and extant) cervid species, as well as trait data (shell diameter and fractal complexity of the first suture) for the ammonite dataset, taken from the recent publication by Raia et al. (2015) in *The American Naturalist*. The goal of this paper was to examine the relationship between ornamental complexity and body size in three very different groups, but the datasets are very relatively large and useful for demonstrating application of comparative methods to fossil trees.

**Format**

The attached datasets consist of three phylogenetic trees as *phylo* objects, a *data.frame* consisting of three traits for ammonites (the third trait is the log of stratigraphic duration), and the two physical traits (shell size and suture complexity) as separate vectors, with taxon names.

**Details**

It appears that the trees were dated with tips at the last appearances, although this doesn’t appear to be explicitly stated in Raia et al.

**Source**


**See Also**

retiolidinae, macroperforateForam
Examples

data(RaiaCopesRule)

# plotting trees
plot(ladderize(ammoniteTreeRaia)); axisPhylo()
plot(ceratopsianTreeRaia); axisPhylo()
plot(cervidTreeRaia); axisPhylo()

# plotting traitgrams for ammonites
plotTraitgram(tree = multi2di(ammoniteTreeRaia), trait = sutureComplexity, 
               conf.int = FALSE, main = "Ammonite Suture Complexity")
plotTraitgram(tree = multi2di(ammoniteTreeRaia), trait = shellSize, 
               conf.int = FALSE, main = "Ammonite Shell Diameter")

# The data set was generated by sourcing the following script:

library(paleotree)

# Let's read in the trees from Raia et al 2015, AmNat
# following is taken from their supplemental appendix, available at AmNat
# they all appear to be trees dated to the last appearance times
# *and* specifically the end-boundary of the interval containing the last appearance

# ammonite genera

ammoniteTreeRaia <- paste0("(((Araxoceras:4,Eraraxoceras:4)Araxoceratidae:26.5,Pseudasp:
" "Dinaritaceae:24.299988,Grambergia:42.5,(Amphiphanoceras:6, Megaphyllites:46.399994) Megaph:
" "yllitaceae:36.5,(Proteusites:11,Nathorstites:21)Nathorstitaceae:31.5,(Inyoites:7,Lanceolit:
Raia et al. are unclear, however... a handful of taxa are known to last occur at the end-Cretaceous mass extinctions. Latest occurring tips are:

```
phylloceras
```

so we can treat distance of Phylloceras from root + end Cretaceous (66.043 Ma) as $\text{root.time} = \text{node.depth.edge.length}(\text{ammoniteTreeRaia})[\text{which}(\text{ammoniteTreeRaia$tip.label == "Phylloceras")}]$

now let's plot it

```
plot(ladderize(\text{ammoniteTreeRaia}));axis(\text{Phylo})
```

and let's load trait data from Raia et al. Appendix B:

```
FD = fractal dimension of first suture (suture complexity)
log D = log of the mean shell diameter per genus (body size)
log dur = log of the stratigraphic duration in million years.
```

```
ammoniteTraitsRaia <- \text{read.table}("ammoniteTraitsRaia.txt", row.names = 1, header = TRUE)
```

```
sutureComplexity <- ammoniteTraitsRaia$FD
shellSize <- ammoniteTraitsRaia$Log_D
```

```
plotTraitgram\text{tree} = \text{multi2di}(\text{ammoniteTreeRaia}), trait = sutureComplexity, conf.int = FALSE, main = "Ammonite Suture Complexity")
```

```
plotTraitgram\text{tree} = \text{multi2di}(\text{ammoniteTreeRaia}), trait = shellSize, conf.int = FALSE, main = "Ammonite Shell Diameter")
```

```
# End(Not run)
```
Raia et al. placed origin of ceratopsians at ~163 Ma, base of Oxfordian.

```r
ceratopsianTreeRaia <- read.tree(text = ceratopsianTreeRaia)
```

```r
cervidTreeRaia <- read.tree(text = cervidTreeRaia)
```

```r
# If you plot the tree there seem to be a lot more taxa that are *almost* as late-occurring
# unclear if this is recently extinct taxa, computational rounding error, or what
# so we can treat distance of Dama dama to root as $root.time
(cervidTreeRaia$root.time <-
node.depth.edge.length(cervidTreeRaia)[which(cervidTreeRaia$tip.label == "Dama_dama")])

plot(cervidTreeRaia);axisPhylo()
save.image("RaiaCopesRule.rdata")

---

### resolveTreeChar

**Resolve Polytomies Using Parsimony-Based Reconstruction of a Discrete Character**

**Description**

This function resolves a set of given topology with less than fully-binary phylogenetic resolution so that lineages are shifted and internal nodes added that minimize the number of independent character transitions needed to explain an observed distribution of discrete character states for the taxa on such a tree, under various maximum-parsimony algorithms of ancestral character reconstruction, powered ultimately by function ancestral.pars in library phangorn. This function is mainly designed for use with poorly resolved trees which are being assessed with the function minCharChange.

**Usage**

```r
resolveTreeChar(tree, trait, orderedChar = FALSE, stateBias = NULL,
iterative = TRUE, cost = NULL, ambiguity = c(NA, "?")
dropAmbiguity = FALSE, polySymbol = "&", contrast = NULL)
```

**Arguments**

- **tree**
  A cladogram of type 'phylo'. Any branch lengths are ignored.

- **trait**
  A vector of trait values for a discrete character, preferably named with taxon names identical to the tip labels on the input tree.

- **orderedChar**
  Is the character of interest given for trait ordered or not? If FALSE (the default), then for each polytomy, all child nodes that appear to have the same state as the ancestor node will remain in the polytomy, and any additional states held by child nodes will each be grouped into their own unique polytomy that forms from a descendant node of the original polytomy. If TRUE, then the character will be reconstructed with a cost (step) matrix of a linear, ordered character, and polytomies will be resolved so that lineages with different states will be placed into a nested ladder that reflects the ordered character. As with the unordered option, child nodes with a state equivalent to the ancestral node will remain in the polytomy, while more primitive or more derived states will be sorted into their own separate ladders composed of paraphyletic groups, ordered so to move 'away' state-by-state from the ancestral node’s inferred character state.
**resolveTreeChar**

**stateBias**
This argument controls how `resolveTreeChar` handles ancestral node reconstructions that have multiple states competing for the maximum weight of any state (i.e., if states 0 and 1 both have 0.4 of the weight). The default, where `stateBias = NULL` causes uncertainty at nodes among states to be treated as a single ‘group’ identical to any states within it. Essentially, this means that for the example polytomy where the ancestor has maximum weight for both 0 and 1, any child nodes with 0, 1 or both of these states will be considered to have an identical state for the purpose of grouping nodes for the purpose of further resolving polytomies. If and only if `orderedChar = TRUE`, then additional options of `stateBias = 'primitive'` and `stateBias = 'derived'` become available, which instead force uncertain node assignments to either be the most primitive (i.e., the minimum) or the most derived (i.e., the maximum) among the maximum-weight states. In particular, `stateBias = 'primitive'` should favor gains and bias any analysis of character transitions against finding reversals.

**iterative**
A logical argument which, if TRUE (the default), causes the function to repeat the polytomy-resolving functionality across the entire tree until the number of nodes stabilizes. If FALSE, polytomies are only passed a single time.

**cost**
A matrix of the cost (i.e., number of steps) necessary to change between states of the input character trait. If NULL (the default), the character is assumed to be unordered with equal cost to change from any state to another. Cost matrices only impact the “MPR” algorithm; if a cost matrix is given but `type = "ACCTRAN"`, an error is issued.

**ambiguity**
A vector of values which indicate ambiguous (i.e., missing or unknown) character state codings in supplied trait data. Taxa coded ambiguously as treated as being equally likely to be any state coding. By default, NA values and "?" symbols are treated as ambiguous character codings, in agreement with behavior of functions in packages phangorn and Claddis. This argument is designed to mirror an hidden argument with an identical name in function phyDat in package phangorn.

**dropAmbiguity**
A logical. If TRUE (which is not the default), all taxa with ambiguous codings as defined by argument ambiguity will be dropped prior to ancestral nodes being inferred. This may result in too few taxa.

**polySymbol**
A single symbol which separates alternative states for polymorphic codings; the default symbol is "&", following the output by Claddis’s ReadMorphNexus function, where polymorphic taxa are indicated by default with a string with state labels separated by an "&" symbol. For example, a taxon coded as polymorphic for states 1 or 2, would be indicated by the string "1&2". `polySymbol` is used to break up these strings and automatically construct a fitting contrast table for use with this data, including for ambiguous character state codings.

**contrast**
A matrix of type integer with cells of 0 and 1, where each row is labeled with a string value used for indicating character states in trait, and each column is labeled with the formal state label to be used for assign taxa to particular character states. A value of 1 indicates that the respective coding string for that row should be interpreted as reflecting the character state listed for that column. A coding could reflect multiple states (such as might occur when taxa are polymorphic for some morphological character), so the sums of rows and columns can sum
to more than 1. If contrast is not NULL (the default), the arguments will nullify. This argument is designed to mirror an hidden argument with an identical name in function phyDat in package phangorn. This structure is based on the contrasts tables used for statistical evaluation of factors. See the phangorn vignette "Special features of phangorn" for more details on its implementation within phangorn including an example. See examples below for the construction of an example contrast matrix for character data with polymorphisms, coded as character data output by Cladis's ReadMorphNexus function, where polymorphic taxa are indicated with a string with state labels separated by an "&" symbol.

Details

As shown in the example code below, this function offers a wide variety of options for manipulating the maximum-parsimony algorithm used (i.e. MPR versus ACCTRAN), the ordering (or not) of character states, and potential biasing of uncertainty character state reconstructions (when ordered characters are assessed). This allows for a wide variety of possible resolutions for a given tree with polytomies and a discrete character. In general, the author expects that use of this function will be optimal when applied to ordered characters using one of the stateBias options, perhaps stateBias = "primitive" (based on theoretical expectations for slow evolving characters). However, anecdotal use of this function with various simulation datasets suggests that the results are quite variable, and so the best option needs to be assessed based on the prior assumptions regarding the data and the performance of the dataset with the various arguments of this function.

Value

Returns the resulting tree, which may be fully resolved, partly more resolved or not more resolved at all (i.e. have less polytomies) depending on what was possible, as constrained by ambiguities in character reconstructions. Applying multi2di is suggested as a post-step to obtain a fully-resolved cladogram, if one is desired.

Author(s)

David W. Bapst

References


See Also

ancPropStateMat which is used internally by this function. This function was intentionally designed for use with minCharChange.

Examples

```r
# let's write a quick & dirty ancestral trait plotting function
quickAncPlot <- function(tree, trait, cex, orderedChar = FALSE, type = "MPR", cost = NULL)(
  ancData <- ancPropStateMat(tree = tree, trait = trait, orderedChar = orderedChar)
  ancCol <- c(1:ncol(ancData)) + 1
  plot(tree, show.tip.label = FALSE, no.margin = TRUE, direction = "upwards")
  tipLabels(pch = 16, pie = ancData[(1:ntip(tree)),], cex = cex, piecol = ancCol,
  col = 0)
  nodelabels(pie = ancData[-(1:ntip(tree)),], cex = cex, piecol = ancCol)
)

# examples with simulated data
set.seed(2)

tree <- rtree(50)
# simulate under a likelihood model
trait <- rTraitDisc(tree, k = 3, rate = 0.7)

tree <- degradeTree(tree, prop_collapse = 0.6)

tree <- ladderize(tree, right = FALSE)

# a bunch of type = MPR (default) examples

treeUnord <- resolveTreeChar(tree, trait, orderedChar = FALSE)
treeOrd <- resolveTreeChar(tree, trait, orderedChar = TRUE, stateBias = NULL)
treeOrdPrim <- resolveTreeChar(tree, trait, orderedChar = TRUE, stateBias = "primitive")
treeOrdDer <- resolveTreeChar(tree, trait, orderedChar = TRUE, stateBias = "derived")

# compare number of nodes
Nnode(tree) # original
Nnode(treeUnord) # unordered, biasStates = NULL, MPR
Nnode(treeOrd) # ordered, biasStates = NULL
Nnode(treeOrdPrim) # ordered, biasStates = 'primitive'
Nnode(treeOrdDer) # ordered, biasStates = 'derived'

# let's compare original tree with unordered-resolved tree layout(1:2)
quickAncPlot(tree, trait, orderedChar = FALSE, cex = 0.3)
text(x = 43, y = 10, "Original", cex = 1.5)
quickAncPlot(treeUnord, trait, orderedChar = FALSE, cex = 0.3)
text(x = 43, y = 10, "orderedChar = FALSE", cex = 1.5)
# some resolution gained
```
now let's compare the original and ordered, both `biasStates` = NULL
layout(1:2)
quickAncPlot(tree,trait,orderedChar = FALSE,cex = 0.3)
text(x = 43,y = 10,"Original",cex = 1.5)
quickAncPlot(treeOrd,trait,orderedChar = TRUE,cex = 0.3)
text(x = 43,y = 10,"orderedChar = TRUE",cex = 1.5)

#now let's compare the three ordered trees
layout(1:3)
quickAncPlot(treeOrd,trait,orderedChar = TRUE,cex = 0.3)
text(x = 41,y = 8,"ordered, biasStates = NULL",cex = 1.5)
quickAncPlot(treeOrdPrim,trait,orderedChar = TRUE,cex = 0.3)
text(x = 41.5,y = 8,"ordered, biasStates = 'primitive'",cex = 1.5)
quickAncPlot(treeOrdDer,trait,orderedChar = TRUE,cex = 0.3)
text(x = 42,y = 8,"ordered, biasStates = 'derived'",cex = 1.5)

#let's compare unordered with ordered, `biasStates` = 'primitive'
layout(1:2)
quickAncPlot(treeUnord,trait,orderedChar = FALSE,cex = 0.3)
text(x = 41,y = 8,"orderedChar = FALSE",cex = 1.5)
quickAncPlot(treeOrdPrim,trait,orderedChar = TRUE,cex = 0.3)
text(x = 40,y = 11,"orderedChar = TRUE",cex = 1.5)
text(x = 40,y = 4,"biasStates = 'primitive'",cex = 1.5)

#these comparisons will differ greatly between datasets
# need to try them on your own
layout(1)

### retiolitinae  
*Cladogram and Range Data for the Retiolitinae*

**Description**

The majority rule consensus cladogram for 22 genera from the Retiolitinae, a clade of Silurian retiolitids, along with discrete time interval data taken from the same publication (Bates et al., 2005). Additional character state data are included for three major, binary-state morphological traits.

**Format**

This dataset is composed of three objects:

- `retioTree` An ape `phylo` object containing the consensus cladogram.
retiolitinae

retioRanges A list containing two matrices. The first matrix describes the first and last interval times for 20 Silurian graptolite zones and the second matrix describes when the various genera on the cladogram first and last appear in those graptolite zones. (In other words, retioRanges has the ‘timeList’ format called by some paleotree functions).

retioChar A matrix containing binary presence-absence character states for these 22 Retiolitinae genera for three characters which they vary in: the presence of extrathecal threads (note only one taxon lacks this character state), the presence of determinant growth and the secondary loss of a nema via resorption. Note these character do not vary within these genera.

Details

Interval dates were taken from Sadler et al. (2009). These zones were not a 1-1 match to those in Bates et al., so it took some merging and splitting by the package author, so buyer beware.

Character data are from an in prep manuscript containing character data for certain major morphological innovations of graptoloids, coded for a large number of genera based on an extensive survey of the published descriptions. The character data presented here is a small subset of the full dataset.

Source

Source for cladogram and zonal ranges for genera:

Source for interval dates for graptolite zones:

Source for morphological character data:
Collected for Bapst and Mitchell, in prep.

See Also

For more example graptolite datasets, see graptDisparity

Examples

```r
#load data
data(retiolitinae)

taxicDivDisc(retioRanges)

plot(retioTree)
tiplabels(pch = 16,col = (retioChar[,2]+1),adj = 0.25)

#Use basic time-scaling (terminal branches only go to FADs)
tree <- bin_timePaleoPhy(tree = retioTree,timeList = retioRanges,type = "basic",)
```
reverseList

Description

Takes a list and reverses the list structure, such that list composed of five elements with eight sub-elements is restructured to have eight elements with five sub-elements each, with the order of elements and sub-elements being retained despite their reversal in hierarchical position.

Usage

reverseList(list, simplify = FALSE)

Arguments

- list: A list composed of multiple elements, with each element a vector or list of equal length
- simplify: Should the result be simplified, as the argument in sapply

Details

The function will fail and return an error if all sub-elements are not vectors or lists of equal length. This function can be useful for instances when each element of a list is by-sample, composed of multiple, different tests on that sample, but where for further analysis/plotting, it would be beneficial to have a list where each element represented values from the same test performed across multiple samples (i.e. plotting a box-plot).

Value

Returns a list with a reversed structure relative to the input, see above.

Author(s)

David W. Bapst
Examples

```r
list1 <- list(list(1:3), list(1:3), list(1:3))
reverseList(list1, simplify = FALSE)
reverseList(list1, simplify = TRUE)
```

---

**rootSplit**  
**Split Tip Taxa by Root Divergence**

**Description**

Sorts terminal taxa into groups descended from each lineage splitting off of the root node.

**Usage**

```r
rootSplit(tree)
```

**Arguments**

- `tree`  
  A phylo object

**Details**

This function can be useful for studying the timing in the order of appearance of descended from different lineages descended from the first bifurcation.

**Value**

Returns a list with each element a character vector containing the names of terminal taxa descended from each lineage splitting off of the root node.

**Author(s)**

David W. Bapst

**Examples**

```r
tree <- rtree(100)
rootSplit(tree)
```
Description

A function for simulating the effect of incomplete sampling of the fossil record.

Usage

```r
c7ampleRanges(taxad, r, alpha = 1, beta = 1, rTimeRatio = 1,
    modern.samp.prob = 1, min.taxa = 2, ranges.only = TRUE, minInt = 0.01,
    merge.cryptic = TRUE, randLiveHat = TRUE, alt.method = FALSE,
    plot = FALSE)
```

Arguments

taxad A two-column matrix of per-taxon ranges. The five-column matrix output of `simFossilRecord`, post transformation with `fossilRecord2fossilTaxa` can also be supplied, which will be common in simulation usages.

r Instantaneous average sampling rate per lineage time units; given as a vector of length one or length equal to the number of taxa.

alpha Alpha parameter of beta distribution; given as a vector of length one or length equal to the number of taxa.

beta Beta parameter of beta distribution; given as a vector of length one or length equal to the number of taxa.

rTimeRatio Ratio of most recent sampling rate over earliest sampling rate; given as a vector of length one or length equal to the number of taxa.

modern.samp.prob Probability of sampling living taxa at the present day (time = 0), see below.

min.taxa Minimum number of taxa sampled. The default is 2.

ranges.only If TRUE, gives taxon first and last occurrences only. If FALSE, gives the time of all sampling events as a list.

minInt Minimum interval size used for simulating complex models

merge.cryptic If TRUE, sampling events for cryptic species will be merged into one taxon.

randLiveHat If TRUE, taxa still alive at modern day have the end-point of their 'hat' chosen from a uniform distribution.

alt.method If TRUE, use the alternative method of discretizing time even if a simple model of sampling is being simulated.

plot If TRUE, plots the sampling models for each taxon against time.
Details

This function implements a range of sampling models in continuous time. By default, sampling is simulated under the simplest model, where sampling occurs as a Poisson process under an instantaneous sampling rate ($r$) which is homogeneous through time and across lineages (Foote, 1997). Under this model, the waiting times to sampling events are exponentially distributed, with an average waiting time of $1/r$. This useful property allows sampling to be rapidly simulated for many taxa under this simple model in sampleRanges, by repeatedly drawing waiting times between sampling events from an exponential distribution. This is the model that is run when alpha, beta and rTimeRatio.

In addition to this simple model, sampleRanges also can consider a range of additional models, including the hatP and incP options of Liow et al. (2010). To describe the behavior of these models, users alter the default values for alpha, beta and rTimeRatio. These parameters, and $r$, can either be a single value which describes the behavior of the entire dataset or as a vector, of same length as the number of taxa, which describes the per-taxon value. When any rTimeRatio, alpha or beta value is not equal to one, then the sampling rate will vary across the duration of a taxon’s temporal range. In general, setting alpha and beta equal to a value above 2 produce “hat” or bell-shaped curves where sampling rates peak at the midpoint of taxon ranges, while setting them unequal will produce asymmetric bell curves according to the beta function (Liow et al., 2010; Liow et al. set alpha = 4 and beta = 4). rTimeRatio is the ratio of the sampling rate of the latest/most recent time divided by the earliest/oldest time.

The input r values will be interpreted differently based on whether one r value or per-taxon values were used. If one value was input, then it is assumed that r represent the grand mean of r for the entire dataset for purposes of time-varying r, such that if rTimeRatio is not equal to 1, taxa near the end and start of the dataset will have very different per-taxon mean sampling rate. If per-taxon values of r were input, then each r is consider the per-taxon mean sampling rate. These will not be changed, but any within-lineage variation is distributed so that the mean is still the input per-taxon value. This also changes the interpretation of rTimeRatio, such that when a single r value and rTimeRatio is given, it is assumed the ratio describes the change in sampling rates from the start of the dataset to the end, while if multiple values are given for either r or rTimeRatio will instead see the value as describing the ratio at the first and last times of each taxon. For the pure hat model, this interpretation of ‘r’ as a grand mean sampling means that taxa will have a sampling rate of $2^r$ at the mid-peak of their range, which will have considerable implications for taxonomic incompleteness.

The particular distinctions about these parameter values are important: all models simulated in sampleRanges are structured to be effectively nested inside a most general model with parameters r, alpha, beta and rTimeRatio.

Note that the modeling of sampling in this function is independent and secondary of the actual simulation of the ranges, which are (generally) produced by the models of simFossilRecord with argument r (sampling rate) not set. Thus, ‘hat-shaped range distributions’ are only contained within single morphotaxa; they do not cross multiple morphotaxa in the case of anagenesis. Cryptic taxa each have their own hat and do not share a single hat; by default the ranges of cryptic taxa are merged to produce the range of a single observed morphotaxon.

‘Hats’ are constrained to start and end with a taxon’s range, representing the rise and fall of taxa in terms of abundance and geographic range (Liow et al., 2010). However, for still-living taxa at the modern day, it is unknown how much longer they may be alive (for memoryless Poisson models, there is no age-dependent extinction). The treatment of these taxa with regards to their ‘hat’ (i.e. the beta distribution) is controlled by randLivehat: when randLiveHat = FALSE, the beta distribution
is fit so that the last appearance of still-alive taxa at the modern day is treated as a last appearance for calculating the hat. When TRUE, the default option, the still-alive taxa are considered to have gotten some distance between 0 and 1 through the beta distribution, as of the modern day. This point of progression is stochastically selected for each taxon by pulling a number from a uniform distribution, and used for calculating the hat.

Because sampling rate varies over morphotaxon ranges under any of these more complex models, sampling events cannot be quickly simulated as waiting times pulled from an exponential distribution. Instead, the taxon durations are discretized into a large number of small time intervals of length minInt (see above; minInt should be small enough that only one sampling event could feasibly happen per interval). The probability of an event occurring within each interval is calculated and used to stochastically simulate sampling events. For each interval, a number between 0 and 1 is randomly pulled from a uniform distribution and to the per-interval sampling probability to test if a sampling event occurred (if the random number is less than the probability, a sampling event is recorded). In general, this method is slower but otherwise comparable to the quicker waiting times method. See the examples below for a small test of this.

As with many functions in the paleotree library, absolute time is always decreasing, i.e. the present day is zero.

If min.taxa is set to zero, the simulation may produce output in which no taxa were ever sampled.

If `modernsamp.prob` is set to 1.0 (the default), then living taxa will always be sampled at least at the present day (if there are any living taxa). If the probability is less than 1, they will be sampled with that probability at the modern day.

By default, this function will merge sampling events from morphologically cryptic taxa, listing them as occurrences for the earliest member of that group. To change this behavior, set `merge.cryptic = FALSE`.

Conditioning on sampling some minimum number of taxa may create strange simulation results for some analyses, such as simulation analyses of birth-death processes. Set `min.taxa = 0` to remove this conditioning.

Value

If `ranges.only` is TRUE, then the output is a two-column per-taxon matrix of first and last appearances in absolute time. NAs mean the taxon was never sampled in the simulation.

If `ranges.only = FALSE` (the default), the output is a list, where each element is a vector of sampling events the timing of sampling events, each corresponding to a different taxon in the input. Elements that are NA are unsampled taxa.

Author(s)

David W. Bapst

References


sampleRanges

See Also

simFossilRecord, binTimeData

Examples

```r
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(30,40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
layout(1:2)
#let's see what the 'true' diversity curve looks like in this case
taxicDivCont(taxa)
#simulate a fossil record with imperfect sampling with sampleRanges
rangesCont <- sampleRanges(taxa, r = 0.5)
#plot the diversity curve based on the sampled ranges
taxicDivCont(rangesCont)
#compare the true history to what we might observe!

#let's try more complicated models!

#a pull-to-the-recent model with x5 increase over time similar to Liow et al.'s incP
layout(1:2)
rangesCont1 <- sampleRanges(taxa, r = 0.5,rTimeRatio = 5,plot = TRUE)
taxicDivCont(rangesCont1)

#a hat-shaped model
layout(1:2)
rangesCont1 <- sampleRanges(taxa, r = 0.5, alpha = 4, beta = 4,plot = TRUE)
taxicDivCont(rangesCont1)

#a combination of these
layout(1:2)
rangesCont1 <- sampleRanges(taxa, r = 0.5, alpha = 4, beta = 4,rTimeRatio = 5,plot = TRUE)
taxicDivCont(rangesCont1)

#testing with cryptic speciation
layout(1)
recordCrypt <- simFossilRecord(p = 0.1, q = 0.1, prop.cryptic = 0.5, nruns = 1,
nTotalTaxa = c(20,30), nExtant = 0)
taxaCrypt <- fossilRecord2FossilTaxa(recordCrypt)
rangesCrypt <- sampleRanges(taxaCrypt, r = 0.5)
taxicDivCont(rangesCrypt)

#an example of hat-shaped models (beta distributions) when there are live taxa
set.seed(444)
recordLive <- simFossilRecord(p = 0.1, q = 0.05, nruns = 1,
nTotalTaxa = c(5,100), nExtant = c(10,100))
taxaLive <- fossilRecord2FossilTaxa(recordLive)
#with end-points of live taxa at random points in the hat
rangesLive <- sampleRanges(taxaLive, r = 0.1, alpha = 4, beta = 4,randLiveHat = TRUE,plot = TRUE)
#with all taxa end-points at end-point of hat
```
rangesLive <- sampleRanges(taxaLive, r = 0.1, alpha = 4, beta = 4, randLiveHat = FALSE, plot = TRUE)

# simulate a model where sampling rate evolves under brownian motion
tree <- tax2phylo(taxa, obs = taxa[,3])
sampRateBM <- rTraitCont(tree)
sampRateBM <- sampRateBM-min(sampRateBM)
layout(1:2)
rangesCont1 <- sampleRanges(taxa, r = sampRateBM, plot = TRUE)
taxicDivCont(rangesCont1)

# evolving sampling rate, hat model and pull of the recent
layout(1:2)
rangesCont1 <- sampleRanges(taxa, r = sampRateBM, alpha = 4, beta = 4, rTimeRatio = 5, plot = TRUE)
taxicDivCont(rangesCont1)
layout(1)

# the simpler model is simulated by pulling waiting times from an exponential
# more complicated models are simulated by discretizing time into small intervals
# are these two methods comparable?
# let's look at the number of taxa sampled under both methods
summary(replicate(100, sum(!is.na(sampleRanges(taxa, r = 0.5, alt.method = FALSE)[,1]))))
summary(replicate(100, sum(!is.na(sampleRanges(taxa, r = 0.5, alt.method = TRUE)[,1]))))
# they look pretty similar!

SamplingConv

Converting Sampling Estimates

Description

Various functions for converting between estimates of sampling in the fossil record.

Usage

sProb2sRate(R, int.length = 1)
sRate2sProb(r, int.length = 1)
pqsRate2sProb(r, p, q, int.length = 1)
qsProb2Comp(R, q, p = NULL, mode = "budding", nrep = 10000)
qsRate2Comp(r, q)
SamplingConv

Arguments

<table>
<thead>
<tr>
<th>Argument</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>R</td>
<td>Per-interval probability of sampling a taxon at least once</td>
</tr>
<tr>
<td>int.length</td>
<td>Length of Time Intervals</td>
</tr>
<tr>
<td>r</td>
<td>Instantaneous rate of sampling</td>
</tr>
<tr>
<td>p</td>
<td>Instantaneous rate of speciation (lambda). If the underlying model assumed is anagenetic (e.g. taxonomic change within a single lineage, 'phyletic evolution') with no branching of lineages, then p will be used as the rate of anagenetic differentiation.</td>
</tr>
<tr>
<td>q</td>
<td>Instantaneous rate of extinction (mu)</td>
</tr>
<tr>
<td>mode</td>
<td>Mode of morphotaxon differentiation, based on definitions in Foote, 1996. Can be pure cladogenetic budding (“budding”), pure cladogenetic bifurcating (“bifurcating”) or pure anagenetic within-lineage change (“anagenesis”; i.e. Foote’s ‘phyletic change’). Default mode is “budding”.</td>
</tr>
<tr>
<td>nrep</td>
<td>Number of repetitions to run in functions which are meant to sum over infinity. Default is arbitrarily high.</td>
</tr>
</tbody>
</table>

Details

This is a family of functions which all convert from some estimate of sampling to another estimate of sampling. Some of these also require estimates of an rate associated with taxonomic diversification, such as the speciation/origin rate or extinction rate. Diversification rates used in these functions should always be the instantaneous rates, often called the per-capita rates by paleontologists (Foote, 2000).

As with many models used in the paleotree library, it is generally assumed that the fossil record of interest is composed of discrete relatively-static taxonomic units which diversify mainly by budding cladogenesis, and that sampling events are rare and approximated by a Poisson model of exponentially-distributed waiting times between sampling events. The veracity of those assumptions is difficult to test and the sensitivity of these analyses to relaxing those assumptions probably varies.

sProb2sRate and sRate2sProb give rough conversions for the probability of sampling once per time interval (R or “sProb” in this package as used in the references below) and the instantaneous rate of sampling per lineage/time unit (“sRate” or r). If you have estimates of the speciation and extinction rate, use pqsRate2sProb instead for a more accurate estimate of R.

qsProb2Comp and qsRate2Comp are different calculations for "Pp" or the probability/proportion of taxa sampled in a clade. Theoretically, one could use it to extrapolate out the 'true' diversity, assuming the sampling rate model was correct. (See Foote and Raup, 1996.)

See the references below for a more detailed explanation of the methods and formulae used. The relevant equations are generally found in the appendices of those papers.

Value

The converted sampling estimate, depending on the function used. See details above.

Author(s)

David W. Bapst, with advice from Michael Foote.
References


See Also

`sampleranges`, `make_durationFreqDisc`, `make_durationFreqCont`, `probAnc`, `pqr2Ps`.

Examples

```r
sRate2sProb(r = 0.5)
sProb2sRate(R = 0.1)
pqsRate2sProb(r = 0.5, p = 0.1, q = 0.1)

# different modes can be tried
qsProb2Comp(R = 0.1, q = 0.1, mode = "budding")
qsProb2Comp(R = 0.1, q = 0.1, mode = "bifurcating")
qsRate2Comp(r = 0.1, q = 0.1)
```

```r
seqTimeList

Construct a Stochastic Sequenced Time-List from an Unsequenced Time-List

Description

This function randomly samples from a timeList object (i.e. a list composed of a matrix of interval start and end dates and a matrix of taxon first and last intervals), to find a set of taxa and intervals that do not overlap, output as a new timeList object.

Usage

`seqTimeList(timeList, nruns = 100, weightSampling = FALSE)`
Arguments

- **timelist**
  A list composed of two matrices, giving interval start and end dates and taxon first and last occurrences within those intervals. Some intervals are expected to overlap (thus necessitating the use of this function), and datasets lacking overlapping intervals will return an error message.

- **nruns**
  Number of new timeList composed of non-overlapping intervals produced.

- **weightsampling**
  If TRUE, weight sampling of new intervals toward smaller intervals. FALSE by default.

Details

Many analyses of diversification and sampling in the fossil record require a dataset composed of sequential non-overlapping intervals, but the nature of the geologic record often makes this difficult, with taxa from different regions, environments and sedimentary basins having first and last appearances placed in entirely in-congruent systems of chronostratigraphic intervals. While one option is to convert such occurrences to a single, global stratigraphic system, this may still result in overlapping intervals when fossil collections are poorly constrained stratigraphically. (For example, this may often be the case in global datasets.) This function offers an approach to avoid this issue in large datasets by randomly subsampling the available taxa and intervals to produce stochastic sets of ranges composed of data drawn from non-overlapping intervals.

This function is stochastic and thus should be set for many runs to produce many such solutions. Additionally, all solutions found are returned, and users may wish to sort amongst these to maximize the number of intervals and number of taxa returned. A single solution which maximizes returned taxa and intervals may not be a precise enough approach to estimating sampling rates, however, given the uncertainty in data. Thus, many runs should always be considered.

By default, solutions are searched for without consideration to the length of intervals used (i.e. the selection of intervals is 'unweighted'). Alternatively, we can 'weight' selection toward the smallest intervals in the set, using the argument weightsampling. Smaller intervals presumably overlap less and thus should retain more taxa and intervals of more equal length. However, in practice with empirical datasets, the package author finds these approaches do not seem to produce very different estimates.

For some datasets, many solutions found using seqTimeList may return infinite sampling values. This is often due to saving too many taxa found in single intervals to the exclusion of longer-ranging taxa (see the example). This excess of single interval taxa is a clear artifact of the randomized seqTimeList procedure and such solutions should probably be ignored.

Value

A list, composed of three elements: nIntervals which is a vector of the number of intervals in each solution, nTaxa which is a vector of the number of taxa in each solution and timelists which is a list composed of each new timeList object as an element.

Author(s)

David W. Bapst
See Also

Resulting time-lists can be analyzed with `freqRat`, `durationFreq`, etc.

Additionally, `binTimeData` can be useful for simulating interval data.

Examples

```r
# Simulate some fossil ranges with simFossilRecord
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(60, 80), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
# simulate a fossil record with imperfect sampling with sampleRanges()
rangesCont <- sampleRanges(taxa, r = 0.1)

# Now let's use binTimeData to get ranges in discrete overlapping intervals
# via pre-set intervals input
presetIntervals <- cbind(c(1000, 995, 990, 980, 970, 960, 950, 940,
                           930, 900, 890, 880, 875, 870),
                          c(995, 989, 980, 975, 960, 950, 940, 930,
                           920, 900, 890, 880, 875, 870))
rangesDisc1 <- binTimeData(rangesCont, int.times = presetIntervals)

seqLists <- seqTimeList(rangesDisc1, nruns = 10)
seqLists$nTaxa
seqLists$nIntervals

# apply freqRat as an example analysis
sapply(seqLists$timeLists, freqRat)

# notice the zero and infinite freqRat estimates? What's going on?
freqRat(seqLists$timeLists[[4]], plot = TRUE)

# too few taxa of two or three interval durations for the ratio to work properly
# perhaps ignore these estimates

# with weighted selection of intervals
seqLists <- seqTimeList(rangesDisc1, nruns = 10, weightSampling = TRUE)

seqLists$nTaxa
seqLists$nIntervals
sapply(seqLists$timeLists, freqRat)

# didn't have much effect in this simulated example
```

setRootAge

Place a Non-Ultrametric Tree of Fossil Taxa on Absolute Time
Description

This function uses a table of fixed dates for operational-taxon-units (tip taxa) to calculate the absolute age of the root divergence for a tree with branch lengths, and then appends this root age to the tree as a $root_time element, and then outputs the tree. Function setRootAges is a wrapper for setRootAge for use with multiple trees in a "multiPhylo" object. This function was mainly written for dealing with trees of extinct taxa dated in units of absolute time from Bayesian analyses, such as with MrBayes, with trees scaled to time units by functions such as obtainDatedPosteriorTreesMrB.

Usage

setRootAge(tree, fixedAges = NULL)

setRootAges(trees, fixedAges = NULL)

Arguments

tree A phylogeny with branch lengths of type "phylo".

fixedAges A table of fixed ages for tip taxa, generally as a dataframe where the first column is of type character, and the second column is of type numeric. Such a table is automatically generated as an attribute of the output from obtainDatedPosteriorTreesMrB, when argument getFixedTimes = TRUE.

trees A list of type "multiPhylo" consisting of multiple phylogenetic trees with branch lengths.

Details

Trees of fossil taxa come with one issue rarely encountered by those dealing with molecular phylogenies: the absolute timing of when tips and divergences is not certain. With the vast majority of molecular phylogenies, it can be assumed the youngest tips occur at 0 time – i.e., the modern. This knowledge gives the tree an anchor for figuring out the absolute timing of events. Many programs and other software designed for depicting and analyzing phylogenetic hypotheses assumes such an apparent absolute time-scale (in R and elsewhere). A tree of Paleozoic brachiopods that include no extant members has no such anchor at time = 0, and such a default assumption in available software can be misleading. The $root_time protocol is intended to grant this absolute time-scale to a dated tree of fossil taxa, and is appended by most of the dating functions in package paleotree. However, trees dated by other approaches, such as via tip-dating in programs such as MrBayes and BEAST, will not have $root_time elements when read into R.

Value

The input tree is output, with a new $root_time element.

Author(s)

David W. Bapst
simFossilRecord

Full-Scale Simulations of the Fossil Record with Birth, Death and Sampling of Morphotaxa

Description

A complete birth-death-sampling branching simulator that captures morphological-taxon identity of lineages, as is typically discussed in models of paleontological data. This function allows for the use of precise point constraints to condition simulation run acceptance and can interpret complex character strings given as rate values for use in modeling complex processes of diversification and sampling.

Usage

```
simFossilRecord(p, q, r = 0, anag.rate = 0, prop.bifurc = 0,
    prop.cryptic = 0, modern.samp.prob = 1, startTaxa = 1, nruns = 1,
    maxAttempts = Inf, totalTime = c(0, 1000), nTotalTaxa = c(1, 1000),
    nExtant = c(0, 1000), nSamp = c(0, 1000), tolerance = 10^-4,
    maxStepTime = 0.01, shiftRoot4TimeSlice = "withExtantOnly",
    count.cryptic = FALSE, negRatesAsZero = TRUE, print.runs = FALSE,
    sortNames = FALSE, plot = FALSE)
```

Arguments

- `p, q, r, anag.rate`
  
  These parameters control the instantaneous ('per-capita') rates of branching, extinction, sampling and anagenesis, respectively. These can be given as a number equal to or greater than zero, or as a character string which will be interpreted as an algebraic equation. These equations can make use of three quantities which
will/may change throughout the simulation: the standing richness is N, the current time passed since the start of the simulation is T, the present duration of a given still-living lineage since its origination time is codeD, and the current branching rate is p (corresponding to the argument name p). Note that p cannot be used in equations for the branching rate itself; it is for making other rates relative to the branching rate. By default, the rates r and anag.rate are set to zero, so that the default simulator is a birth-death simulator. Rates set to Inf are treated as if 0. When a rate is set to 0, this event type will not occur in the simulation. Setting certain processes to zero, like sampling, may increase simulation efficiency, if the goal is a birth-death or pure-birth model. See documentation for argument negRatesAsZero about the treatment of rates that decrease below zero. Notation of branching, extinction and sampling rates as p, q, r follows what is typical for the paleobiology literature (e.g. Foote, 1997), not the Greek letters lambda, mu, phi found more typically in the biological literature (e.g. Stadler, 2009; Heath et al., 2014; Gavryushkina et al., 2014).

prop.cryptic, prop.bifurc
These parameters control (respectively) the proportion of branching events that have morphological differentiation, versus those that are cryptic (prop.cryptic) and the proportion of morphological branching events that are bifurcating, as opposed to budding. Both of these proportions must be a number between 0 and 1. By default, both are set to zero, meaning all branching events are events of budding cladogenesis. See description of the available models of morphological differentiation in the Description section.

modern.samp.prob
The probability that a taxon is sampled at the modern time (or, for timeSliceFossilRecord, the time at which the simulation data is slice). Must be a number between 0 and 1. If 1, all taxa that survive to the modern day (to the sliceTime) are sampled, if 0, none are.

startTaxa
Number of initial taxa to begin a simulation with. All will have the simulation start date listed as their time of origination.

nruns
Number of simulation datasets to accept, save and output. If nruns = 1, output will be a single object of class fossilRecordSimulation, and if nruns is greater than 1, a list will be output composed of nruns objects of class fossilRecordSimulation.

maxAttempts
Number of simulation attempts allowed before the simulation process is halted with an error message. Default is Inf.

totalTime, nTotalTaxa, nExtant, nSamp
These arguments represent stopping and acceptance conditions for simulation runs. They are respectively totalTime, the total length of the simulation in time-units, nTotalTaxa, the total number of taxa over the past evolutionary history of the clade, nExtant, the total number of extant taxa at the end of the simulation and nSamp the total number of sampled taxa (not counting extant taxa sampled at the modern day). These are used to determine when to end simulation runs, and whether to accept or reject them as output. They can be input as a vector of two numbers, representing minimum and maximum values of a range for accepted simulation runs (i.e. the simulation length can be between 0 and 1000 time-steps, by default), or as a single number, representing a
point condition (i.e. if $n_{\text{Samp}} = 100$ then the only simulation runs with exactly 100 taxa sampled will be output). Note that it is easy to set combinations of parameters and run conditions that are impossible to produce satisfactory input under, in which case `simFossilRecord` would run in a nonstop loop. How cryptic taxa are counted for the sake of these conditions is controlled by argument `count.cryptic`.

**tolerance**
A small number which defines a tiny interval for the sake of placing run-sampling dates before events and for use in determining whether a taxon is extant in `simFossilRecordMethods`.

**maxStepTime**
When rates are time-dependent (i.e. when parameters 'D' or 'T' are used in equations input for one of the four rate arguments), then protocol used by `simFossilRecord` of drawing waiting times to the next event could produce a serious mismatch of resulting process to the defined model, because the possibility of new events is only considered at the end of these waiting times. Instead, any time a waiting time greater than `maxStepTime` is selected, then instead no event occurs and a time-step equal to `maxStepTime` occurs instead, thus effectively discretizing the progression of time in the simulations run by `simFossilRecord`. Decreasing this value will increase accuracy (as the time-scale is effectively more discretized) but increase computation time, as the computer will need to stop and check rates to see if an event happened more often. Users should toggle this value relative to the time-dependent rate equations they input, relative to the rate of change in rates expected in time-dependent rates.

**shiftRoot4TimeSlice**
Should the dating of events be shifted, so that the date given for `slicetime` is now 0, or should the dates not be shifted, so that they remain on the same scale as the input? This argument accepts a logical TRUE or FALSE, but also accepts the string "withExtantOnly", which will only 'shift' the time-scale if living taxa are present, as determined by having ranges that overlap within `tolerance` of `sliceTime`.

**count.cryptic**
If TRUE, cryptic taxa are counted as separate taxa for conditioning limits that count a number of taxon units, such as `nTotalTaxa`, `nExtant` and `nSamp`. If FALSE (the default), then each cryptic complex (i.e. each distinguishable morphotaxon) is treated as a single taxon. See examples.

**negRatesAsZero**
A logical. Should rates calculated as a negative number cause the simulation to fail with an error message ( = FALSE) or should these be treated as zero (" = TRUE", the default). This is equivalent to saying that the `rate.as.used` = max(0, `rate.as.given`).

**print.runs**
If TRUE, prints the proportion of simulations accepted for output to the terminal.

**sortNames**
If TRUE, output taxonomic lists are sorted by the taxon names (thus sorting cryptic taxa together) rather than by taxon ID number (i.e. the order they were simulated in).

**plot**
If TRUE, plots the diversity curves of accepted simulations, including both the diversity curve of the true number of taxa and the diversity curve for the 'observed' (sampled) number of taxa.

**Details**
`simFossilRecord` simulates a birth-death-sampling branching process (ala Foote, 1997, 2000;
Stadler, 2009) in which lineages of organisms may branch, go extinct or be sampled thought a continuous time-interval, with the occurrence of these events modeled as Poisson process controlled by some set of instantaneous rates. This model is ultimately based on the birth-death model (Kendall, 1948; Nee, 2006), which is widely implemented in many R packages. Unlike other such typical branching simulators, this function enmeshes the lineage units within explicit models of how lineages are morphologically differentiated (Bapst, 2013). This is key to allow comparison to datasets from the fossil record, as morphotaxa are the basic units of paleontological diversity estimates and phylogenetic analyses.

Models of Morphological Differentiation and Branching (Cladogenesis and Anagenesis)

These models of morphological differentiation do not involve the direct simulation of morphological traits. Instead, morphotaxon identity is used as a proxy of the distinctiveness of lineages on morphological grounds, as if there was some hypothetical paleontologist attempting to taxonomically sort collections of specimens of these simulated lineages. Two lineages are either identical, and thus share the same morphotaxon identity, or they are distinct, and thus have separate morphotaxon identities. Morphological differentiation is assumed to be an instantaneous process for the purposes of this model, such that no intermediate could be uncovered.

Specifically, \texttt{simFossilRecord} allows for three types of binary branching events (here grouped under the term ‘cladogenesis’: ‘budding cladogenesis’, ‘bifurcating cladogenesis’ and ‘cryptic cladogenesis’, as well as for a fourth event-type, ‘anagenesis’ (see Wagner and Erwin, 1995; Foote, 1996, and Bapst, 2013, for further details). Budding, bifurcation and cryptic cladogenetic events all share in common that a single geneological lineage splits into two descendant lineages, but differ in the morphological differentiation of these child lineages relative to their parent. Under budding cladogenesis, only one of the child lineages becomes morphologically distinguishable from the parent, and thus the ancestral morphotaxon persists through the branching event as the child lineage that does not differentiate. Under bifurcating cladogenesis, both child lineages become immediately distinct from the ancestor, and thus two new morphotaxa appear while the ancestor terminates in an event known as ‘pseudoextinction’. Cryptic cladogenesis has no morphological differentiation: both child lineages are presumed to be indistinct from the ancestor and from each other, which means a hypothetical differentiation independent of any branching, such that a morphotaxon instantaneously transitions to a new morphotaxon identity, resulting in the pseudoextinction of the ancestral morphotaxon and the immediate ‘pseudospeciation’ of the child morphotaxon. The two morphotaxa do not overlap in time at all, as modeled here (contra to the models described by Ezard et al., 2012). For ease of following these cryptic lineages, cryptic cladogenetic events are treated in terms of data structure similarly to budding cladogenetic events, with one child lineage treated as a persistence of the ancestral lineage, and the other as a new morphologically indistinguishable lineage. This model of cryptic cladogenesis is ultimately based on the hierarchical birth-death model used by many authors for modeling patterns across paraphyletic higher taxa and the lower taxon units within them (e.g. Patzkowsky, 1995; Foote, 2012).

The occurrence of the various models is controlled by multiple arguments of \texttt{simFossilRecord}. The overall instantaneous rate of branching (cladogenesis) is controlled by argument \( p \), and the proportion of each type of cladogenesis controlled by arguments \( \text{prop.bifurc} \) and \( \text{prop.cryptic} \). \( \text{prop.cryptic} \) controls the overall probability that any branching event will be cryptic versus involving any morphological differentiation (budding or bifurcating). If \( \text{prop.cryptic} = 1 \), all branching events will be cryptic cladogenesis, and if \( \text{prop.cryptic} = 0 \), all branching events will involve morphological differentiation and none will be cryptic. \( \text{prop.bifurc} \) controls how many branching events that involve morphological differentiation (i.e. the inverse of \( \text{prop.cryptic} \)) are bifurcating, as opposed to budding cladogenesis. If \( \text{prop.bifurc} = 1 \), all morphologically-
differentiating branching events will be bifurcating cladogenesis, and if \( \text{prop.bifurc} = 0 \), all morphologically-differentiating branching events will be budding cladogenesis. Thus, for example, the probability of a given cladogenesis event being budding is given by:

\[
\text{Prob(budding cladogenesis at a branching event)} = (1 - \text{prop.cryptic}) \times (1 - \text{prop.bifurc})
\]

By default, \( \text{prop.cryptic} = 0 \) and \( \text{prop.bifurc} = 0 \), so all branching events by default will be instances of budding cladogenesis. Anagenesis is completely independent of these, controlled as its own Poisson process with an instantaneous rate defined by the argument \( \text{anag.rate} \). By default, this rate is set to zero and thus there is no anagenetic events without user intervention.

**Stopping Conditions and Acceptance Criteria for Simulations**

How forward-time simulations are generated, halted and whether they are accepted or not for output is a critical component of simulation design. Most uses of \text{simFossilRecord} will involve iteratively generating and analyzing multiple simulation runs. Runs are only accepted for output if they meet the conditioning criteria defined in the arguments, either matching point constraints or falling within range constraints. However, this requires separating the processes of halting simulation runs and accepting a run for output, particularly to avoid bias related to statistical sampling issues.

Hartmann et al. (2011) recently discovered a potential statistical artifact when branching simulations are conditioned on some number of taxa. Previously within \text{paleotree}, this was accounted for in the deprecated function \text{simFossilTaxa} by a complex arrangement of minimum and maximum constraints, and an (incorrect) presumption that allowing simulations to continue for a short distance after constraints were reached. This strategy is not applied here. Instead, \text{simFossilRecord} applies the General Sampling Algorithm presented by Hartmann et al. (or at least, a close variant). A simulation continues until extinction or some maximum time-constraint is reached, evaluated for intervals that match the set run conditions (e.g. \( n\text{Extant}, n\text{TotalTime} \)) and, if some interval or set of intervals matches the run conditions, a date is randomly sampled from within this interval/interval. The simulation is then cut at this date using the \text{timeSliceFossilRecord} function, and saved as an accepted run. The simulation data is otherwise discarded and then a new simulation initiated (thus, at most, only one simulated dataset is accepted from one simulation run).

Thus, accepted simulations runs should reflect unbiased samples of evolutionary histories that precisely match the input constraints, which can be very precise, unlike how stopping and acceptance conditions were handled in the previous (deprecated) \text{simFossilTaxa} function. Of course, selecting very precise constraints that are very unlikely or impossible given other model parameters may take considerable computation time to find acceptable simulation runs, or effectively never find any acceptable simulation runs.

**On Time-Scale Used in Output**

Dates given in the output are on an reversed absolute time-scale; i.e. time decreases going from the past to the future, as is typical in paleontological uses of time (as time before present) and as for most function in package \text{paleotree}. The endpoints of the time-scale are decided by details of the simulation and can be modified by several arguments. By default (with \text{shiftRoot4TimeSlice} = "withExtantOnly"), any simulation run that is accepted with extant taxa will have zero as the end-time (i.e. when those taxa are extant), as zero is the typical time assigned to the modern day in empirical studies. If a simulation ends with all taxa extinct, however, then instead the start-time of a run (i.e. when the run initiates with starting taxa) will be maximum value assigned to the conditioning argument \text{totalTime}. If \text{shiftRoot4TimeSlice} = FALSE then the start-time of the run will always be this maximum value for \text{totalTime}, and any extant taxa will stop at some time greater than zero.
Value

simFossilRecord returns either a single object of class fossilRecordSimulation or a list of multiple such objects, depending on whether \( nruns \) was 1 or more.

An object of class fossilRecordSimulation consists of a list object composed of multiple elements, each of which is data for 'one taxon'. Each data element for each taxon is itself a list, composed of two elements: the first describes vital information about the taxon unit, and the second describes the sampling times of each taxon.

The first element of the list (named $taxa.data) is a distinctive six-element vector composed of numbers (some are nominally integers, but not all, so all are stored as double-precision integers) with the following field names:

- **taxon.id** The ID number of this particular taxon-unit.
- **ancestor.id** The ID number of the ancestral taxon-unit. The initial taxa in a simulation will be listed with NA as their ancestor.
- **orig.time** True time of origination for a taxon-unit in absolute time.
- **ext.time** True time of extinction for a taxon-unit in absolute time. Extant taxa will be listed with an ext.time of the run-end time of the simulation run, which for simulations with extant taxa is 0 by default (but this may be modified using argument shiftRoot4TimeSlice).
- **still.alive** Indicates whether a taxon-unit is 'still alive' or not: '1' indicates the taxon-unit is extant, '0' indicates the taxon-unit is extinct
- **looks.like** The ID number of the first morphotaxon in a dataset that 'looks like' this taxon-unit; i.e. belongs to the same multi-lineage cryptic complex. Taxa that are morphologically distinct from any previous lineage will have their taxon.id match their looks.like. Thus, this column is rather uninformative unless cryptic cladogenesis occurred in a simulation.

The second element for each taxon-unit is a vector of sampling times, creatively named $sampling.times, with each value representing a data in absolute time when that taxon was sampled in the simulated fossil record. If a taxon was never sampled, this vector is an empty numeric vector of length = 0.

As is typical for paleontological uses of absolute time, absolute time in these simulations is always decreasing toward the modern; i.e. an absolute date of 50 means a point in time which is 50 time-units before the present-day, if the present-day is zero (the default, but see argument shiftRoot4TimeSlice).

Each individual element of a fossilRecordSimulation list object is named, generally of the form "t1" and "t2", where the number is the taxon.id. Cryptic taxa are instead named in the form of "t1.2" and "t5.3", where the first number is the taxon which they are a cryptic descendant of (looks.like) and the second number, after the period, is the order of appearance of lineage units in that cryptic complex. For example, for "t5.3", the first number is the taxon.id and the second number communicates that this is the third lineage to appear in this cryptic complex.

Author(s)

David W. Bapst, inspired by code written by Peter Smits.
References


**See Also**

*simFossilRecordMethods*

This function essentially replaces and adds to all functionality of the deprecated paleotree functions *simFossilTaxa, simFossilTaxaSRCond, simPaleoTrees*, as well as the combined used of *simFossilTaxa and sampleRanges* for most models of sampling.
Examples

set.seed(2)

# quick birth-death-sampling run with 1 run, 50 taxa

record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, nruns = 1,
nTotalTaxa = 50, plot = TRUE)

# examining multiple runs of simulations

# example of repeated pure birth simulations over 50 time-units
records <- simFossilRecord(p = 0.1, q = 0, nruns = 10,
totalTime = 50, plot = TRUE)
# plot multiple diversity curves on a log scale
records <- lapply(records, fossilRecord2fossilTaxa)
multiDiv(records, plotMultCurves = TRUE, plotLogRich = TRUE)
# histogram of total number of taxa
hist(sapply(records, nrow))

# example of repeated birth-death-sampling simulations over 50 time-units
records <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, nruns = 10,
totalTime = 50, plot = TRUE)
records <- lapply(records, fossilRecord2fossilTaxa)
multiDiv(records, plotMultCurves = TRUE)

# like above, but conditioned instead on having 10 extant taxa
# between 1 and 100 time-units
set.seed(1)

records <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, nruns = 10,
totalTime = c(1,300), nExtant = 10, plot = TRUE)
records <- lapply(records, fossilRecord2fossilTaxa)
multiDiv(records, plotMultCurves = TRUE)

# How probable were the runs I accepted?
# The effect of conditions

set.seed(1)

# Let's look at an example of a birth-death process
# with high extinction relative to branching
# use default run conditions (barely any conditioning)
# use print.runs to look at acceptance probability
records <- simFossilRecord(p = 0.1, q = 0.8, nruns = 10,
print.runs = TRUE, plot = TRUE)
# 10 runs accepted from a total of 10 !

# now let's give much more stringent run conditions
# require 3 extant taxa at minimum, 5 taxa total minimum
records <- simFossilRecord(p = 0.1, q = 0.8, nrungs = 10, nExtant = c(3,100), nTotalTaxa = c(5,100), print.runs = TRUE, plot = TRUE)
# thousands of simulations to just obtain 10 acceptable runs!
# most ended in extinction before minimums were hit

# beware analysis of simulated where acceptance conditions
# are too stringent: your data will be a 'special case'
# of the simulation parameters
# it will also take you a long time to generate reasonable
# numbers of replicates for whatever analysis you are doing

# TLDR: You should look at print.runs = TRUE

##############################
# Using the rate equation-input for complex diversification models
# First up... Diversity Dependent Models!
# Let’s try Diversity-Dependent Branching over 50 time-units

# first, let’s write the rate equation
# We’ll use the diversity dependent rate equation model
# from Ettienne et al. 2012 as an example here
# Under this equation, p = q at carrying capacity K
# Many others are possible!
# Note that we don’t need to use max(0,rate) as negative rates
# are converted to zero by default, as controlled by
# the argument negRatesAsZero

# From Ettienne et al.
# lambda = lambda0 - (lambda0 - mu)*(n/K)
# lambda and mu are branching rate and extinction rate
# lambda and mu == p and q in paleotree (i.e. Foote convention)
# lambda0 is the branching rate at richness = 0
# K is the carrying capacity
# n is the richness

# 'N' is the algebra symbol for standing taxonomic richness
# for simFossilRecord’s simulation capabilities
# also branching rate cannot reference extinction rate
# we’ll have to set lambda0, mu and K in the rate equation directly

lambda0 <- 0.3 # branching rate at 0 richness in Ltu
K <- 40 # carrying capacity
mu <- 0.1 # extinction rate will 0.1 Ltu ( = 1/3 of lambda0)

# technically, mu here represents the lambda at richness = K
# i.e. lambdaK
# Ettienne et al. are just implicitly saying that the carrying capacity
# is the richness at which lambda == mu

# construct the equation programmatically using paste0
branchingRateEq <- paste0(lamda0, "-\(\frac{\mu}{K}\)\), \(\mu\)\), \(\mu_0\)\), \(\mu_0\)\), \(\mu_0\)\), \(\mu_0\)
# and take a look at it...
branchingRateEq
# its a thing of beauty, folks

# now let's try it
records <- simFossilRecord(p = branchingRateEq, q = mu, nruns = 3, totalTime = 100, plot = TRUE, print.runs = TRUE)
records <- lapply(records, fossilRecord2fossilTaxa)
multiDiv(records, plotMultCurves = TRUE)
# those are some happy little diversity plateaus!

# now let's do diversity-dependent extinction

# let's slightly modify the model from Ettiene et al.
# \(\mu = \mu_0 + (\mu_0 - \mu_K)\)\)
mu0 <- 0.001 # \(\mu\) at \(n = 0\)
muk <- 0.1 # \(\mu\) at \(n = K\) (should be equal to lambda at \(K\))
K <- 40
lambda <- muk # equal to muk

# construct the equation programmatically using paste0
extrateEq <- paste0(mu0, "-\(\mu_0\)\), \(\mu_0\)\), \(\mu_0\)\), \(\mu_0\)
extrateEq

# now let's try it
records <- simFossilRecord(p = lambda, q = extrateEq, nruns = 3, totalTime = 100, plot = TRUE, print.runs = TRUE)
records <- lapply(records, fossilRecord2fossilTaxa)
multiDiv(records, plotMultCurves = TRUE)

# these plateaus looks a little more spiky
#( maybe there is more turnover at \(K?\) )
# also, it took a longer for the rapid rise to occur

# Now let's try an example with time-dependent origination
# and extinction constrained to equal origination

# Note! Use of time-dependent parameters "D" and "T" may
# result in slower than normal simulation run times
# as the time-scale has to be discretized; see
# info for argument maxTimeStep above

# First, let's define a time-dependent rate equation
# "T" is the symbol for time passed
timeEquation <- "0.4-\(0.007\times T\)"

#in this equation, 0.4 is the rate at time = 0
# and it will decrease by 0.007 with every time-unit
# at time = 50, the final rate will be 0.05
# We can easily make it so extinction is always equal to branching rate
# "P" is the algebraic equivalent for "branching rate" in simFossilRecord

# now let's try it
records <- simFossilRecord(p = timeEquation, q = "P", nruns = 3,
totalTime = 50, plot = TRUE, print.runs = TRUE)
records <- lapply(records, fossilRecord2fossilTaxa)
multiDiv(records, plotMultCurves = TRUE)

# high variability that seems to then smooth out as turnover decreases

# And duration what about duration-dependent processes?
# let's do a duration-dep extinction equation:
durDepExt <- "0.01+(0.01*D)"

# okay, let's take it for a spin
records <- simFossilRecord(p = 0.1, q = durDepExt, nruns = 3,
totalTime = 50, plot = TRUE, print.runs = TRUE)
records <- lapply(records, fossilRecord2fossilTaxa)
multiDiv(records, plotMultCurves = TRUE)

# creates runs full of short lived taxa

# Some more stuff to do with rate formulae!

# The formulae input method for rates allows
# for the rate to be a random variable

# For example, we could constantly redraw
# the branching rate from an exponential

record <- simFossilRecord(
  p = "rexp(n = 1,rate = 10)",
  q = 0.1, r = 0.1, nruns = 1,
nTotalTaxa = 50, plot = TRUE)

# Setting up specific time-variable rates can be laborious though
# e.g. one rate during this 10 unit interval,
# another during this interval, etc
# The problem is setting this up within a fixed function

# Worked Example
# What if we want to draw a new rate from a
# lognormal distribution every 10 time units?

# Need to randomly draw these rates *before* running simFossilTaxa
# This means also that we will need to individually do each simFossilTaxa run
# since the rates are drawn outside of simFossilTaxa

# Get some reasonable log normal rates:
rates <- 0.1+rlnorm(100,meanlog = 1,sdlog = 1)/100

# Now paste it into a formulae that describes a function that
# will change the rate output every 10 time units
rateEquation <- paste0("c("paste0(rates,collapse = ","))[1+(T/%10)]")
# and let's run it
record <- simFossilRecord(p = rateEquation, q = 0.1, r = 0.1, nruns = 1, 
totalTime = c(30,40), plot = TRUE)

# Speciation Modes
# Some examples of varying the 'speciation modes' in simFossilRecord

# The default is pure budding cladogenesis
# anag.rate = prop.bifurc = prop.cryptic = 0
# let's just set those for the moment anyway
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, 
anag.rate = 0, prop.bifurc = 0, prop.cryptic = 0, 
nruns = 1, nTotalTaxa = c(20,30), nExtant = 0, plot = TRUE)

#convert and plot phylogeny
# note this will not reflect the 'budding' pattern
# branching events will just appear like bifurcation
# it's a typical convention for phylogeny plotting
converted <- fossilRecord2FossilTaxa(record)
tree <- taxa2phylo(converted,plot = TRUE)

#now, an example of pure bifurcation
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, 
anag.rate = 0, prop.bifurc = 1, prop.cryptic = 0, 
nruns = 1, nTotalTaxa = c(20,30), nExtant = 0)
tree <- taxa2phylo(fossilRecord2FossilTaxa(record), plot = TRUE)

# all the short branches are due to ancestors that terminate
# via pseudoextinction at bifurcation events

# an example with anagenesis = branching
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, 
anag.rate = 0.1, prop.bifurc = 0, prop.cryptic = 0, 
nruns = 1, nTotalTaxa = c(20,30), nExtant = 0)
tree <- taxa2phylo(fossilRecord2FossilTaxa(record), plot = TRUE)
# lots of pseudoextinction

# an example with anagenesis, pure bifurcation
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, 
anag.rate = 0.1, prop.bifurc = 1, prop.cryptic = 0, 
nruns = 1, nTotalTaxa = c(20,30), nExtant = 0)
tree <- taxa2phylo(fossilRecord2FossilTaxa(record), plot = TRUE)
# lots and lots of pseudoextinction

# an example with half cryptic speciation
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, 
anag.rate = 0, prop.bifurc = 0, prop.cryptic = 0.5, 
nruns = 1, nTotalTaxa = c(20,30), nExtant = 0)
tree <- taxa2phylo(fossilRecord2FossilTaxa(record), plot = TRUE)
# notice that the tree has many more than the maximum of 30 tips:
# that's because the cryptic taxa are not counted as
# separate taxa by default, as controlled by count.cryptic

# an example with anagenesis, bifurcation, cryptic speciation
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1,
anag.rate = 0.1, prop.bifurc = 0.5, prop.cryptic = 0.5,
nruns = 1, ntotalTaxa = c(20, 30), nExtant = 0)
tree <- taxa2phylo(fossilRecord2fossilTaxa(record), plot = TRUE)
# note in this case, 50% of branching is cryptic
# 25% is bifurcation, 25% is budding

# an example with anagenesis, pure cryptic speciation
# morphotaxon identity will thus be entirely indep of branching!
# I wonder if this is what is really going on, sometimes...
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1,
anag.rate = 0.1, prop.bifurc = 0, prop.cryptic = 1,
nruns = 1, ntotalTaxa = c(20, 30), nExtant = 0)
tree <- taxa2phylo(fossilRecord2fossilTaxa(record), plot = TRUE)

# merging cryptic taxa when all speciation is cryptic
set.seed(1)
record <- simFossilRecord(p = 0.1,
q = 0.1, r = 0.1,
prop.crypt = 1,
totalTime = 50, plot = TRUE)
# there looks like there is only a single taxon, but...
length(record) # actual number of cryptic lineages

#################
# playing with count.cryptic with simulations of pure cryptic speciation

# can choose to condition on total morphologically-distinguishable taxa
# or total taxa including cryptic taxa with count.cryptic = FALSE

# an example with pure cryptic speciation with count.cryptic = TRUE
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1,
anag.rate = 0, prop.bifurc = 0, prop.cryptic = 1,
nruns = 1, totalTime = 50, ntotalTaxa = c(10, 100), count.cryptic = TRUE)
tree <- taxa2phylo(fossilRecord2fossilTaxa(record))
plot(tree); axisPhylo()
# notice how the tip labels indicate all are the same morphotaxon

# we'll replace the # of taxa constraints with a time constraint
# or else the count.cryptic = FALSE simulation will never end!

# an example with pure cryptic speciation with count.cryptic = FALSE
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1,
anag.rate = 0, prop.bifurc = 0, prop.cryptic = 1,
nruns = 1, totalTime = 50, count.cryptic = FALSE)
tree <- taxa2phylo(fossilRecord2fossilTaxa(record))
plot(tree); axisPhylo()
# let's look at numbers of taxa returned when varying count.cryptic  
# with prop.cryptic = 0.5

# simple simulation going for 50 total taxa

# first, count.cryptic = FALSE (default)
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1,  
anag.rate = 0, prop.bifurc = 0, prop.cryptic = 0.5,  
nruns = 1, nTotalTaxa = 50, count.cryptic = FALSE)
taxa <- fossilRecord2FossilTaxa(record)
ntaxa <- nrow(taxa)  # number of lineages (inc. cryptic)
length(unique(taxa[,6]))  # number of morph-distinguishable taxa

# and count.cryptic = TRUE
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1,  
anag.rate = 0, prop.bifurc = 0, prop.cryptic = 0.5,  
nruns = 1, nExtant = 10, totalTime = c(1,100), count.cryptic = TRUE)
taxa <- fossilRecord2FossilTaxa(record)
sum(taxa[,5])  # number of still-living lineages (inc. cryptic)
length(unique(taxa[taxa[,5] == 1,6]))  # number of still-living morph-dist. taxa

# okay...
# now let's try with 50 extant taxa

# first, count.cryptic = FALSE (default)
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1,  
anag.rate = 0, prop.bifurc = 0, prop.cryptic = 0.5,  
nruns = 1, nExtant = 10, totalTime = c(1,100), count.cryptic = FALSE)
taxa <- fossilRecord2FossilTaxa(record)
sum(taxa[,5])  # number of still-living lineages (inc. cryptic)
length(unique(taxa[taxa[,5] == 1,6]))  # number of still-living morph-dist. taxa

# and count.cryptic = TRUE
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1,  
anag.rate = 0, prop.bifurc = 0, prop.cryptic = 0.5,  
nruns = 1, nExtant = 10, totalTime = c(1,100), count.cryptic = TRUE)
taxa <- fossilRecord2FossilTaxa(record)
sum(taxa[,5])  # number of still-living lineages (inc. cryptic)
length(unique(taxa[taxa[,5] == 1,6]))  # number of still-living morph-dist. taxa

# an example using startTaxa to have more initial taxa
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, nruns = 1,  
nTotalTaxa = 100, startTaxa = 20, plot = TRUE)

# Using run conditions

# Users can generate datasets that meet multiple conditions:  
# such as time, number of total taxa, extant taxa, sampled taxa
# These can be set as point conditions or ranges

# let's set time = 10-100 units, total taxa = 30-40, extant = 10  
# and look at acceptance rates with print.run
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, nruns = 1,  
totalTime = c(10,100), nTotalTaxa = c(30,40), nExtant = 10,  
print.runs = TRUE, plot = TRUE)

# let's make the constraints on total taxa a little tighter 
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, nruns = 1,  
totalTime = c(50,100), nTotalTaxa = 30, nExtant = 10,  
print.runs = TRUE, plot = TRUE)
# still okay acceptance rates

# alright, now let's add a constraint on sampled taxa 
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, nruns = 1,  
totalTime = c(50,100), nTotalTaxa = 30, nExtant = 10,  
nSamp = 15, print.runs = TRUE, plot = TRUE)
# still okay acceptance rates

# we can be really odd and condition on having a single taxon 
set.seed(1)
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, nruns = 1,  
totalTime = c(10,20), plot = TRUE)

# Simulations of entirely extinct taxa

# Typically, a user may want to condition on a precise 
# number of sampled taxa in an all-extinct simulation
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, nruns = 1,  
nTotalTaxa = c(1,100), nExtant = 0, nSamp = 20,  
print.runs = TRUE, plot = TRUE)

# Note that when simulations don't include 
# sampling or extant taxa, the plot 
# functionality changes
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0, nruns = 1,  
nExtant = 0, print.runs = TRUE, plot = TRUE)
# something similar happens when there is no sampling
# and there are extant taxa but they aren't sampled
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0, nruns = 1,  
nExtant = 10, nTotalTaxa = 100, modern.samp.prob = 0,  
print.runs = TRUE, plot = TRUE)

# We can set up a test to make sure that no extant taxa somehow get
# returned in many simulations with extinct-only conditioning:
res <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1,  
nTotalTaxa = 10, nExtant = 0, nruns = 1000, plot = TRUE)
anyLive <- any(sapply(res,function(z)
Methods for Editing or Converting Output from *simFossilRecord*

**Description**

These are a set of functions available for manipulating, translating and editing the objects of class *fossilRecordSimulation* output from function *simFossilRecord*.

**Usage**

```r

timeSliceFossilRecord(fossilRecord, sliceTime, shiftRoot4TimeSlice = FALSE, modern.samp.prob = 1, tolerance = 10^-4)

fossilRecord2fossilTaxa(fossilRecord)

fossilTaxa2fossilRecord(fossilTaxa)

fossilRecord2fossilRanges(fossilRecord, merge.cryptic = TRUE, ranges.only = TRUE)
```

**Arguments**

- **fossilRecord**
  A list object output by *simFossilRecord*, often composed of multiple elements, each of which is data for ’one taxon’, with the first element being a distinctive six-element vector composed of numbers, corresponding to the six fields in tables output by the deprecated function *simFossilTaxa*.

- **sliceTime**
  The date to slice the *simFossilRecord* output at, given in time-units before the modern, on the same scale as the input *fossilRecord*.

- **shiftRoot4TimeSlice**
  Should the dating of events be shifted, so that the date given for *sliceTime* is now 0, or should the dates not be shifted, so that they remain on the same scale as the input? This argument accepts a logical TRUE or FALSE, but also accepts the string ”withExtantOnly”, which will only ’shift’ the time-scale if living taxa are present, as determined by having ranges that overlap within tolerance of *sliceTime*. 

```r

any(sapply(z,function(x) x[[1]][5] == 1)))
# test if any are still alive
if(anyLive){
  stop("Runs have extant taxa under conditioning for none")
}
```
modern.samp.prob

The probability that a taxon is sampled at the modern time (or, for timeSliceFossilRecord, the time at which the simulation data is slice). Must be a number between 0 and 1. If 1, all taxa that survive to the modern day (to the sliceTime) are sampled, if 0, none are.

tolerance

A small number which sets a range around the sliceTime within which taxa will be considered extant for the purposes of output.

fossilTaxa

A fossilTaxa object, composed of a table containing information on the true first and last appearance times of taxa, as well as their ancestor-descendant relationships.

merge.cryptic

If TRUE, sampling events for cryptic taxon-units (i.e. those in the same cryptic complex) will be merged into sampling events for a single taxon-unit (with the name of the first taxon in that cryptic complex).

ranges.only

If TRUE (the default), fossilRecord2fossilRanges will return the dates of the first and last sampled occurrences of each taxon-unit (i.e. the stratigraphic range of each taxon). If FALSE, instead a list will be output, with each element being a vector of dates for all sampling events of each taxon-unit.

Details

These functions exist to manipulate fossilRecordSimulation objects output from simFossilRecord, particularly so that they can be interfaced with functions in library paleotree in the same way that output from the deprecated 'legacy' simulation function simFossilTaxa was used.

timeSliceFossilRecord takes a given fossilRecordSimulation object and 'slices' the data to remove any events that occur after the given sliceTime and make it so any taxa still alive as of sliceTime are now listed as extant.

fossilRecord2fossilTaxa converts a fossilRecordSimulation object to the flat table format of taxon data as was originally output by deprecated function simFossilTaxa, and can be taken as input by a number of paleotree functions such as sampleRanges, taxa2phylo and taxa2cladogram.

fossilTaxa2fossilRecord does the reverse, converting a simFossilTaxa table into a fossilRecordSimulation list object, but returns a fossilRecordSimulation object that considers each species as un-sampled (as sampling information is not contained within a simFossilTaxa table).

fossilRecord2fossilRanges converts a fossilRecordSimulation object to the flat table format of observed taxon ranges, as is typically output by processing simFossilRecord simulation output with paleotree function sampleRanges.

Value

Depends on the function and the arguments given. See Details.

Author(s)

David W. Bapst

See Also

simFossilRecord
Examples

```r
set.seed(44)
record <- simFossilRecord(p = 0.1, q = 0.1, r = 0.1, nrugs = 1,
                          nTotalTaxa = c(20, 30), nExtant = 0, plot = TRUE)

# time-slicing

# let's try slicing this record at 940 time-units
slicedRecord <- timeSliceFossilRecord(fossilRecord = record, sliceTime = 940)
# and let's plot it
divCurveFossilRecordSim(slicedRecord)

# now with shiftRoot4TimeSlice = TRUE to shift the root age
slicedRecord <- timeSliceFossilRecord(fossilRecord = record, sliceTime = 940,
                                       shiftRoot4TimeSlice = TRUE)
# and let's plot it
divCurveFossilRecordSim(slicedRecord)

# plot look a little different due to how axis limits are treated...
# notice that in both, 'modern' (extant) taxa are sampled with probability = 1
# let's try it again, make that probability = 0

# now with shiftRoot4TimeSlice = TRUE
slicedRecord <- timeSliceFossilRecord(fossilRecord = record, sliceTime = 940,
                                       shiftRoot4TimeSlice = TRUE, modern.samp.prob = 0)
# and let's plot it
divCurveFossilRecordSim(slicedRecord)

# converting to taxa objects and observed ranges

# convert to taxa data
taxa <- fossilRecord2FossilTaxa(record)
# convert to ranges
ranges <- fossilRecord2FossilRanges(record)

# plot diversity curves with multiDiv
multiDiv(list(taxa, ranges), plotMultiCurves = TRUE)
# should look a lot like what we got earlier

# get the cladogram we'd obtain for these taxa with taxa2cladogram
cladogram <- taxa2cladogram(taxa, plot = TRUE)

# now get the time-scaled phylogenies with taxa2phylo

# first, with tips extending to the true times of extinction
treeExt <- taxa2phylo(taxa, plot = TRUE)

# now, with tips extending to the first appearance dates (FADs) of taxa
# get the FADs from the ranges
```
SongZhangDicrano

Cladistic Data for Dicranograptid Graptolites from Song and Zhang (2014)

Description

Character matrix and two cladograms for 13 dicranograptid (and outgroup) graptoloids, taken from Song and Zhang (2014). Included here for use with functions related to character change.

Format

Loading this dataset adds two objects to the R environment. charMatDicrano is a data.frame object composed of multiple factors, with NA values representing missing values (states coded as '?'), read in with readNexus from package phylobase. cladogramDicranoX12 and cladogramDicranoX13 are both cladograms, formatted as phylo class objects for use with package ape, without branch-lengths (as these was are, respectively, consensus tree and a maximum-parsimony tree from separate maximum-parsimony analyses).

Details

This example dataset is composed of a small cladistic character data for 13 taxa and 24 characters, taken from Song and Zhang (2014). Note that character 22 is a biostratigraphic character, which was not included in all analyses by Song and Zhang.

The first included cladogram cladogramDicranoX12 is the majority-rule consensus of a maximum-parsimony analysis on 12 taxa (excluding on taxa with incompletely known anatomy) with 24 characters, including a biostratigraphic character. This tree is included here as, among the four trees depicted, it appeared to be the basis for the majority of Song and Zhang’s discussion of dicranograptid systematics.

The second cladogram cladogramDicranoX13 is a maximum-parsimony tree found by a maximum-parsimony analysis of 13 taxa with 24 characters, including a biostratigraphic character. This tree is much more resolved than the alternative majority-rule cladogram for 12 taxa.

The matrix and both trees were entered by hand from their flat graphic depiction in Song and Zhang’s manuscript.

Source

Examples

data(SongZhangDicrano)

# Examining morphospace with a distance matrix

# calculate a distance matrix from the morph character data
char <- charMatDicrano[,,-22] # remove strat character
cDist <- matrix(nrow(char),nrow(char))
rownames(cDist) <- colnames(charDist) <- rownames(char)
for(i in 1:nrow(char)){for(j in 1:nrow(char)){
  charDist[i,j] <- length(charDist)/length(charDist)
}
}

###
# PCO of character distance matrix

# can apply PCO (use lingoes correction to account for negative values
# resulting from non-euclidean matrix
pco_res <- pcoa(charDist,correction = "lingoes")

# relative corrected eigenvalues
rel_corr_eig <- pco_res$values$Rel_corr_eig
layout(1:2)
plot(rel_corr_eig)
# cumulative
plot(cumsum(rel_corr_eig))

# well let's look at those PCO axes anyway
layout(1)
pco_axes <- pco_res$vectors
plot(pco_axes[,1],pco_axes[,2],pch = 16,
  xlab = paste("PCO Axis 1, Rel. Corr. Eigenvalue = ",round(rel_corr_eig[1,3]),

####
# plot 12 taxon majority rule tree from Song and Zhang
plot(cladogramDicranoX12,
  main = "MajRule_24charX12Taxa_wBiostratChar")

# plot 13 taxon MPT
plot(cladogramDicranoX13,
taxa2cladogram

Convert Simulated Taxon Data into a Cladogram

Description

Convert ancestor-descendant relationships of taxa into an 'ideal' unscaled cladogram, where taxa that could share true synapomorphies are arranged into nested clades.

Usage

taxa2cladogram(taxad, drop.cryptic = FALSE, plot = FALSE)

Arguments

taxad A five-column matrix of taxonomic data, as output by simFossilRecord, after transformation with function fossilRecord2fossilTaxa.
drop.cryptic Should cryptic species be dropped (except for the first)? Not dropped by default.
plot Should the result be plotted?

Details

This function simulates an ideal cladistic process, where the relationships of a set of morphologically static taxa is resolved into a set of nested hierarchical relationships (a standard cladogram), as much as would be expected given the input relationships among those taxa. taxa2cladogram uses information on the ancestor-descendant relationships of a bunch of taxa and constructs an unscaled cladogram of the hierarchically-nesting relationships among those taxa. There’s no actual cladistics going on, this is just a simulation of that process. If there is any chance that a set of taxa could be
resolved into a set of nested relationships given their ancestor-descendant relationships, they will be resolved so in the output of taxa2cladogram. No morphological characters are considered, we just assume that if there is a nesting relationship, then it could be resolved as such. This makes it the "ideal" cladogram of a simulated clade.

The result will probably not be fully resolved, as including both ancestor and descendant taxa will generally make it impossible to produce a fully nesting system of relationships. For example, consider a set of three morphologically-static taxa where the first is an ancestor (either direct or indirect, ala Foote, 1996) of both the second and third. If we imagine an ideal cladistic analysis of the morphological characters of those three taxa, this set of taxa will be unable to be broken up into bifurcating-nested relationships and thus result in a polytomy. Any set of ancestor-descendant relationships will have many of these, as some ancestors must have more than one descendant for the clade to diversify, as noted by Wagner and Erwin, 1995.

If there are cryptic taxa present in the output from simFossilRecord, these and any of their morphologically distinguishable descendants are collapsed into a polytomy to simulate the expected pattern of lack of phylogenetic resolution. In addition to this merging, cryptic taxa can be dropped via the argument drop.cryptic, such that only the first 'species' of each cryptic taxon assemblage is listed among the tip taxa (what we would actually expect to obtain, as we would not recognize cryptic taxa to be treated as different OTUs). By default, cryptic taxa are not dropped so that the same number of taxa as in the simulated data is retained.

Value

The resulting phylogeny without branch lengths is output as an object of class phylo.

The tip labels are the rownames from the simulation input; see documentation for simFossilRecord and fossilRecord2fossilTaxa documentation for details.

Author(s)

David W. Bapst

References


See Also

simFossilRecord, taxa2phylo, fossilRecord2fossilTaxa

Examples

set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
APolynomial = c(30,40), nExtant = 0)
taxa <- fossilRecord2fossilTaxa(record)
#Let's use taxa2cladogram to get the 'ideal' cladogram of the taxa
layout(1:2)
taxa2phylo <- taxa2cladogram(taxa, plot = TRUE)
# compare the "real" time-scaled plot
# to the 'ideal' cladogram

tree <- taxa2phylo(taxa, plot = TRUE)

# testing with cryptic speciation
recordCrypt <- simFossilRecord(p = 0.1, q = 0.1, prop.cryptic = 0.5, nruns = 1,
nTotalTaxa = c(30,40), nExtant = 0)
taxaCrypt <- fossilRecord2FossilTaxa(recordCrypt)
layout(1:2)
parOrig <- par(no.readonly = TRUE)
par(mar = c(0,0,0,0))
clandoCrypt1 <- taxa2cladogram(taxaCrypt, drop.cryptic = FALSE)
plot(clandoCrypt1)
clandoCrypt2 <- taxa2cladogram(taxaCrypt, drop.cryptic = TRUE)
plot(clandoCrypt2)

# reset plotting
par(parOrig)
layout(1)

---

taxa2phylo

**Convert Simulated Taxon Data into a Phylogeny**

**Description**

Converts temporal and ancestor-descendant relationships of taxa into a time-scaled phylogeny.

**Usage**

taxa2phylo(taxad, obs_time = NULL, plot = FALSE)

**Arguments**

taxad A five-column matrix of taxonomic data, as output by fossilRecord2fossilTaxa via simulations produced using simFossilRecord

obs_time A vector of per-taxon times of observation which must be in the same order of taxa as in the object taxad; if NULL, the LADs (column 4) in taxad2 are used

plot Plot the resulting phylogeny?

**Details**

As described in the documentation for taxa2cladogram, the relationships among morphotaxa in the fossil record are difficult to describe in terms of traditional phylogenies. One possibility is to arbitrarily choose particular instantaneous points of time in the range of some taxa and describe the time-scaled relationships of the populations present at those dates. This is the tactic used by taxa2phylo.
By default, the dates selected ('obs-time' argument) are the last occurrences of the taxon, so a simple use of this function will produce a time-scaled tree which describes the relationships of the populations present at the last occurrence of each taxon in the sampled data. Alternatively, obs_time can be supplied with different dates within the taxon ranges.

All data relating to when static morpho-taxon appear or disappear in the record is lost; branching points will be the actual time of speciation, which (under budding) will often be in the middle of the temporal range of a taxon.

Cryptic taxa are not dropped or merged as can be done with taxa2cladogram. The purpose of taxa2phylo is to obtain the 'true' pattern of evolution for the observation times, independent of what we might actually be able to recover, for the purpose of comparing in simulation analyses.

As with many functions in the paleotree library, absolute time is always decreasing, i.e. the present day is zero.

Value

The resulting phylogeny with branch lengths is output as an object of class phylo. This function will output trees with the element $root.time, which is the time of the root divergence in absolute time.

The tip labels are the rownames from the simulation input; see documentation for simFossilRecord and fossilRecord2fossilTaxa documentation for details.

Note

DO NOT use this function to time-scale a real tree for a real dataset. It assumes you know the divergence/speciation times of the branching nodes and relationships perfectly, which is almost impossible given the undersampled nature of the fossil record. Use timePaleoPhy or cal3TimePaleoPhy instead.

DO use this function when doing simulations and you want to make a tree of the 'true' history, such as for simulating trait evolution along phylogenetic branches.

Unlike taxa2cladogram, this function does not merge cryptic taxa in output from simFossilRecord (via fossilRecord2fossilTaxa) and I do not offer an option to secondarily drop them. The tip labels should provide the necessary information for users to drop such taxa, however. See simFossilRecord.

Author(s)

David W. Bapst

See Also

simFossilRecord, taxa2cladogram, fossilRecord2fossilTaxa

Examples

```r
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
                          nTotalTaxa = c(30,40), nExtant = 0)
```
taxa <- fossilRecord2fossilTaxa(record)
#let's use taxa2cladogram to get the 'ideal' cladogram of the taxa
tree <- taxa2phylo(taxa)
phyloDiv(tree)

#now a phylogeny with tips placed at the apparent time of extinction for each taxon
rangesCont <- sampleRanges(taxa, r = 0.5)
tree <- taxa2phylo(taxa, obs_time = rangesCont[,2])
phyloDiv(tree, drop.ZLB = FALSE)
#note that it drops taxa which were never sampled!

#testing with cryptic speciation
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, prop.cryptic = 0.5, nruns = 1,
nTotalTaxa = c(30,40), nExtant = 0, count.cryptic = TRUE)
taxaCrypt <- fossilRecord2fossilTaxa(record)
treeCrypt <- taxa2phylo(taxaCrypt)
layout(1)
plot(treeCrypt)
axisPhylo()

---

taxonSortPBDBocc  Sorting Unique Taxa of a Given Rank from Paleobiology Database Occurrence Data

**Description**

Functions for sorting out unique taxa from Paleobiology Database occurrence downloads, which should accept several different formats resulting from different versions of the PBDB API and different vocabularies available from the API.

**Usage**

```r
taxonSortPBDBocc(data, rank, onlyFormal = TRUE, cleanUncertain = TRUE,
cleanResoValues = c(NA, "\"", "", "n. sp.", "n. gen.", "", ""))
```

**Arguments**

- **data**: A table of occurrence data collected from the Paleobiology Database.
- **rank**: The selected taxon rank; must be one of ‘species’, ‘genus’, ‘family’, ‘order’, ‘class’ or ‘phylum’.
- **onlyFormal**: If TRUE (the default) only taxa formally accepted by the Paleobiology Database are returned. If FALSE, then the identified name fields are searched for any additional ‘informal’ taxa with the proper taxon. If their taxon name happens to match any formal taxa, their occurrences are merged onto the formal taxa. This argument generally has any appreciable effect when rank = species.
cleanUncertain  If TRUE (the default) any occurrences with an entry in the respective 'resolution' field that is *not* found in the argument cleanResoValue will be removed from the dataset. These are assumed to be values indicating taxonomic uncertainty, i.e. 'cf.' or '?'.

cleanResoValues  The set of values that can be found in a 'resolution' field that do not cause a taxon to be removed, as they do not seem to indicate taxonomic uncertainty.

Details

Data input for taxonSortPBDBocc are expected to be from version 1.2 API with the 'pbdb' vocabulary. However, datasets are passed to internal function translatePBDBocc, which attempts to correct any necessary field names and field contents used by taxonSortPBDBocc.

This function can pull either just the 'formally' identified and synonymized taxa in a given table of occurrence data or pull in addition occurrences listed under informal taxa of the sought taxonomic rank. Only formal taxa are sorted by default; this is controlled by argument onlyFormal. Pulling the informally-listed taxonomic occurrences is often necessary in some groups that have received little focused taxonomic effort, such that many species are linked to their generic taxon ID and never received a species-level taxonomic ID in the PBDB. Pulling both formal and informally listed taxonomic occurrences is a hierarchical process and performed in stages: formal taxa are identified first, informal taxa are identified from the occurrences that are 'leftover', and informal occurrences with name labels that match a previously sorted formally listed taxon are concatenated to the 'formal' occurrences for that same taxon, rather than being listed under separate elements of the list as if they were separate taxa. This function is simpler than similar functions that inspired it by using the input "rank" to both filter occurrences and directly reference a taxon's accepted taxonomic placement, rather than a series of specific if() checks.

Unlike some similar functions in other packages, such as version 0.3 paleobiodb’s pbdb_temp_range, taxonSortPBDBocc does not check if sorted taxa have a single 'taxon_no' ID number. This makes the blanket assumption that if a taxon's listed name in relevant fields is identical, the taxon is identical, with the important caveat that occurrences with accepted formal synonymies are sorted first based on their accepted names, followed by taxa without formal taxon IDs. This should avoid mistakingly linking the same occurrences to multiple taxa or assigning occurrences listed under separate formal taxa to the same taxon based on their 'identified' taxon name, as long as all formal taxa have unique names (which is an untested assumption). In some cases, this procedure is helpful, such as when taxa with identical generic and species names are listed under separate taxon ID numbers because of a difference in the listed subgenus for some occurrences (example, "Pseudoclimacograptus (Metaclimacograptus) hughesi‘ and ‘Pseudoclimacograptus hughesi’ in the PBDB as of 03/01/2015). Presumably any data that would be affected by differences in this procedure is very minor.

Occurrences with taxonomic uncertainty indicators in the listed identified taxon name are removed by default, as controlled by argument cleanUncertain. This is done by removing any occurrences that have an entry in primary_reso (was "genus_reso" in v1.1 API) when rank is a supraspecific level, and species_reso when rank = species, if that entry is not found in cleanResoValues. In some rare cases, when onlyFormal = FALSE, supraspecific taxonomic names may be returned in the output that have various 'cruft' attached, like 'n.sp.'.

Empty values in the input data table (""") are converted to NAs, as they may be due to issues with using read.csv to convert API-downloaded data.
Value

Returns a list where each element is different unique taxon obtained by the sorting function, and named with that taxon name. Each element is composed of a table containing all the same occurrence data fields as the input (potentially with some fields renamed and some field contents change, due to vocabulary translation).

Author(s)

David W. Bapst, but partly inspired by Matthew Clapham’s cleanTaxon (found at this location on github) and R package paleobioDB’s pbdb_temp_range function (found at this location on github).

See Also

coccData2timeList, plotOccData and the example graptolite dataset at grap2PDB

Examples

# load example graptolite PBDB occ dataset
data(graptPDB)

# get formal genera
occGenus <- taxonSortPBDBocc(graptOccPBDB, rank = "genus")
length(occGenus)

# get formal species
occSpeciesFormal <- taxonSortPBDBocc(graptOccPBDB, rank = "species")
length(occSpeciesFormal)

# yes, there are fewer 'formal' graptolite species in the PBDB then genera

# get formal and informal species
occSpeciesInformal <- taxonSortPBDBocc(graptOccPBDB, rank = "species", onlyFormal = FALSE)
length(occSpeciesInformal)

# way more graptolite species are 'informal' in the PBDB

# get formal and informal species
# including from occurrences with uncertain taxonomy
# basically everything and the kitchen sink
occSpeciesEverything <- taxonSortPBDBocc(graptOccPBDB, rank = "species", onlyFormal = FALSE, cleanUncertain = FALSE)
length(occSpeciesEverything)

## Not run:

# simple function for getting occurrence data from API v1.1
easyGetPBDBocc <- function(taxa, show = c("ident","phylo")){
  # cleans PBDB occurrence downloads of warnings
  taxa <- paste(taxa,collapse = ",")
  taxa <- paste(unlist(strsplit(taxa,"_")),collapse = "%20")
show <- paste(show,collapse = "",)
command <- paste0("http://paleobiodb.org/data1.2/occs/list.txt?base_name = ",
taxa,"&show = ",show,"&limit = all",
collapse = "")
command <- paste(unlist(strsplit(command,split = " ")),collapse = "\%20")
downData <- readLines(command)
if(length(grep("Warning",downData)) != 0){
  start <- grep("Records",downData)
  warn <- downData[1:(start-1)]
  warn <- sapply(warn, function(x)
    paste0(unlist(strsplit(unlist(strsplit(x,"\"")),"")),collapse = ""))
  warn <- paste0(warn,collapse = "\n")
  names(warn) <- NULL
  mat <- downData[-(1:start)]
  mat <- read.csv(textConnection(mat))
} else{
  mat <- downData
  mat <- read.csv(textConnection(mat))
}
return(mat)

#try a PBDB API download with lots of synonymization
#this should have only 1 species
#old way:
#acoData <- read.csv(paste0("http://paleobiodb.org/data1.1/occs/list.txt?base_name = Acosarina%20minuta&show = ident,phylo&limit = all"))
# with easyGetPBDBocc:
acoData <- easyGetPBDBocc("Acosarina minuta")
x <- taxonSortPBDBocc(acoData, rank = "species", onlyFormal = FALSE)
names(x)

#make sure works with API v1.1
dicelloData <- read.csv(paste0("http://paleobiodb.org","/data1.1/occs/list.txt?base_name = Dicellograptus","&show = ident,phylo&limit = all"))
dicelloOcc2 <- taxonSortPBDBocc(dicelloData, rank = "species", onlyFormal = FALSE)
names(dicelloOcc2)

#make sure works with compact vocab v1.1
dicelloData <- read.csv(paste0("http://paleobiodb.org","/data1.1/occs/list.txt?base_name = Dicellograptus","&show = ident,phylo&limit = all&vocab = com"))
dicelloOccCom1 <- taxonSortPBDBocc(dicelloData, rank = "species", onlyFormal = FALSE)
names(dicelloOccCom1)
head(dicelloOccCom1[[1]])[,1:7]

#make sure works with compact vocab v1.2
dicelloData <- read.csv(paste0("http://paleobiodb.org","/data1.2/occs/list.txt?base_name = Dicellograptus","&show = ident,phylo&limit = all&vocab = com"))
dicelloOccCom1 <- taxonSortPBDBocc(dicelloData, rank = "species", onlyFormal = FALSE)
Create a Taxonomy-Based Phylogeny ('Taxon Tree') from a Hierarchical Table of Taxonomy Memberships

This function takes a matrix of taxon names, indicating a set of hierarchical taxonomic relationships conveyed as nested placements for a set of tip-taxa (listed in the last column of the matrix) and returns a 'taxonomy-tree' phylogeny object of class 'phylo'.

Usage

```r
taxonTable2taxonTree(taxonTable, cleanTree = TRUE)
```

Arguments

taxonTable: A matrix of type character and multiple rows and columns, containing the tip taxa in the last column, one per row, with progressively larger taxa listed in prior columns (reading left-to-right). Invariant columns (i.e. taxa that all tip taxa are in) are allowed, but all but the most 'shallow' of such invariant taxa are dropped prior to transformation to a taxon-tree phylogeny object.

cleanTree: By default, the tree is run through a series of post-processing, including having singles collapsed, nodes reordered and being written out as a Newick string and read back in, to ensure functionality with ape functions and ape-derived functions. If FALSE, none of this post-processing is done and users should beware, as such trees can lead to hard-crashes of R.

Details

This function can deal with empty entries in cells of `taxonTable` by assuming these are lower-level taxa which are 'floating' freely somewhere in taxa several levels higher.

Value

A phylogeny of class 'phylo', where each tip is a taxon listed in the last column of the input taxonTable. Edges are scaled so that the distance from one taxon rank to another 1, then merged to remove singleton nodes. As not all taxa have parents at the immediate taxon level above, this leads to some odd cases. For example, two genera emanating from a node representing a class but with a very short (length = 1) branch and a long branch (length = 3) means one genus is simply placed in the class, with no family or order listed while the one on the long branch is within an order and family that is otherwise monogeneric.
The names of higher taxa than the tips should be appended as the element `node.label` for the internal nodes.

**Author(s)**

David W. Bapst

**See Also**

`makePBDBtaxonTree`, `parentChild2taxonTree`

**Examples**

```r
# let's create a small, really cheesy example
pokeTable <- rbind(cbind("Pokezooa","Shelloidea","Squirtadae"),
c("Squirtle","Blastoise","Wartortle"),
c("Pokezooa","Shelloidea","La"),
c("Pokezooa","La"),
cbind("Pokezooa","Hirsutamon","Rodentapokemorpha"),
c("Linoone","Sandshrew","Pikachu"),
c("Pokezooa","Hirsutamon",NA,"Ursaring")

pokeTree <- taxonTable2taxonTree(pokeTable)
plot(pokeTree); nodelabels(pokeTree$node.label)
```

**Description**

This function simulates the diversification of clades composed of monophyletic terminal taxa, which are distinguished in a fashion completely alternative to way taxa are defined in the simulation functions `simFossilRecord`, `taxa2cladogram` and `taxa2phylo`.

**Usage**

```r
simTermTaxa(ntaxa, sumRate = 0.2)

simTermTaxaAdvanced(p = 0.1, q = 0.1, mintaxa = 1, maxtaxa = 1000,
imtime = 1, maxtime = 1000, minExtant = 0, maxExtant = NULL,
min.cond = TRUE)

trueTermTaxaTree(TermTaxaRes, time.obs)

deadTree(ntaxa, sumRate = 0.2)
```
Arguments

- `ntaxa`: Number of monophyletic 'terminal' taxa (tip terminals) to be included on the simulated tree.
- `sumRate`: The sum of the instantaneous branching and extinction rates; see below.
- `p`: Instantaneous rate of speciation/branching.
- `q`: Instantaneous rate of extinction.
- `mintaxa`: Minimum number of total taxa over the entire history of a clade necessary for a dataset to be accepted.
- `maxtaxa`: Maximum number of total taxa over the entire history of a clade necessary for a dataset to be accepted.
- `mintime`: Minimum time units to run any given simulation before stopping.
- `maxtime`: Maximum time units to run any given simulation before stopping.
- `minExtant`: Minimum number of living taxa allowed at end of simulations.
- `maxExtant`: Maximum number of living taxa allowed at end of simulations.
- `min. cond`: If TRUE, the default, simulations are stopped when they meet all minimum conditions. If FALSE, simulations will continue until they hit maximum conditions, but are only accepted as long as they still meet all minimum conditions in addition.
- `TermTaxaRes`: The list output produced by `simTermTaxa`.
- `time.obs`: A per-taxon vector of times of observation for the taxa in `TermTaxaRes`.

Details

deadTree generates a time-scaled topology for an entirely extinct clade of a specific number of tip taxa. Because the clade is extinct and assumed to have gone extinct in the distant past, many details of typical birth-death simulators can be ignored. If a generated clade is already conditioned upon the (a) that some number of taxa was reached and (b) then the clade went extinct, the topology (i.e. the distribution of branching and extinction events) among the branches should be independent of the actual generating rate. The frequency of nodes is a simple mathematical function of the number of taxa (i.e. number of nodes is the number of taxa -1) and their placement should completely random, given that we generally treat birth-death processes as independent Poisson processes. Thus, in terms of generating the topology, this function is nothing but a simple wrapper for the ape function rtree, which randomly places splits among a set of taxa using a simple algorithm (see Paradis, 2012). To match the expectation of a birth-death process, new branch lengths are calculated as an exponential distribution with mean 1/sumRate, where sumRate represents the sum of the branching and extinction rates. Although as long as both the branching rate and extinction rates are more than zero, any non-ultrametric tree is possible, only when the two rates are non-zero and equal to each other will there be a high chance of getting an extinct clade with many tips. Any analyses one could do on a tree such as this will almost certainly give estimates of equal branching and extinction rates, just because all taxa are extinct.

simTermTaxa produces 'terminal-taxon' datasets; datasets of clades where the set of distinguishable taxa are defined as intrinsically monophyletic. (In version 1.6, I referred to this as the 'candle' mode, so named from the 'candling' horticultural practice and the visual conceptualization of the model.) On theoretical terms, terminal-taxon datasets are what would occur if (a) only descendant lineages
can be sample and (b) all taxa are immediately differentiated as of the last speciation event and continue to be so differentiated until they go extinct. In practice, this means the taxa on such a tree would represent a sample of all the terminal branches, which start with some speciation event and end in an extinction event. These are taken to be the true original ranges of these taxa. No further taxa can be sampled than this set, whatsoever. Note that the differentiation here is a result of a posteriori consideration of the phylogeny: one can’t even know what lineages could be sampled or the actual start points of such taxa until after the entire phylogeny of a group of organisms is generated.

Because all evolutionary history prior to any branching events is unsampled, this model is somewhat agnostic about the general model of differentiation among lineages. The only thing that can be said is that synapomorphies are assumed to be potentially present along every single branch, such that in an ideal scenario every clade could be defined. This would suggest very high anagenesis or bifurcation.

Because the set of observable taxa is a limited subset of the true evolution history, the true taxon ranges are not a faithful reproduction of the true diversity curve. See an example below.

simTermTaxa uses deadTree to make a phylogeny, so the only datasets produced are of extinct clades. simTermTaxaAdvanced is an alternative to simTermTaxa which uses simFossilRecord to generate the underlying pattern of evolutionary relationships and not deadTree. The arguments are thus similar to simFossilRecord, with some differences (as simTermTaxaAdvanced originally called the deprecated function simFossilTaxa). In particular, simTermTaxaAdvanced can be used to produce simulated datasets which have extant taxa.

trueTermTaxaTree is analogous to the function of taxa2phylo, in that it outputs the time-scaled-phylogeny for a terminal-taxon dataset for some times of observations. Unlike with the use of taxa2phylo on the output on simFossilRecord (via fossilRecord2fossilTaxa, there is no need to use trueTermTaxaTree to obtain the true phylogeny when times of extinction are the times of observation; just get the $tree element from the result output by simTermTaxa.

Also unlike with taxa2phylo, the cladistic topology of relationships among morphotaxa never changes as a function of time of observation. For obtaining the ‘ideal cladogram’ of relationships among the terminal taxa, merely take the $tree element of the output from simTermTaxaData and remove the branch lengths (see below for an example).

As with many functions in the paleotree library, absolute time is always decreasing, i.e. the present day is zero.

Value

deadTree gives time-scaled phylo object, with a $root.time element. As discussed above, the result is always an extinct phylogeny of exactly ntaxa.

simTermTaxa and simTermTaxaAdvanced both produce a list with two components: $taxonRanges which is a two-column matrix where each row gives the true first and last appearance of observable taxa and $tree which is a time-scaled phylogeny with end-points at the true last appearance time of taxa.

trueTermTaxaTree produces a time-scaled tree as a phylo object, which describes the relationships of populations at the times of observation given in the time.obs argument.

Author(s)

David W. Bapst
References


See Also

deadtree is simply a wrapper of the function rtree in ape.

For a very different way of simulating diversification in the fossil record, see simFossilRecord, fossilRecord2fossilTAXA, taxa2phylo and taxa2cladogram.

Examples

```r
set.seed(444)
# example for 20 taxa
termTaxaRes <- simTermTaxa(20)

# let look at the taxa...
taxa <- termTaxaRes$taxonRanges
taxicDivCont(taxa)
# because ancestors don't even exist as taxa
# the true diversity curve can go to zero
# kinda bizarre!

# the tree should give a better idea
tree <- termTaxaRes$tree
phyloDiv(tree)
# well, okay, its a tree.

# get the 'ideal cladogram' ala taxa2cladogram
  # much easier with terminal-taxa simulations as no paraphyletic taxa
cladogram <- tree
cladogram$edge.length <- NULL
plot(cladogram)

# trying out trueTermTaxaTree
# random times of observation: uniform distribution
time.obs <- apply(taxa,1,function(x) unif(1,x[2],x[1]))
tree1 <- trueTermTaxaTree(termTaxaRes,time.obs)
layout(1:2)
plot(tree)
plot(tree1)
layout(1)

# let's look at the change in the terminal branches
plot(tree$edge.length,tree1$edge.length)
# can see some edges are shorter on the new tree, cool

# let's now simulate sampling and use FADs
layout(1:2)
```
testEdgeMat is a small simple function which tests the $edge matrix of phylo objects for inconsistencies that can cause downstream analytical problems. The associated function, cleanNewPhylo puts an input phylo object, presumably freshly created or reconstituted by some function, through a series of post-processing. This includes having singles collapsed, nodes reordered and being written out as a Newick string and read back in, to ensure functionality with ape functions and ape-derived functions.
Usage

testEdgeMat(tree)

cleanNewPhylo(tree)

Arguments

tree A phylogeny object of type phylo

Details

Useful when doing complex manipulations and reconstitutions of phylo objects (or their de novo construction), and thus is used by a number of paleotree functions.

Value

For testEdgeMat, if all the checks in the function pass correctly, the logical TRUE is returned.
For cleanNewPhylo, an object of class phylo is returned.

Author(s)

David W. Bapst, with a large number of tests incorporated from Emmanuel Paradis’s checkValidPhylo function, provided at his github repository here, which was released GPL v>2:
https://github.com/emmanuelparadis/checkValidPhylo

Examples

set.seed(444)
tree <- rtree(10)
# should return TRUE
testEdgeMat(tree)

# should also work on star trees
testEdgeMat(stree(10))

# should also work on trees with two taxa
testEdgeMat(rtree(2))

# should also work on trees with one taxon
testEdgeMat(stree(1)))

#running cleanNewPhylo on this tree should have little effect
#beyond ladderizing it...
tree1 <- cleanNewPhylo(tree)

#compare outputs
layout(1:2)
plot(tree)
plot(tree1)
layout(1)
timeLadderTree
 Resolve Polytomies by Order of First Appearance

Description

Resolves polytomies in trees with lineages arranged in a pectinate pattern (i.e. a ladder-like subtree), ordered by the time of first appearance (FAD) for each lineage.

Usage

timeLadderTree(tree, timeData)

Arguments

tree          A phylo object

timeData     Two-column matrix of per-taxon first and last occurrences in absolute continuous time.

Details

This method of resolving polytomies assumes that the order of stratigraphic appearance perfectly depicts the order of branching. This may not be a good assumption for poorly sampled fossil records.

This function is for resolving trees when a continuous time-scale is known. For discrete time-scales, see the function bin_timePaleoPhy.

Taxa with the same identical first appearance date will be ordered randomly. Thus, the output is slightly stochastic, but only when ties exist. This is probably uncommon with real data on continuous time-scales.

Taxa not shared between the tree and the timeData matrix, or listed as having a FAD or LAD of NA in timeData will be dropped and will not be included in the output tree.

See this blog post for more information:
http://nemagruptus.blogspot.com/2012/07/resolving-polytomies-according-to.html

Value

Returns the modified tree as an object of class phylo, with no edge lengths.

Author(s)

David W. Bapst

See Also

di2multi
Examples

```r
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1, nTotalTaxa = c(100,200))
taxa <- fossilRecord2fossilTaxa(record)
tree <- taxa2cladogram(taxa)
ranges <- sampleRanges(taxa,r = 0.5)
tree1 <- timeLadderTree(tree,ranges)
layout(1:2)
plot(ladderize(tree),show.tip.label = FALSE)
plot(ladderize(tree1),show.tip.label = FALSE)

# an example with applying timeLadderTree to discrete time data
rangeData <- binTimeData(ranges,int.len = 5)  # sim discrete range data
tree2 <- bin_timePaleoPhy(tree,timeList = rangeData,timeres = TRUE)
plot(ladderize(tree),show.tip.label = FALSE)
plot(ladderize(tree2),show.tip.label = FALSE)
axisPhylo()

layout(1)
```

Description

Functions for manipulating data where the first and last appearances of taxa are known from bounded intervals of time. The two main functions listed here are for converting between (1) a data structure consisting of a single ‘flat’ table where each taxon is listed as a set of four dates (a fourDate data type), and (2) a list format where each taxon is listed as its first and last intervals, with an associated table of age bounds for the intervals referred to in the first table (referred to as a timeList data structure by many paleoTree functions).

Usage

```r
timeList2fourDate(timeList)

fourDate2timeList(fourDate)
```

Arguments

- **timeList**: A list composed of two matrices with two columns each, the first giving interval start and end date bounds, and the second giving taxon first and last interval appearances in reference to the intervals listed in the first matrix.
A four column matrix where each row is a different taxon, the first two columns are the lower and upper bounds on the time of first appearance for that taxon and the third and fourth columns are respectively the lower and upper bounds on the time of last appearance for that taxon, all in time before present.

Details

timelist2fourdate is for converting from a timelist format to a fourdate format. fourdate2timelist is for converting from a fourdate format to a timelist format.

Value

A converted data object, respective to the function applied.

Author(s)

David W. Bapst

References

See my recent blog post on temporal datasets in paleontology for some details:

See Also

bin_timePaleoPhy and taxicDivDisc for common applications; binTimeData for a simulation function for such data objects

Examples

# timelist object from the retiolinae dataset
data(retiolitinae)
str(retioRanges)
taxicDivDisc(retioRanges)
fourDateRet <- timelist2fourDate(retioRanges)

# total uncertainty in retio first and last appearances?
sum(((fourDateRet[,1]-fourDateRet[,2])+)(fourDateRet[,3]-fourDateRet[,4]))

# convert back
newTimelist <- fourDate2timelist(fourDateRet)
taxicDivDisc(retioRanges)
Typical 'a posteriori' Time-Scaling Approaches For Paleontological Phylogenies

Description

Time-scales an unscaled cladogram of fossil taxa using information on their temporal ranges, using various methods. Also can resolve polytomies randomly and output samples of randomly-resolved trees. As simple methods of time-scaling phylogenies of fossil taxa can have biasing effects on macroevolutionary analyses (Bapst, 2014, Paleobiology), this function is largely retained for legacy purposes and plotting applications. The time-scaling methods implemented by the functions listed here do not return realistic estimates of divergence dates, users should investigate other time-scaling methods such as cal3TimePaleoPhy.

Usage

timePaleoPhy(tree, timeData, type = "basic", vartime = NULL, ntrees = 1, randres = FALSE, timeres = FALSE, add.term = FALSE, inc.term.adj = FALSE, dateTreatment = "firstLast", node.mins = NULL, noisyDrop = TRUE, plot = FALSE)

bin_timePaleoPhy(tree, timeList, type = "basic", vartime = NULL, ntrees = 1, nonstoch.bin = FALSE, randres = FALSE, timeres = FALSE, sites = NULL, point.occurs = FALSE, add.term = FALSE, inc.term.adj = FALSE, dateTreatment = "firstLast", node.mins = NULL, noisyDrop = TRUE, plot = FALSE)

Arguments

tree
An unscaled cladogram of fossil taxa, of class phylo. Tip labels must match the taxon labels in the respective temporal data.

timeData
Two-column matrix of first and last occurrences in absolute continuous time, with row names as the taxon IDs used on the tree. This means the first column is very precise FADs (first appearance dates) and the second column is very precise LADs (last appearance dates), reflect the precise points in time when taxa first and last appear. If there is stratigraphic uncertainty in when taxa appear in the fossil record, it is preferable to use the 'bin' time-scaling functions; however, see the argument dateTreatment.

type
Type of time-scaling method used. Can be "basic", "equal", "equal_paleotree_legacy", "equal_date.phylo_legacy" "aba", "zbla" or "mbl". Type = "basic" by default. See details below.

vartime
Time variable; usage depends on the method ‘type’ argument. Ignored if type = "basic".

ntrees
Number of time-scaled trees to output. If ntrees is greater than one and both randres is false and dateTreatment is neither ‘minMax’ or ‘randObs’, the function will fail and a warning is issued, as these arguments would simply produce multiple identical time-scaled trees.
randres Should polytomies be randomly resolved? By default, timePaleoPhy does not resolve polytomies, instead outputting a time-scaled tree that is only as resolved as the input tree. If randres = T, then polytomies will be randomly resolved using multi2di from the package ape. If randres = T and ntrees = 1, a warning is printed that users should analyze multiple randomly-resolved trees, rather than a single such tree, although a tree is still output.

timeres Should polytomies be resolved relative to the order of appearance of lineages? By default, timePaleoPhy does not resolve polytomies, instead outputting a time-scaled tree that is only as resolved as the input tree. If timeres = TRUE, then polytomies will be resolved with respect to time using the paleotree function timeladdertree. See that functions help page for more information; the result of time-order resolving of polytomies generally does not differ across multiple uses, unlike use of multi2di.

add.terms If true, adds terminal ranges. By default, this function will not add the ranges of taxa when time-scaling a tree, so that the tips correspond temporally to the first appearance datums of the given taxa. If add.terms = T, then the ‘terminal ranges’ of the taxa are added to the tips after tree is time-scaled, such that the tips now correspond to the last appearance datums.

inc.terms.adj If true, includes terminal ranges in branch length estimates for the various adjustment of branch lengths under all methods except ‘basic’ (i.e., a terminal length branch will not be treated as zero length is this argument is TRUE if the taxon at this tip has a non-zero duration). By default, this argument is FALSE and this function will not include the ranges of taxa when adjusting branch lengths, so that zero-length branches before first appearance times will be extended. An error is returned if this argument is true but type = "basic" or add.terms = FALSE, as this argument is inconsistent with those argument options.

dateTreatment This argument controls the interpretation of timeData. The default setting ‘first-Last’ treats the dates in timeData as a column of precise first and last appearances, such that first appearances will be used to date nodes and last appearances will only be called on if add.terms = TRUE. A second option, added by great demand, is ‘minMax’ which treats these dates as minimum and maximum bounds on single point dates. Under this option, all taxa in the analysis will be treated as being point dates, such that the first appearance is also the last. These dates will be pulled under a uniform distribution. If ‘minMax’ is used, add.terms becomes meaningless, and the use of it will return an error message. A third option is ‘randObs’. This assumes that the dates in the matrix are first and last appearance times, but that the desired time of observation is unknown. Thus, this is much like ‘firstLast’ except the effective time of observation (the taxon’s LAD under ‘firstLast’) is treated an uncertain date, and is randomly sampled between the first and last appearance times. The FAD still is treated as a fixed number, used for dating the nodes. In previous versions of paleotree, this was called in timePaleoPhy using the argument rand.obs, which has been removed for clarity. This temporal uncertainty in times of observation might be useful if a user is interested in applying phylogeny-based approaches to studying trait evolution, but have per-taxon measurements of traits that come from museum specimens with uncertain temporal placement. With both arguments ‘minMax’
and 'randObs', the sampling of dates from random distributions should compel users to produce many time-scaled trees for any given analytical purpose. Note that 'minMax' returns an error in 'bin' time-scaling functions; please use 'points.occur' instead.

**node.mins**

The minimum dates of internal nodes (clades) on a phylogeny can be set using node.mins. This argument takes a vector of the same length as the number of nodes, with dates given in the same order as nodes are ordered in the `tree$edge` matrix. Note that in `tree$edge`, terminal tips are given the first set of numbers (1:ntip(tree)), so the first element of node.mins is the first internal node (the node numbered `ntip(tree)`+1, which is generally the root for most `phylo` objects read by `read.tree`). Not all nodes need be given minimum dates; those without minimum dates can be given as NA in node.mins, but the vector must be the same length as the number of internal nodes in `tree`. These are minimum date constraints, such that a node will be forced to be at least as old as this date, but the final date may be even older depending on the taxon dates used, the time-scaling method applied, the `vartime` used and any other minimum node dates given (e.g. if a clade is given a very old minimum date, this will (of course) over-ride any minimum dates given for clades that that node is nested within). Although `vartime` does adjust the node age downwards when the equal method is used, if a user has a specific date they’d like to constrain the root to, they should use node.mins instead because the result is more predictable.

**noisyDrop**

If TRUE (the default), any taxa dropped from `tree` due to not having a matching entry in the time data will be listed in a system message.

**plot**

If TRUE, plots the input and output phylogenies.

**timelist**

A list composed of two matrices giving interval times and taxon appearance dates. The rownames of the second matrix should be the taxon IDs, identical to the `tip.labels` for `tree`. See details.

**nonstoch.bin**

If TRUE, dates are not stochastically pulled from uniform distributions. See below for more details.

**sites**

Optional two column matrix, composed of site IDs for taxon FADs and LADs. The sites argument allows users to constrain the placement of dates by restricting multiple fossil taxa whose FADs or LADs are from the same very temporally restricted sites (such as fossil-rich Lagerstatten) to always have the same date, across many iterations of time-scaled trees. To do this, simply give a matrix where the "site" of each FAD and LAD for every taxon is listed, as corresponding to the second matrix in timelist. If no sites matrix is given (the default), then it is assumed all fossil come from different "sites" and there is no shared temporal structure among the events.

**point.occur**

If true, will automatically produce a 'sites' matrix which forces all FADs and LADs to equal each other. This should be used when all taxa are only known from single 'point occurrences', i.e. each is only recovered from a single bed/horizon, such as a Lagerstatten.

**Details**

*Time-Scaling Methods*
These functions are an attempt to unify and collect previously used and discussed 'a posteriori' methods for time-scaling phylogenies of fossil taxa. Unfortunately, it can be difficult to attribute some time-scaling methods to specific references in the literature.

There are five main a posteriori approaches that can be used by timePaleoPhy. Four of these main types use some value of absolute time, chosen a priori, to time-scale the tree. This is handled by the argument var.time, which is NULL by default and unused for type "basic".

"basic" This most simple of time-scaling methods ignores var.time and scales nodes so they are as old as the first appearance of their oldest descendant (Smith, 1994). This method produces many zero-length branches (Hunt and Carrano, 2010).

"equal" The 'equal' method defined by G. Lloyd and used in Brusatte et al. (2008) and Lloyd et al. (2012). Originally usable in code supplied by G. Lloyd, the algorithm is recreated here as closely as possible. This method works by increasing the time of the root divergence by some amount and then adjusting zero-length branches so that time on early branches is re-apportioned out along those later branches equally. Branches are adjusted in order relative to the number of nodes separating the edge from the root, going from the furthest (most shallow) edges to the deepest edges. The choice of ordering algorithm can have an unanticipated large effect on the resulting time-scaled trees created using "equal" and it appears that paleotree and functions written by G. Lloyd were not always consistent. The default option described here was only introduced into either software sources in August 2014. Thus, two legacy 'equal' methods are included in this function, so users can emulate older ordering algorithms for 'equal' which are now deprecated, as they do not match the underlying logic of the original 'equal' algorithm and do not minimize down-passes when adjusting branch lengths on the time-scaled tree.

The root age can be adjusted backwards in time by either increasing by an arbitrary amount (via the var.time argument) or by setting the root age directly (via the node.mins argument); conversely, the function will also allow a user to opt to not alter the root age at all.

"equal_paleotree_legacy" Exactly like 'equal' above, except that edges are ordered instead by their depth (i.e. number of nodes from the root). This minor modified version was referred to as 'equal' for this timePaleoPhy function in paleotree until February 2014, and thus is included here solely for legacy purposes. This ordering algorithm does not minimize branch adjustment cycles, like the newer default offered under currently 'equal'.

"equal_date.phylo_legacy" Exactly like 'equal' above, except that edges are ordered relative to their time (i.e. total edge length) from the root following the application of the 'basic' time-scaling method, exactly as in G. Lloyd’s original application. This was the method for sorting edges in the "equal" algorithm in G. Lloyd’s date.phylo script and DatePhylo in package strap until August 2014, and was the default "equal" algorithm in paleotree’s timePaleoPhy function from February 2014 until August 2014. This ordering algorithm does not minimize branch adjustment cycles, like the newer default offered under currently 'equal'. Due to how the presence of zero-length branches can make ordering branches based on time to be very unpredictable, this version of the 'equal' algorithm is highly not recommended.

"aba" All branches additive. This method takes the "basic" tree and adds var.time to all branches. Note that this time-scaling method can warp the tree structure, leading to tips to originate out of order with the appearance data used.

"zlb" Zero-length branches additive. This method adds var.time to all zero-length branches in the "basic" tree. Discussed (possibly?) by Hunt and Carrano, 2010. Note that this time-scaling
method can warp the tree structure, leading to tips to originate out of order with the appearance data used.

"mbl" Minimum branch length. Scales all branches so they are greater than or equal to vartime, and subtract time added to later branches from earlier branches in order to maintain the temporal structure of events. A version of this was first introduced by Laurin (2004).

These functions cannot time-scale branches relative to reconstructed character changes along branches, as used by Lloyd et al. (2012). Please see DatePhylo in R package strap for this functionality.

These functions will intuitively drop taxa from the tree with NA for range or are missing from timeData or timeList. Taxa dropped from the tree will be will be listed in a message output to the user. The same is done for taxa in the timeList object not listed in the tree.

As with many functions in the paleotree library, absolute time is always decreasing, i.e. the present day is zero.

As of August 2014, please note that the branch-ordering algorithm used in 'equal' has changed to match the current algorithm used by DatePhylo in package strap, and that two legacy versions of 'equal' have been added to this function, respectively representing how timePaleoPhy and DatePhylo (and its predecessor dateNphylo) applied the 'equal' time-scaling method.

Interpretation of Taxon Ages in timePaleoPhy

`timePaleoPhy` is primarily designed for direct application to datasets where taxon first and last appearances are precisely known in continuous time, with no stratigraphic uncertainty. This is an uncommon form of data to have from the fossil record, although not an impossible form (micropaleontologists often have very precise range charts, for example). Instead, most data has some form of stratigraphic uncertainty. However, for some groups, the more typical 'first' and 'last' dates found in the literature or in databases represent the minimum and maximum absolute ages for the fossil collections that a taxon is known is known from. Presumably, the first and last appearances of that taxon in the fossil record is at unknown dates within these bounds.

As of paleotree version 2.0. the treatment of taxon ages in timePaleoPhy is handled by the argument datetreatment. By default, this argument is set to 'firstLast' which means the matrix of ages are treated as precise first and last appearance dates (i.e. FADs and LADs). The earlier FADs will be used to calibrate the node ages, which could produce fairly nonsensical results if these are 'minimum' ages instead and reflect age uncertainty. Alternatively, datetreatment can be set to 'minMax' which instead treats taxon age data as minimum and maximum bounds on a single point date. These point dates, if the minimum and maximum bounds option is selected, are chose under a uniform distribution. Many time-scaled trees should be created to approximate the uncertainty in the dates. Additionally, there is a third option for datetreatment: users may also make it so that the 'times of observation' of trees are uncertain, such that the tips of the tree (with terminal ranges added) should be randomly selected from a uniform distribution. Essentially, this third option treats the dates as first and last appearances, but treats the first appearance dates as known and fixed, but the 'last appearance' dates as unknown. In previous versions of paleotree, this third option was enacted with the argument rand.obs, which has been removed for clarity.

Interpretation of Taxon Ages in bin_timePaleoPhy

As an alternative to using timePaleoPhy, bin_timePaleoPhy is a wrapper of timePaleoPhy which produces time-scaled trees for datasets which only have interval data available. For each output tree, taxon first and last appearance dates are placed within their listed intervals under a uniform distribution. Thus, a large sample of time-scaled trees will approximate the uncertainty in the actual timing of the FADs and LADs. In some ways, treating taxonomic age uncertainty may be more
logical via bin_timePaleoPhy, as it is tied to specific interval bounds, and there are more options available for certain types of age uncertainty, such as for cases where specimens come from the same fossil site.

The input timeList object for bin_timePaleoPhy can have overlapping (i.e. non-sequential) intervals, and intervals of uneven size. Taxa alive in the modern should be listed as last occurring in a time interval that begins at time 0 and ends at time 0. If taxa occur only in single collections (i.e. their first and last appearance in the fossil record is synchronous, the argument point_occur will force all taxa to have instantaneous durations in the fossil record. Otherwise, by default, taxa are assumed to first and last appear in the fossil record at different points in time, with some positive duration. The sites matrix can be used to force only a portion of taxa to have simultaneous first and last appearances.

By setting the argument nonstoch.bin to TRUE for bin_timePaleoPhy, the dates are NOT stochastically pulled from uniform bins but instead FADs are assigned to the earliest time of whichever interval they were placed in and LADs are placed at the most recent time in their placed interval. This option may be useful for plotting. The sites argument becomes arbitrary if nonstoch.bin = TRUE.

If timeData or the elements of timeList are actually data.frames (as output by read.csv or read.table), these will be coerced to a matrix.

**Tutorial**

A tutorial for applying the time-scaling functions in paleotree, along with an example using real (graptolite) data, can be found here: http://nemagraptus.blogspot.com/2013/06/a-tutorial-to-cal3-time-scaling-using.html

**Value**

The output of these functions is a time-scaled tree or set of time-scaled trees, of either class phylo or multiphylo, depending on the argument ntrees. All trees are output with an element $root.time. This is the time of the root on the tree and is important for comparing patterns across trees. Note that the $root.time element is defined relative to the earliest first appearance date, and thus later tips may seem to occur in the distant future under the 'aba' and 'zbla' time-scaling methods.

Trees created with bin_timePaleoPhy will output with some additional elements, in particular $ranges.used, a matrix which records the continuous-time ranges generated for time-scaling each tree. (Essentially a pseudo-timeData matrix.)

**Note**

Please account for stratigraphic uncertainty in your analysis. Unless you have exceptionally resolved data, select an appropriate option in dateTreatmet within timePaleoPhy, use the more sophisticated bin_timePaleoPhy or code your own wrapper function of timePaleoPhy that accounts for stratigraphic uncertainty in your dataset.

**Author(s)**

David W. Bapst, heavily inspired by code supplied by Graeme Lloyd and Gene Hunt.

**References**


See Also

cal3TimePaleoPhy, binTimeData, multi2di

For an alternative time-scaling function, which includes the ‘ruta’ method that weights the time-scaling of branches by estimates of character change along with implementations of the ‘basic’ and ‘equal’ methods described here, please see function DatePhylo in package strap.

Examples

```r
# examples with empirical data

data(retiolitinae)

#Can plot the unscaled cladogram
plot(retioTree)

#Can plot discrete time interval diversity curve with retioRanges
taxicDivDisc(retioRanges)

#Use basic time-scaling (terminal branches only go to FADs)
tree <- bin_timePaloPhy(tree = retioTree, timeList = retioRanges, type = "basic", ntrees = 1, plot = TRUE)

#Use basic time-scaling (terminal branches go to LADs)
tree <- bin_timePaloPhy(tree = retioTree, timeList = retioRanges, type = "basic", add.term = TRUE, ntrees = 1, plot = TRUE)

#minimum branch length time-scaling (terminal branches only go to FADs)
tree <- bin_timePaloPhy(tree = retioTree, timeList = retioRanges, type = "mbl", vartime = 1, ntrees = 1, plot = TRUE)
```

----------
# examples with simulated data

#Simulate some fossil ranges with simFossilRecord
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1,
nTotalTaxa = c(30,40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
#simulate a fossil record with imperfect sampling with sampleRanges
rangesCont <- sampleRanges(taxa, r = 0.5)
#let's use taxa2cladogram to get the 'ideal' cladogram of the taxa
cladogram <- taxa2cladogram(taxa,plot = TRUE)
#Now let's try timePaleoPhy using the continuous range data
tree <- timePaleoPhy(cladogram, rangesCont, type = "basic", plot = TRUE)
#plot diversity curve
phylodiv(tree)

#that tree lacked the terminal parts of ranges (tips stops at the taxon FADs)
#let's add those terminal ranges back on with addTerm

#can randomly resolve trees using the argument randres
#each resulting tree will have polytomies randomly resolved in different ways using multi2di
tree <- timePaleoPhy(cladogram, rangesCont, type = "basic", ntrees = 1, randres = TRUE,
addTerm = TRUE, plot = TRUE)
#notice well the warning it prints!
#we would need to set ntrees to a large number to get a fair sample of trees

#if we set ntrees>1, timePaleoPhy will make multiple time-trees
trees <- timePaleoPhy(cladogram, rangesCont, type = "basic", ntrees = 9, randres = TRUE,
addTerm = TRUE, plot = TRUE)
#let's compare nine of them at once in a plot
layout(matrix(1:9,3,3))
parOrig <- par(no.readonly = TRUE)
par(mar = c(1,1,1,1))
for(i in 1:9){plot(ladderize(trees[[i]]), show.tip.label = FALSE, no.margin = TRUE)}
#they are all a bit different!

#we can also resolve the polytomies in the tree according to time of first appearance
#via the function timeLadderTree, by setting the argument 'timeres' to TRUE
tree <- timePaleoPhy(cladogram, rangesCont, type = "basic", ntrees = 1, timeres = TRUE,
addTerm = TRUE, plot = TRUE)

#can plot the median diversity curve with multiDiv
layout(1);par(parOrig)
multiDiv(trees)

#compare different methods of timePaleoPhy
layout(matrix(1:6,3,2))
parOrig <- par(no.readonly = TRUE)
par(mar = c(3,2,1,2))
plot(ladderize(timePaleoPhy(cladogram, rangesCont, type = "basic", vartime = NULL, add.term = TRUE)))
axisPhylo(); text(x = 50, y = 23, "type = basic", adj = c(0,0.5), cex = 1.2)
plot(ladderize(timePaleoPhy(cladogram, rangesCont, type = "equal", vartime = 10, add.term = TRUE)))
axisPhylo(); text(x = 55, y = 23, "type = equal", adj = c(0,0.5), cex = 1.2)
plot(ladderize(timePaleoPhy(cladogram, rangesCont, type = "aba", vartime = 1, add.term = TRUE)))
axisPhylo(); text(x = 55, y = 23, "type = aba", adj = c(0,0.5), cex = 1.2)
plot(ladderize(timePaleoPhy(cladogram, rangesCont, type = "zlba", vartime = 1, add.term = TRUE)))
axisPhylo(); text(x = 55, y = 23, "type = zlba", adj = c(0,0.5), cex = 1.2)
plot(ladderize(timePaleoPhy(cladogram, rangesCont, type = "mb1", vartime = 1, add.term = TRUE)))
axisPhylo(); text(x = 55, y = 23, "type = mb1", adj = c(0,0.5), cex = 1.2)
layout(1); par(parOrig)

# using nodemins
# let's say we have (molecular??) evidence that node #5 is at least 1200 time-units ago
# to use nodemins, first need to drop any unshared taxa
droppers <- cladogram$tip.label[is.na(cladogram$tip.label)]
cladodrop <- drop.tip(cladogram, droppers)
# now make vector same length as number of nodes
nodeDates <- rep(NA, Nnode(cladodrop))
nodeDates[5] <- 1200
ttree1 <- timePaleoPhy(cladodrop, rangesCont, type = "basic",
   randres = FALSE, node.mins = nodeDates, plot = TRUE)
ttree2 <- timePaleoPhy(cladodrop, rangesCont, type = "basic",
   randres = TRUE, node.mins = nodeDates, plot = TRUE)

# Using bin_timePaleoPhy to time-scale with discrete interval data
# first let's use binTimeData() to bin in intervals of 1 time unit
rangesDisc <- binTimeData(rangesCont, int.length = 1)
ttreeB1 <- bin_timePaleoPhy(cladogram, rangesDisc, type = "basic", ntrees = 1, randres = TRUE, add.term = TRUE, plot = FALSE)
# notice the warning it prints!
phylodiv(ttreeB1)

# with time-order resolving via timeLadderTree
ttreeB2 <- bin_timePaleoPhy(cladogram, rangesDisc, type = "basic", ntrees = 1, timeres = TRUE, add.term = TRUE, plot = FALSE)

# can also force the appearance timings not to be chosen stochastically
ntreeB3 <- bin_timePaleoPhy(cladogram, rangesDisc, type = "basic", ntrees = 1, nonstoch.bin = TRUE, randres = TRUE, add.term = TRUE, plot = FALSE)

# testing node.mins in bin_timePaleoPhy
ntree <- bin_timePaleoPhy(cladodrop, rangesDisc, type = "basic", ntrees = 1, add.term = TRUE, randres = FALSE, node.mins = nodeDates, plot = TRUE)
# with randres = TRUE
ntree <- bin_timePaleoPhy(cladodrop, rangesDisc, type = "basic", ntrees = 1, add.term = TRUE, randres = TRUE, node.mins = nodeDates, plot = TRUE)

# simple three taxon example for testing inc.term.adj
ranges1 <- cbind(c(3,4,5), c(2,3,1)); rownames(ranges1) <- paste("t",1:3,sep = "")
clad1 <- read.tree(file = NA, text = "(t1,(t2,t3));")
**timeSliceTree**

*Time-Slicing a Phylogeny*

**Description**

Removes the portion of a tree after a set point in time, as if the tree after that moment had been sliced away.

**Usage**

```r
timeSliceTree(ttree, sliceTime, drop.extinct = FALSE, plot = TRUE)
```

**Arguments**

- **ttree**: A time-scaled phylogeny of class phylo
- **sliceTime**: Time to 'slice' the tree at. See details below.
- **drop.extinct**: If true, drops tips that go extinct before timeSlice.
- **plot**: If true, plots input and output trees for comparison.

**Details**

The function assumes that ttree will generally have an element called $root.time, which is the time before present that the root divergence occurred. If $root.time is not present as an element of ttree, then it is assumed the tip furthest from the root is at time 0 (present-day) and a new $root.time is calculated (a warning will be issued in this case).

The sliceTime is always calculated as on the same scale as ttree$root.time. In other words, if root.time = 100, then timeSlice = 80 will slice the tree 20 time units after the root.

If drop.extinct = T, then extinct tips are dropped and (if present) the $root.time of ttree is adjusted. This is done using the function dropExtinct.

**Value**

Returns the modified phylogeny as an object of class phylo.

Tip labels for cut branches which held multiple descendant tips will be the label for the earliest appearing tip descendant of that branch. This is somewhat arbitrary; the actual morphotaxon present at that time might have been a different taxon. For simulated datasets, use taxa2phylo.
treeContradiction

**Measure the Contradiction Difference Between Two Phylogenetic Topologies**

**Description**

An alternative measure of pair-wise dissimilarity between two tree topologies which ignores differences in phylogenetic resolution between the two, unlike typical metrics (such as Robinson-Foulds distance). The metric essentially counts up the number of splits on both trees that are directly contradicted by a split on the contrasting topology (treating both as unrooted). By default, this 'contradiction difference' value is then scaled to between 0 and 1, by dividing by the total number of splits that could have been contradicted across both trees (\( \frac{2 \times (\text{Number of shared tips} - 2)}{\text{Total possible splits}} \)).

**Author(s)**

David W. Bapst, with modification of code by Klaus Schliep to avoid use of function `distNodes` which has difficulty with large trees and greatly benefiting function runtime.

**See Also**

`phylodiv, dropExtinct, dropExtant`

Also see the function `treeSlice` in the library phytools, which will slice a tree at some point in and return all the subtrees which remain after the slicing time. (Effectively the reverse of `timeSliceTree`.)

**Examples**

```r
# a neat example of using phylodiv with timeSliceTree
# to simulate doing extant-only phylogeny studies
# of diversification... in the past!
set.seed(444)
record <- simFossilRecord(p = 0.1, q = 0.1, nruns = 1, nTotalTaxa = c(30, 40), nExtant = 0)
taxa <- fossilRecord2FossilTaxa(record)
taxicDivCont(taxa)
# that's the whole diversity curve

# with timeSliceTree we could look at the lineage accumulation curve
# we'd get of species sampled at a point in time
tree <- tax2phylo(taxa)
# use timeSliceTree to make tree of relationships up until time = 950
tree950 <- timeSliceTree(tree, sliceTime = 950, plot = TRUE, drop.extinct = FALSE)

# use drop.extinct = T to only get the tree of lineages extant at time = 950

# now it's an ultrametric tree with many fewer tips...
# lets plot the lineage accumulation plot on a log scale
phylodiv(tree950, plotLogRich = TRUE)
```
On this scaled, 0 represents no conflicting relationships and 1 reflects two entirely conflicting topologies, similar to the rescaling in Colless’s consensus fork index.

Usage

\[
\text{treeContradiction(tree1, tree2, rescale = TRUE)}
\]

Arguments

- **tree1, tree2**: Two phylogenies, with the same number of tips and an identical set of tip labels, both of class `phylo`.
- **rescale**: A logical. If `FALSE`, the raw number of contradicted splits across both trees is reported. If `TRUE` (the default), the contradiction difference value is returned rescaled to the total number of splits across both input trees that could have contradicted.

Details

Algorithmically, conflicting splits are identified by counting the number of splits (via `ape`'s `prop.part`) on one tree that disagree with at least one split on the other tree: for example, split (AB)CD would be contradicted by split (AC)BD. To put it another way, all we need to test for is whether the taxa segregated by that split were found to be more closely related to some other taxa, not so segregated by the considered split.

This metric was designed mainly for use with trees that differ in their resolution, particularly when it is necessary to compare between summary trees (such as consensus trees of half-compatibility summaries) from separate phylogenetic analyses. Note that comparing summary trees can be problematic in some instances, and users should carefully consider their question of interest, and whether it may be more ideal to consider whole samples of trees (e.g., the posterior sample, or the sample of most parsimonious trees).

The contradiction difference is *not* a metric distance: most notably, the triangle inequality is not held and thus the 'space' it describes between topologies is not a metric space. This can be shown most simply when considering any two different but fully-resolve topologies and a third topology that is a star tree. The star tree will have a zero pair-wise CD with either fully-resolved phylogeny, but there will be a positive CD between the fully-resolved trees. An example of this is shown in the examples below.

The CD also suggest very large differences when small numbers of taxa shift greatly across the tree, a property shared by many other typical tree comparisons, such as RF distances. See examples below.

Value

The contradiction difference between two trees is reported as a single numeric variable.

Author(s)

David W. Bapst. This code was produced as part of a project funded by National Science Foundation grant EAR-1147537 to S. J. Carlson.
References

This contradiction difference measure was introduced in:

See Also

See phangorn’s function for calculating the Robinson-Foulds distance: treecontradiction.
Graeme Lloyd’s metatree package, currently not on CRAN, also contains the function MultiTreeDistance for calculating both the contradiction difference measure and the Robinson-Foulds distance. This function is optimized for very large samples of trees or very large trees, and thus may be faster than treecontradiction. Also see the function MultiTreeContradiction in the same package.

Examples

# let’s simulate two trees
set.seed(1)
treeA <- rtree(30, br = NULL)
treeB <- rtree(30, br = NULL)

# visualise the difference between these two trees
library(phytools)
plot(cophyllo(treeA, treeB))

# what is the Robinson-Foulds (RF) distance between these trees?
library(phangorn)
treedist(treeA, treeB)

# The RF distance is less intuitive when
# we consider a tree that isn’t well-resolved

# let’s simulate the worst resolved tree possible: a star tree
treeC <- stree(30)

# plot the tanglegram between A and C
plot(cophyllo(treeA, treeC))

# however the RF distance is *not* zero
# even though the only difference is a difference in resolution
treedist(treeA, treeC)

# End(Not run)
twoWayEcologyCluster

R-Mode vs Q-Mode Two-Way Cluster Analyses and Abundance Plot for Community Ecology Data

Description

This mode plots both R-mode (across sites) and Q-mode (across taxa) dendrograms for a community ecology dataset, with branches aligned with a grid of dots representing the relative abundance of taxa at each site in the dataset.

Usage

```r
twoWayEcologyCluster(xDist, yDist, propAbund, clustMethod = "average", marginBetween = 0.1, abundExpansion = 4, xAxisLabel = "Across Sites", yAxisLabel = "Across Taxa")
```

Arguments

**xDist**
The pair-wise distance matrix for the cluster diagram drawn along the horizontal axis of the graphic. Should be a distance matrix, or a matrix that can be coerced to a distance matrix, for the same number of units as rows in *propAbund*.

**yDist**
The pair-wise distance matrix for the cluster diagram drawn along the vertical axis of the graphic. Should be a distance matrix, or a matrix that can be coerced to a distance matrix, for the same number of units as columns in *propAbund*.

**propAbund**
A matrix of abundance data, preferably relative abundance scaled as proportions of the total number of individuals at each site. This data determines the size scale of the taxon/site dots.

**clustMethod**
The agglomerative clustering method used, as with argument *method* with function *hclust*. *clustMethod* must be one of "average" (the default method for this function, also known as average-linkage or as UPGMA), "ward.D", "ward.D2", "single", "complete", "mcquitty" (also known as WPGMA), "median" (also known as WPGMC) or "centroid" (also known as UPGMC).

**marginBetween**
Argument controlling space placed between the cluster diagrams and the abundance plot. Default is 0.1.

**abundExpansion**
An argument that is a multiplier controlling the size of dots plotted for reflecting relative abundance.

**xAxisLabel**
The label placed on the horizontal axis of the plot.

**yAxisLabel**
The label placed on the vertical axis of the plot.

Details

You might be able to apply this to datasets that aren’t community ecology datasets of proportional abundance, but I can’t guarantee pr even predict what will happen.

Value

This function creates a plot, and returns nothing, not even invisible output.

Author(s)

David W. Bapst

References

The function here was designed to emulate previous published 'two-way' cluster diagrams, particularly the one in Miller, 1988:


See Also

Several other functions for community ecology data in paleotree are described at the *communityEcology* help file. Also see the example dataset, *kanto*.
### Examples

```r
## Not run:
library(vegan)

# generate random community ecology data
# using a Poisson distribution
data <- matrix(rpois(5*7,1),5,7)

# relative abundance, distance matrices
propAbundMat <- t(apply(data,1,function(x) x/sum(x)))
rownames(propAbundMat) <- paste0("x", 1:nrow(propAbundMat))
colnames(propAbundMat) <- paste0("y", 1:ncol(propAbundMat))

siteDist <- vegdist(propAbundMat, "bray")
taxaDist <- vegdist(t(propAbundMat), "bray")

dev.new(width=10)
twoWayEcologyCluster(
  xDist = siteDist,
  yDist = taxaDist,
  propAbund = propAbundMat
)

```

```
```

# now let's try an example with some pokemon data
data(kanto)

# get distance matrices for sites and taxa
# based on bray-curtis dist
# standardized to total abundance

# standardize site matrix to relative abundance
siteStand <- decostand(kanto, method = "total")

# calculate site distance matrix (Bray-Curtis)
siteDist <- vegdist(siteStand, "bray")

# calculate taxa distance matrix (Bray-Curtis)
# from transposed standardized site matrix
taxaDist <- vegdist(t(siteStand), "bray")

dev.new(width=10)
twoWayEcologyCluster(
  xDist = siteDist,
  yDist = taxaDist,
  propAbund = siteStand
)

## End(Not run)
```
unitLengthTree **Scale Tree to Unit-Length**

Description

Rescales all edges of a phylogeny to be equal to 1 ("unit-length").

Usage

```
unitLengthTree(tree)
```

Arguments

- `tree`: an object of class phylo

Details

Probably not a good way to scale a tree for comparative studies. What does it mean to scale every edge of the phylogeny to the same length?

Value

Returns the modified phylogeny as an object of class phylo. Any $root.time element is removed.

See Also

As an alternative to using unitLengthTree in comparative studies, see `timePaleoPhy`

See also `speciationTree` in the package geiger, which does essentially the same thing as unitLengthTree.

Examples

```
set.seed(444)
tree <- rtree(10)
layout(1:2)
plot(tree)
plot(unitLengthTree(tree))
layout(1)
```
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