

# Package ‘paleoTS’

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**Title** Analyze Paleontological Time-Series

**Version** 0.6.2

**Description** Facilitates analysis of paleontological sequences of trait values. Functions are provided to fit, using maximum likelihood, simple evolutionary models (including unbiased random walks, directional evolution, stasis, Ornstein-Uhlenbeck, covariate-tracking) and complex models (punctuation, mode shifts).

**Depends** R (>= 3.4.0)

**License** GPL-3

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**NeedsCompilation** no

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---

`akaike.wts`*Compute Akaike weights from AIC scores*

---

**Description**

Compute Akaike weights from AIC scores

**Usage**

```
akaike.wts(aa)
```

**Arguments**

`aa` vector of AIC or AICc scores

**Details**

This function converts a vector of AIC or AICc scores to Akaike weights, a vector that summarizes proportional model support.

**Value**

a vector of Akaike weights

**See Also**

[compareModels](#)

**Examples**

```
akaike.wts(c(10, 14, 20))
```

---

`as.paleoTS`*Make a Paleontological Time-series object*

---

**Description**

Combines information into an object of class paleoTS

**Usage**

```
as.paleoTS(
  mm,
  vv,
  nn,
  tt,
  MM = NULL,
  genpars = NULL,
  label = NULL,
  start.age = NULL,
  oldest = c("first", "last"),
  reset.time = TRUE
)
```

**Arguments**

mm	vector of sample means
vv	vector of sample variances
nn	vector of sample sizes
tt	vector of sample ages
MM	vector of true means (simulated data)
genpars	generating parameters (simulated data)
label	optional, label for time-series
start.age	optional, age of oldest sample
oldest	value indicating if the oldest sample is first or last in the sequence
reset.time	logical; if TRUE, then change time scale to start at t=0 and adjust start.age accordingly

**Details**

This function combines data into a paleoTS object. For empirical data it may be more convenient to use read.paleoTS.

If sample ages decrease through the sequence, as if given in millions of years ago, tt will automatically be converted to time elapsed from the beginning of the sequence as long as reset.time = TRUE.

**Value**

a paleoTS object

**Note**

All model-fitting functions estimate the contribution of sampling noise to the observed differences between samples. They do this assuming that the trait is represented by sample means, which have sampling variances equal to variance divided by sample size,  $vv/nn$ . If one is interested in analyzing

statistics other than the sample mean (medians, quantiles, or other statistics), use the the following procedure: set the statistic in question as the mm values, replace vv with a vector of the squared standard errors for each estimate (generated by other means, for example bootstrapping), and set all values of nn to one. When fitting such time-series, be sure to set the argument pool = FALSE.

### See Also

[read.paleoTS](#)

### Examples

```
x <- as.paleoTS(mm = rnorm(20), vv = rep(1, 20), nn = rep(25, 20), tt=0:19)
plot(x) # easier to use sim.Stasis()
```

---

as.paleoTSfit	<i>Create a paleoTSfit object</i>
---------------	-----------------------------------

---

### Description

Create a paleoTSfit object

### Usage

```
as.paleoTSfit(
  logL,
  parameters,
  modelName,
  method,
  K,
  n,
  se,
  convergence = NULL,
  logLFunction = NULL,
  ...
)
```

### Arguments

logL	model log-likelihood
parameters	model parameter estimates
modelName	model name
method	parameterization, either "AD" or "Joint"
K	number of model parameters
n	sample size
se	standard errors of parameter estimates
convergence	code indicating optimization convergence
logLFunction	name of the log-likelihood function
...	optional additional elements added by some functions

**Value**

a paleoTSfit object

**Note**

All model-fitting functions use this function to package the resulting data-model fits. Users will not need to call this function.

**Examples**

```
# fake example; users won't need to use this unless they make their own model-fitting functions
w <- as.paleoTSfit(logL = 10, parameters = 2, modelName = "StrictStasis",
                  method = "Joint", K = 1, n = 25, se = NULL)
```

---

bootSimpleComplex	<i>Bootstrap test to see if a complex model is significantly better than a simple model.</i>
-------------------	--

---

**Description**

Bootstrap test to see if a complex model is significantly better than a simple model.

**Usage**

```
bootSimpleComplex(
  y,
  simpleFit,
  complexFit,
  nboot = 99,
  minb = 7,
  ret.full.distribution = FALSE,
  parallel = FALSE,
  ...
)
```

**Arguments**

y	a paleoTS object
simpleFit	a paleoTSfit object, representing the model fit of a simple model
complexFit	a paleoTSfit object, representing the model fit of a complex model
nboot	number of replications for parametric bootstrapping
minb	minimum number of populations within each segment
ret.full.distribution	logical, indicating if the null distribution for the likelihood ratio from the parametric bootstrap should be returned
parallel	logical, if TRUE, the bootstrapping is done using parallel computing
...	further arguments, passed to optimization functions

## Details

Simulations suggest that AICc can be overly liberal with complex models with mode shifts or punctuations (Hunt et al., 2015). This function implements an alternative of parametric bootstrapping to compare the fit of a simple model with a complex model. It proceeds in five steps:

1. Compute the observed gain in support from the simple to complex model as the likelihood ratio,  $LR_{obs} = -2(\log L_{simple} - \log L_{complex})$
2. Simulate trait evolution under the specified simple model nboot times
3. Fit to each simulated sequence the specified simple and complex models
4. Measure the gain in support from simple to complex as the bootstrap likelihood ratio for each simulated sequence
5. Compute the P-value as the percentile of the bootstrap distribution corresponding to the observed LR.

Argument `simpleFit` should be a `paleoTS` object returned by the function `fitSimple` or similar functions (e.g., `opt.joint.GRW`, `opt.GRW`, etc.). Argument `complexFit` must be a `paleoTS` object returned by `fitGpunc` or `fitModeShift`.

Calculations can be speeded up by setting `parallel = TRUE`, which uses package `doParallel` to run the bootstrap replicates in parallel, using one fewer than the number of detected cores.

## Value

A list of the observed likelihood ratio statistic, `LRob`s, the P-value of the test, and the number of bootstrap replicates. If `ret.full.distribution = TRUE`, the null distribution of likelihood ratios generated by parametric bootstrapping is also returned.

## References

Hunt, G., M. J. Hopkins and S. Lidgard. 2015. Simple versus complex models of trait evolution and stasis as a response to environmental change. *PNAS* 112(16): 4885-4890.

## See Also

[sim.Stasis.RW](#), [fitModeShift](#)

## Examples

```
## Not run:
x <- sim.Stasis.RW(ns = c(15, 15), omega = 0.5, ms = 1, order = "Stasis-RW")
ws <- fitSimple(x)
wc <- fitModeShift(x, order = "Stasis-RW", rw.model = "GRW")
bootSimpleComplex(x, ws, wc, nboot = 50, minb = 7) # nboot too low for real analysis!

## End(Not run)
```

---

cantius_L	<i>Time-series of the length of lower first molar for the Cantius lineage</i>
-----------	---

---

**Description**

Time-series of the length of lower first molar for the Cantius lineage

**Usage**

```
cantius_L
```

**Format**

a paleoTS object with the data

**Source**

Clyde, W. C. and P. D. Gingerich (1994). Rates of evolution in the dentition of early Eocene Cantius: comparison of size and shape. *Paleobiology* 20(4): 506-522.

---

checkSSMresiduals	<i>Compute and (optionally) plot residuals from SSM model fit</i>
-------------------	---

---

**Description**

Compute and (optionally) plot residuals from SSM model fit

**Usage**

```
checkSSMresiduals(
  y,
  w,
  show.plot = TRUE,
  resid.type = c("standardized", "unstandardized")
)
```

**Arguments**

y	a paleoTS object
w	a paleoTSfit object, where the fitting functions uses SSM
show.plot	logical, if TRUE, several diagnostic plots are shown
resid.type	residual type, either "standardized" or "unstandardized"

**Details**

It is recommended that `resid.type` be set to the default, "standardized", which will scale residuals by their expected standard deviation

**Value**

a vector of residuals, returned invisibly

**Examples**

```
y <- sim.GRW(ns = 50, ms = 0.2)
w <- fitSimple(y, model = "URW", method = "SSM") # wrong model
checkSSMresiduals(y, w, show.plot = TRUE) # positive residuals show model mis-fit
```

---

compareModels

*Compare model fits for a paleontological time-series*

---

**Description**

Takes output from model-fitting functions and compiles model-fit information (log-likelihood, AICc, etc.) into a convenient table

**Usage**

```
compareModels(..., silent = FALSE, sort = FALSE)
```

**Arguments**

<code>...</code>	any number of model fit (as <code>paleoTSfit</code> ) objects
<code>silent</code>	if TRUE, suppresses printing
<code>sort</code>	if TRUE, the table sorts models from best to worst

**Value**

if `silent = FALSE`, the table is printed and nothing is returned. If `silent = TRUE`, printing is suppressed and a list of two objects is returned: the table of model fits, `modelFits`, and a list of parameter estimates, `p1`.

**Examples**

```
x <- sim.GRW(ns = 40, ms = 0.5, vs = 0.1)
m1 <- fitSimple(x, model = "GRW") # the true model
m2 <- fitSimple(x, model = "URW")
plot(x, modelFit = m1)
compareModels(m1, m2)
```

---

dorsal.spines	<i>Time-series of dorsal spine data from a fossil stickleback lineage</i>
---------------	---

---

**Description**

Time-series of dorsal spine data from a fossil stickleback lineage

**Usage**

dorsal.spines

**Format**

a paleoTS object of the mean number of dorsal spines (log-transformed)

**Source**

Bell, M.A., M.P. Travis and D.M. Blouw 2006. Inferring natural selection in a fossil threespine stickleback. *Paleobiology* **32**:562-577.

Hunt, G., M. A. Bell and M. P. Travis (2008). Evolution toward a new adaptive optimum: phenotypic evolution in a fossil stickleback lineage. *Evolution* 62(3): 700-710.

---

ESD	<i>Compute Expected Squared Divergence (ESD) for Evolutionary Models</i>
-----	--

---

**Description**

Computes for a specified model and duration of time the expected squared divergence (ESD), which is a useful measure of the magnitude or rate of change across different models.

**Usage**

```
ESD(
  y,
  dt,
  model = c("GRW", "URW", "Stasis", "allThree"),
  method = c("Joint", "AD"),
  pool = TRUE,
  ...
)
```

**Arguments**

y	a paleoTS object
dt	the time interval to evaluate ESD
model	the model of evolution to assume; see Details
method	Joint or AD parameterization
pool	logical, if TRUE, variances are averaged (pooled) across samples
...	other arguments to the model-fitting functions

**Details**

Hunt (2012) argued that rate metrics make sense only in the context of specific models of evolution. It is thus difficult to meaningfully compare rates across sequences generated by different evolutionary processes. ESD values can be used for a specified model and duration as a comparable measure of the amount of evolutionary change that is expected. Acceptable values for the model argument can be "GRW" for the general random walk (directional change), "URW" for the unbiased random walk, and "Stasis." In addition, one can also specify "allThree", in which case all these models will be fit and the resulting ESD will be the weighted average of them, using model support (Akaike weights) for the weighting (see Hunt [2012], p. 370)

**Value**

the ESD value

**References**

Hunt, G. 2012. Measuring rates of phenotypic evolution and the inseparability of tempo and mode. *Paleobiology* 38:351–373.

**Examples**

```
x<- sim.GRW(ns=20)
esd.urw<- ESD(x, dt=10, model="URW")
esd.all<- ESD(x, dt=10, model="allThree")
```

---

fit.sgs

---

*Fit a model of trait evolution with a protracted punctuation.*


---

**Description**

This function fits a model of punctuated change that is protracted enough that it is captured by multiple transitional populations. Trait evolution starts in stasis, shifts to a general random walk, and then shifts back into stasis.

**Usage**

```
fit.sgs(
  y,
  minb = 7,
  oshare = TRUE,
  pool = TRUE,
  silent = FALSE,
  hess = FALSE,
  meth = "L-BFGS-B",
  model = "GRW"
)
```

**Arguments**

<code>y</code>	a paleoTS object
<code>minb</code>	minimum number of populations within each segment
<code>oshare</code>	logical, if TRUE, variance assumed to be shared (equal) across segments
<code>pool</code>	if TRUE, sample variances are substituted with their pooled estimate
<code>silent</code>	logical, if TRUE, progress updates are suppressed
<code>hess</code>	if TRUE, standard errors computed from the Hessian matrix are returned
<code>meth</code>	optimization method, passes to <code>optim</code>
<code>model</code>	type of random walk: "URW", unbiased random walk, or "GRW", a general (directional) random walk

**Value**

a paleoTSfit object

**See Also**

[fitGpunc](#)

**Examples**

```
x <- sim.sgs(ns = c(10, 10, 10)) # default values OK
w <- fit.sgs(x, minb = 8) # increase minb so example takes less time; not recommended!
plot(x)
abline(v = c(16, 31), lwd = 3) # actual shifts
abline(v = c(w$parameters[6:7]), lwd = 2, lty = 3, col = "red") # inferred shifts
```

---

`fit3models`*Fit a set of standard evolutionary models*

---

**Description**

Fit a set of standard evolutionary models

**Usage**

```
fit3models(y, silent = FALSE, method = c("Joint", "AD", "SSM"), ...)
```

```
fit4models(y, silent = FALSE, method = c("Joint", "AD", "SSM"), ...)
```

**Arguments**

<code>y</code>	a paleoTS object
<code>silent</code>	if TRUE, results are returned as a list and not printed
<code>method</code>	"Joint", "AD", or "SSM"; see <a href="#">fitSimple</a>
<code>...</code>	other arguments passed to model fitting functions

**Details**

Function `fit3models` fits the general (biased) random walk (GRW), unbiased random walk (URW), and Stasis models. In addition to these three, `fit4models` also fits the model of Strict Stasis.

**Value**

if `silent = FALSE`, a table of model fit statistics, also printed to the screen. if `silent = TRUE`, a list of the model fit statistics and model parameter values.

**Functions**

- `fit4models()`: add model of "Strict Stasis" to the three models

**See Also**

[fitSimple](#)

**Examples**

```
x <- sim.GRW(ns = 50, ms = 0.2)
fit4models(x)
```

---

`fit9models`*Fit large set of models to a time-series*

---

### Description

This function fits nine models to a time-series following Hunt et al. (2015). These include the simple models fit by `fit4models` along with mode shift and punctuation models.

### Usage

```
fit9models(y, silent = FALSE, method = c("Joint", "AD"), ...)
```

### Arguments

<code>y</code>	a paleoTS object
<code>silent</code>	logical, if TRUE, progress updates are suppressed
<code>method</code>	parameterization to use: Joint or AD; see Details
<code>...</code>	other arguments, passed to optimization functions

### Value

if `silent = FALSE`, a table of model fit statistics, also printed to the screen. if `silent = TRUE`, a list of the model fit statistics and model parameter values.

### References

Hunt, G., M. J. Hopkins and S. Lidgard. 2015. Simple versus complex models of trait evolution and stasis as a response to environmental change. PNAS 112(16): 4885-4890.

### Examples

```
## Not run:  
x <- sim.Stasis.RW(ns = c(15, 15), omega = 0.5, ms = 1, order = "Stasis-RW")  
plot(x)  
fit9models(x)  
  
## End(Not run)
```

fitGpunc

*Fit trait evolution model with punctuations estimated from the data***Description**

Fit trait evolution model with punctuations estimated from the data

**Usage**

```
fitGpunc(
  y,
  ng = 2,
  minb = 7,
  pool = TRUE,
  oshare = TRUE,
  method = c("Joint", "AD"),
  silent = FALSE,
  hess = FALSE,
  parallel = FALSE,
  ...
)
```

**Arguments**

y	a paleoTS object
ng	number of groups (segments) in the sequence
minb	minimum number of populations within each segment
pool	if TRUE, sample variances are substituted with their pooled estimate
oshare	logical, if TRUE, variance assumed to be shared (equal) across segments
method	parameterization to use: Joint or AD; see Details
silent	logical, if TRUE, progress updates are suppressed
hess	if TRUE, standard errors computed from the Hessian matrix are returned
parallel	logical, if TRUE, the analysis is done in parallel
...	other arguments, passed to optimization functions

**Details**

This function tests all possible shift points for punctuations, subject to the constraint that the number of populations in each segment is always  $\geq \text{minb}$ . The shiftpoint yielding the highest log-likelihood is returned as the solution, along with the log-likelihoods (`all.logl`) of all tested shift points (GG).

The function uses `opt.punc` (if `method = "AD"`) or `opt.joint.punc` (if `method = "Joint"`) to do the fitting.

**Value**

a paleoTSfit object with the results of the model-fitting.

**Note**

Calculations can be speeded up by setting `parallel = TRUE`, which uses package `doParallel` to run the bootstrap replicates in parallel, using one fewer than the number of detected cores.

**See Also**

[fit9models](#), [sim.punc](#)

**Examples**

```
x <- sim.punc(ns = c(15, 15), theta = c(0,3), omega = c(0.1, 0.1))
w.punc <- fitGpunc(x, oshare = TRUE)
plot(x, modelFit = w.punc)
```

---

fitModeShift

*Fit model in which the mode of trait evolution shifts once*

---

**Description**

Trait evolution is modeled as a shift from a random walk (general or unbiased) to stasis, or vice versa.

**Usage**

```
fitModeShift(
  y,
  minb = 7,
  pool = TRUE,
  order = c("Stasis-RW", "RW-Stasis"),
  rw.model = c("URW", "GRW"),
  method = c("Joint", "AD"),
  silent = FALSE,
  hess = FALSE,
  ...
)
```

**Arguments**

<code>y</code>	paleoTS object
<code>minb</code>	minimum number of populations within each segment
<code>pool</code>	if TRUE, sample variances are substituted with their pooled estimate
<code>order</code>	whether stasis or random walk come first, one of Stasis-RW or RW-Stasis

rw.model	whether the random walk segment is an unbiased random walk, URW or a general random walk, GRW
method	parameterization to use: Joint or AD
silent	logical, if TRUE, progress updates are suppressed
hess	if TRUE, standard errors computed from the Hessian matrix are returned
...	other arguments, passed to optimization functions

**Value**

a paleoTSfit object

**See Also**

[sim.Stasis.RW](#)

**Examples**

```
x <- sim.Stasis.RW(ns = c(15, 15), omega = 0.5, ms = 1, order = "Stasis-RW")
plot(x)
w <- fitModeShift(x, order = "Stasis-RW", rw.model = "GRW")
abline(v = x$tt[15], lwd = 3) # actual shift point
abline(v = x$tt[w$par["shift1"]], lty = 3, lwd = 2, col = "red") # inferred shift
```

---

fitMult

*Fit the same simple model across multiple time-series*


---

**Description**

Fit the same simple model across multiple time-series

**Usage**

```
fitMult(
  y1,
  model = c("GRW", "URW", "Stasis", "covTrack"),
  method = c("Joint", "AD"),
  pool = TRUE,
  z1 = NULL,
  hess = FALSE
)
```

**Arguments**

<code>y1</code>	a list of paleoTS objects
<code>model</code>	the model to fit; see Details
<code>method</code>	parameterization to use: Joint or AD
<code>pool</code>	if TRUE, sample variances are substituted with their pooled estimate
<code>z1</code>	for the covTrack model only, a list of covariate vectors, one each paleoTS object in <code>y1</code>
<code>hess</code>	if TRUE, standard errors computed from the Hessian matrix are returned

**Details**

This function fits a model with shared parameters across multiple trait time-series. The most likely application would be to model a common evolutionary dynamic across different sequences, perhaps representing time-series of the same trait and lineage from different localities or time intervals.

Four simple models are currently implemented:

- **GRW**: parameters `mstep` and `vstep` of the general random walk are shared across sequences.
- **URW**: parameter `vstep` of the unbiased random walk is shared across sequences.
- **Stasis**: parameter `omega` of stasis is shared across sequences.
- **covTrack**: parameters `b0`, `b1`, and `evvar` of the covariate-tracking model are shared across sequences.

Under the joint parameterization, `method = "Joint"`, an additional parameter, `anc` is fit, representing the ancestral (starting) trait value. This parameter is estimated separately in each sequence so it is not assumed that they all start at the same trait value.

**Value**

a `paleoTSfit` object with the results of the model-fitting

**Note**

The models are described in the help for `fitSimple` and the functions linked from there.

**See Also**

[fitSimple](#)

**Examples**

```
x1 <- sim.GRW(ms = 1, vs = 0.2)
x2 <- sim.GRW(ms = 1, vs = 0.2)
fitMult(list(x1, x2), model = "GRW")
```

fitSimple

*Fit simple models of trait evolution***Description**

Fit simple models of trait evolution

**Usage**

```
fitSimple(
  y,
  model = c("GRW", "URW", "Stasis", "StrictStasis", "OU", "ACDC", "covTrack"),
  method = c("Joint", "AD", "SSM"),
  pool = TRUE,
  z = NULL,
  hess = FALSE
)
```

**Arguments**

y	a paleoTS object
model	the model to be fit, one of "GRW", "URW", "Stasis", "OU", "ACDC", "covTrack"
method	parameterization to use: Joint, AD or SSM; see Details
pool	if TRUE, sample variances are substituted with their pooled estimate
z	a vector of a covariate, used only for the "covTrack" model
hess	if TRUE, standard errors computed from the Hessian matrix are returned

**Details**

This is a convenience function that calls the specific individual functions for each model and parameterization, such as `opt.GRW` and `opt.joint.GRW`. The models that this function can fit are:

- **GRW**: General Random Walk. Under this model, evolutionary changes, or "steps" are drawn from a distribution with a mean of `mstep` and variance of `vstep`. `mstep` determines directionality and `vstep` determines volatility (Hunt, 2006).
- **URW**: Unbiased Random Walk. Same as GRW with `mstep = 0`, and thus evolution is non-directional. For a URW, `vstep` is the rate parameter.
- **Stasis**: This parameterization follows Sheets & Mitchell (2001), with a constant mean `theta` and variance `omega` (equivalent to white noise).
- **Strict Stasis**: Same as Stasis with `omega = 0`, indicating no real evolutionary differences; all observed variation is sampling error (Hunt et al. 2015).
- **OU**: Ornstein-Uhlenbeck model (Hunt et al. 2008). This model is that of a population ascending a nearby peak in the adaptive landscape. The optimal trait value is `theta`, `alpha` indicates the strength of attraction to that peak (= strength of stabilizing selection around `theta`), `vstep` measures the random walk component (from genetic drift) and `anc` is the trait value at the start of the sequence.

- **ACDC**: Accelerating or decelerating evolution model (Blomberg et al. 2003). This model is that of a population undergoing a random walk with a step variance that increases or decreases over time. The initial step variance is  $v_{step0}$ , and the parameter  $r$  controls its rate of increase (if positive) or decrease (if negative) over time. When  $r < 0$ , the is equivalent to the "Early burst" model of Harmon et al.
- **covTrack**: Covariate-tracking (Hunt et al. 2010). The trait tracks a covariate with slope  $b_1$ , consistent with an adaptive response.  $evar$  is the residual variance, and, under method = "Joint",  $b_0$  is the intercept of the relationship between trait and covariate. model.

### Value

a `paleoTSfit` object with the model fitting results

### Note

For the covariate-tracking model,  $z$  should be a vector of length  $n$  when method = "Joint" and  $n - 1$  when method = "AD", where  $n$  is the number of samples in  $y$ .

Method = "Joint" is a full likelihood approach, considering each time-series as a joint sample from a multivariate normal distribution. Method = "AD" is a REML approach that uses the differences between successive samples. They perform similarly, but the Joint approach does better under some circumstances (Hunt, 2008).

### References

- Hunt, G. 2006. Fitting and comparing models of phyletic evolution: random walks and beyond. *Paleobiology* 32(4): 578-601.
- Hunt, G. 2008. Evolutionary patterns within fossil lineages: model-based assessment of modes, rates, punctuations and process. p. 117-131 *In* From Evolution to Geobiology: Research Questions Driving Paleontology at the Start of a New Century. Bambach, R. and P. Kelley (Eds).
- Hunt, G., M. A. Bell and M. P. Travis. 2008. Evolution toward a new adaptive optimum: phenotypic evolution in a fossil stickleback lineage. *Evolution* 62(3): 700-710.
- Sheets, H. D., and C. Mitchell. 2010. Why the null matters: statistical tests, random walks and evolution. *Genetica* 112– 113:105–125.
- Blomberg, S. P., T. Garland, and A. R. Ives. 2003. Testing for phylogenetic signal in comparative data: behavioural traits are more labile. *Evolution* 57(4):717-745.
- Harmon, L. J. et al. 2010. Early bursts of body size and shape evolution are rare in comparative data. *Evolution* 64(8):2385-2396.

### See Also

[opt.GRW](#), [opt.joint.GRW](#), [opt.joint.OU](#), [opt.covTrack](#)

### Examples

```
y <- sim.Stasis(ns = 20, omega = 2)
w1 <- fitSimple(y, model = "GRW")
w2 <- fitSimple(y, model = "URW")
w3 <- fitSimple(y, model = "Stasis")
```

```
compareModels(w1, w2, w3)
```

---

IC *Compute Information Criteria*

---

**Description**

Compute Information Criteria

**Usage**

```
IC(logL, K, n = NULL, method = c("AICc", "AIC", "BIC"))
```

**Arguments**

logL	log-likelihood
K	number of parameters
n	sample size
method	either "AIC", "AICc", or "BIC"

**Value**

the value of the specified information criterion

**Note**

This function is used internally by the model-fitting functions. It will not generally be called directly by the user.

**Examples**

```
ic1 <- IC(logL = 0, K = 2, method = "AIC") # plain AIC
ic2 <- IC(logL = 0, K = 2, n = 10, method = "AICc")
ic3 <- IC(logL = 0, K = 2, n = 1000, method = "AICc") # converges to AIC with increasing n
print(rbind(ic1, ic2, ic3))
```

---

Kfiltertv

*Time-varying Kalman filter calculations*


---

### Description

Time-varying Kalman filter calculations

### Usage

```
Kfiltertv(num, y, Atv, mu0, Sigma0, Phitv, Ups, Gam, Qtv, Rtv, input)
```

### Arguments

num	the number of samples in the time-series
y	values of the time-series
Atv	q x p x n observation array
mu0	p x 1 vector setting the mean of the system at time zero
Sigma0	p x p variance matrix of the system at time zero
Phitv	p x p x n array reflecting autoregression of the state variables
Ups	p x r matrix with the coefficients/parameters relating the inputs to the system equation
Gam	q x r matrix with the coefficients/parameters relating the inputs to the observation equation
Qtv	p x p x n array of system stochastic variance; user needs to ensure positive definite
Rtv	q x q x n array observation stochastic variance; user needs to ensure positive definite
input	n x r array of the exogenous variables/covariates

### Details

For the dimensions of the argument arrays, n is the length of the time-series, q is the dimension of the observation variable(s), p is the dimension of the state variable(s), and r is the dimension of the input variable(s).

This function is based on the `Kfilter` function of the *astsa* package, modified modified to allow for time-varying terms for the Kalman filter. This modification facilitates fitting a broader array of models and handling non-uniform temporal spacing of samples. See the documentation for that function, and the reference below for additional information.

**Value**

A list of the following elements:

- xp one-step-ahead prediction of the state
- Pp mean square prediction error
- xf filter value of the state
- Pf mean square filter error
- like log-likelihood
- innov innovation series
- sig innovation covariances
- Kn last value of the gain, needed for smoothing

**Note**

This function is used in the internal SSM log-likelihood functions for the models. The user will not need to use this they create their own model-fitting functions.

**Author(s)**

John Fricks (jfricks@asu.edu)

**References**

Shumway, R. H., and D. S. Stoffer. 2017. Time Series Analysis and its Applications (4th Ed.) Springer International.

**Examples**

```
y <- sim.GRW(ms = 0, vs = 1, vp = 0)
n <- length(y)
kf <- Kfiltertv(n ,y = y$mm, Atv = array(1, dim = c(1,1,n)), mu0 = y$mm[1],
               Sigma0 = y$vv[1]/y$nn[1], Phitv = array(1, dim = c(1,1,n)),
               Ups = NULL, Gam = NULL, Qtv = array(1, dim = c(1,1,n)),
               Rtv = array(0, dim = c(1,1,n)), input = NULL)
```

---

In.paleoTS

*Approximate log-transformation of time-series data*

---

**Description**

Approximate log-transformation of time-series data

**Usage**

```
In.paleoTS(y)
```

**Arguments**

`y` a paleoTS object

**Details**

For a random variable  $x$ , its approximate mean on a natural log scale is the log of its untransformed mean. The approximate variance on a log scale is equal to the squared coefficient of variation.

**Value**

the converted paleoTS object

**Note**

This transformation only makes sense for variables with dimension and a true zero point, such as lengths and areas.

**References**

Hunt, G. 2006. Fitting and comparing models of phyletic evolution: random walks and beyond. *Paleobiology* 32:578-601.  
Lewontin, R. 1966. On the measurement of relative variability. *Systematic Zoology* 15:141-142.

**Examples**

```
x <- sim.Stasis(ns = 10, theta = 20, omega = 1)
plot(x)
x1 <- ln.paleoTS(x)
plot(x1)
```

---

LRI

*Log-rate, Log-interval (LRI) method of Gingerich*


---

**Description**

Gingerich (1993) introduced a method that plots on log-log scale, the rate and interval for each pair of samples in an evolutionary sequence. On this plot, the slope is interpreted as an indicator of evolutionary mode (-1 for stasis, 0.5 for random walk, 0 for directional), and the intercept is interpreted as a measure of the rate of evolution over one generation.

**Usage**

```
LRI(y, gen.per.t = 1e+06, draw = TRUE)
```

**Arguments**

`y` a paleoTS object  
`gen.per.t` the number of generations per unit time  
`draw` logical, if TRUE, a plot is produced

**Details**

Following Gingerich (1993), a robust line is fit through the points by minimizing the sum of absolute deviations. If generations are one year long and time is measured in Myr,  $\text{gen.per.t} = 1e6$ .

**Value**

A named vector with three elements: Intercept, slope, and GenerationalRate

**Note**

This method was important in early attempts to disentangle evolutionary tempo and mode. Likelihood-based methods have a more sound statistical basis, and in particular the estimation of 'Generational Rates' using LRI is compromised by sampling error; see Hunt (2012) and the example below.

**References**

Gingerich, P.D. 1993. Quantification and comparison of evolutionary rates. *American Journal of Science* 293-A:453–478.

Hunt, G. 2012. Measuring rates of phenotypic evolution and the inseparability of tempo and mode. *Paleobiology* 38:351–373.

**See Also**

[lynchD](#)

**Examples**

```
set.seed(1)
xFast <- sim.GRW(ns = 20, ms = 0.5, vs = 0.2) # fast evolution
xSlow <- sim.Stasis(ns = 20, omega = 0)      # strict stasis (zero rates)
lri.Fast <- LRI(xFast, draw = FALSE)
lri.Slow <- LRI(xSlow, draw = FALSE)
print(lri.Fast[3], 4)
print(lri.Slow[3], 4) # LRI thinks strict stasis rates are MUCH faster!
```

---

lynchD

---

*Compute Lynch's Delta rate metric*


---

**Description**

This function computes D, the rate metric proposed by Lynch (1990). This metric derives from the random walk model, with  $D = V_{\text{step}} / (2V_p)$ , where  $V_{\text{step}}$  is the step variance of the unbiased random walk, and  $V_p$  is the within sample variance, pooled among samples. Under mutation - drift equilibrium, D is expected to range approximately between  $5e-5$  and  $5e-3$ .

**Usage**

```
lynchD(y, gen.per.t = 1e+06, pool = TRUE, method = c("Joint", "AD"), ...)
```

**Arguments**

y	a paleoTS object
gen.per.t	the number of generations per unit time
pool	logical, whether variances should be pooled over samples
method	parameterization to use: based on ancestor-descendant (AD) differences, or Joint consideration of all samples
...	further arguments, passed to opt.URW or opt.joint.URW

**Value**

D	value of rate metric
pooled.var	value of pooled within-sample variance
gen.per.t	number of generations per unit time
vstep	computed Vstep, at the original time scale of y
drift.range	expected minimum and maximum values of D consistent with neutral evolution
result	conclusion reached about the plausibility of neutral evolution

**References**

Lynch (1990). The rate of morphological evolution in mammals from the standpoint of the neutral expectation. *The American Naturalist* 136:727-741. Hunt, G. 2012. Fitting and comparing models of phyletic evolution: random walks and beyond. *Paleobiology* 38:351-373.

**Examples**

```
y <- sim.GRW(ns = 20, ms = 0, vs = 1e-4, tt=seq(0, 1e6, length.out=20)) # per-year simulation
lynchD(y, gen.per.t = 1)
```

---

mle.GRW

*Analytical ML estimator for random walk and stasis models*


---

**Description**

Analytical ML estimator for random walk and stasis models

**Usage**

```
mle.GRW(y)
```

```
mle.URW(y)
```

```
mle.Stasis(y)
```

**Arguments**

`y` a paleoTS object

**Value**

a vector of `mstep` and `vstep` for `mle.GRW`, `vstep` for `mle.URW`, and `theta` and `omega` for `mle.Stasis`

**Functions**

- `mle.URW()`: ML parameter estimates for URW model
- `mle.Stasis()`: ML parameter estimates for Stasis model

**Note**

These analytical solutions assume even spacing of samples and equal sampling variance in each, which will usually be violated in real data. They are used here mostly to generate initial parameter estimates for numerical optimization; they not likely to be called directly by the user.

**See Also**

[fitSimple](#)

**Examples**

```
y <- sim.GRW(ms = 1, vs = 1)
w <- mle.GRW(y)
print(w)
```

---

opt.covTrack

*Fit a model in which a trait tracks a covariate*

---

**Description**

Fit a model in which a trait tracks a covariate

**Usage**

```
opt.covTrack(
  y,
  z,
  pool = TRUE,
  cl = list(fnscale = -1),
  meth = "L-BFGS-B",
  hess = FALSE
)

opt.joint.covTrack(
```

```

y,
z,
pool = TRUE,
cl = list(fnscale = -1),
meth = "L-BFGS-B",
hess = FALSE
)

```

### Arguments

y	a paloeTS object
z	a vector of covariate values
pool	if TRUE, sample variances are substituted with their pooled estimate
cl	optional control list, passed to <code>optim()</code>
meth	optimization algorithm, passed to <code>optim()</code>
hess	if TRUE, return standard errors of parameter estimates from the hessian matrix

### Details

In this model, changes in a trait are linearly related to changes in a covariate with a slope of  $b$  and residual variance  $evar$ :  $dx = b * dz + eps$ , where  $eps \sim N(0, evar)$ . This model was described, and applied to an example in which body size changes tracked changes in temperature, by Hunt et al. (2010).

For the AD version (`opt.covTrack`), a trait sequence of length  $ns$ , the covariate,  $z$ , can be of length  $ns - 1$ , interpreted as the vector of *changes*,  $dx$ . If  $z$  is of length  $ns$ , differences are taken and these are used as the  $dx$ 's, with a warning issued.

The Joint version (`opt.joint.covTrack`),  $z$  should be of length  $ns$  and there is an additional parameter that is the intercept of the linear relationship between trait and covariate. See warning below about using the Joint version.

### Value

a `paleoTSfit` object with the results of the model fitting

### Functions

- `opt.joint.covTrack()`: fits the `covTrack` model using the joint parameterization

### Warning

The Joint parameterization of this model can be fooled by temporal autocorrelation and, especially, trends in the trait and the covariate. The latter is tested for, but the AD parameterization is generally safer for this model.

## References

Hunt, G. S. Wicaksono, J. E. Brown, and K. G. Macleod. 2010. Climate-driven body size trends in the ostracod fauna of the deep Indian Ocean. *Palaeontology* 53(6): 1255-1268.

## See Also

[fitSimple](#)

## Examples

```
set.seed(13)
z <- c(1, 2, 2, 4, 0, 8, 2, 3, 1, 9, 4, 3)
x <- sim.covTrack(ns = 12, z = z, b = 0.5, evar = 0.2)
w.urw <- opt.URW(x)
w.cov <- opt.covTrack(x, z)
compareModels(w.urw, w.cov)
```

---

opt.GRW

*Fit evolutionary model using "AD" parameterization*

---

## Description

Fit evolutionary model using "AD" parameterization

## Usage

```
opt.GRW(  
  y,  
  pool = TRUE,  
  cl = list(fnscale = -1),  
  meth = "L-BFGS-B",  
  hess = FALSE  
)
```

```
opt.URW(  
  y,  
  pool = TRUE,  
  cl = list(fnscale = -1),  
  meth = "L-BFGS-B",  
  hess = FALSE  
)
```

```
opt.Stasis(  
  y,  
  pool = TRUE,  
  cl = list(fnscale = -1),  
  meth = "L-BFGS-B",
```

```
    hess = FALSE
  )

  opt.StrictStasis(y, pool = TRUE, cl = list(fnscale = -1), hess = FALSE)
```

### Arguments

y	a paleoTS object
pool	if TRUE, sample variances are substituted with their pooled estimate
cl	optional control list, passed to <code>optim()</code>
meth	optimization algorithm, passed to <code>optim()</code>
hess	if TRUE, return standard errors of parameter estimates from the hessian matrix

### Details

These functions use differences between consecutive populations in the time series in order to remove temporal autocorrelation. This is referred to as the "Ancestor-Descendant" or "AD" parameterization by Hunt [2008], and it is a REML approach (like phylogenetic independent contrasts). A full ML approach, called "Joint" was found to have generally better performance (Hunt, 2008) and generally should be used instead.

### Value

a paleoTSfit object with the model fitting results

### Functions

- `opt.URW()`: fit the URW model by the AD parameterization
- `opt.Stasis()`: fit the Stasis model by the AD parameterization
- `opt.StrictStasis()`: fit the Strict Stasis model by the AD parameterization

### Note

It is easier to use the convenience function `fitSimple`.

### References

Hunt, G. 2006. Fitting and comparing models of phyletic evolution: random walks and beyond. *Paleobiology* 32(4): 578-601.

### See Also

[fitSimple](#), [opt.joint.GRW](#)

**Examples**

```
x <- sim.GRW(ns = 20, ms = 1) # strong trend
plot(x)
w.grw <- opt.GRW(x)
w.urw <- opt.URW(x)
compareModels(w.grw, w.urw)
```

---

opt.GRW.shift

*Fit random walk model with shift(s) in generating parameters*


---

**Description**

Fit random walk model with shift(s) in generating parameters

**Usage**

```
opt.GRW.shift(y, ng = 2, minb = 7, model = 1, pool = TRUE, silent = FALSE)
```

**Arguments**

y	a paloeTS object
ng	number of segments in the sequence
minb	minimum number of populations in each segment
model	numeric, specifies exact evolutionary model; see Details
pool	if TRUE, sample variances are substituted with their pooled estimate
silent	logical, if TRUE, progress updates are suppressed

**Details**

Fits a model in which a sequence is divided into two or more segments and trait evolution proceeds as a general random walk, with each segment (potentially) getting its own generating parameters (mstep, vstep).

This function tests for shifts after each population, subject to the constraint that the number of populations in each segment is always  $\geq$  minb. The shiftpoint yielding the highest log-likelihood is returned as the solution, along with the log-likelihoods (all.logl of all tested shift points (GG).

Different variants of the model can be specified by the model argument:

- model = 1: mstep is separate across segments; vstep is shared
- model = 2: mstep is shared across segments; vstep is separate
- model = 3: mstep is set to zero (unbiased random walk); vstep is separate across segments
- model = 4: mstep and vstep are both separate across segments

**Value**

a paleoTSfit object

**See Also**

[sim.GRW.shift](#)

**Examples**

```
x <- sim.GRW.shift(ns = c(15,15), ms = c(0, 1), vs = c(0.1,0.1))
w.sep <- opt.GRW.shift(x, ng = 2, model = 4)
w.sameVs <- opt.GRW.shift(x, ng = 2, model = 1)
compareModels(w.sep, w.sameVs)
plot(x)
abline(v = x$tt[16], lwd = 3) # actual shift point
abline(v = x$tt[w.sameVs$par["shift1"]], lty = 3, col = "red", lwd = 2) # inferred shift point
```

---

opt.joint.GRW

*Fit evolutionary models using the "Joint" parameterization*

---

**Description**

Fit evolutionary models using the "Joint" parameterization

**Usage**

```
opt.joint.GRW(
  y,
  pool = TRUE,
  cl = list(fnscale = -1),
  meth = "L-BFGS-B",
  hess = FALSE
)
```

```
opt.joint.URW(
  y,
  pool = TRUE,
  cl = list(fnscale = -1),
  meth = "L-BFGS-B",
  hess = FALSE
)
```

```
opt.joint.Stasis(
  y,
  pool = TRUE,
  cl = list(fnscale = -1),
  meth = "L-BFGS-B",
```

```

    hess = FALSE
  )

  opt.joint.StrictStasis(y, pool = TRUE, cl = list(fnscale = -1), hess = FALSE)

```

### Arguments

y	a paleoTS object
pool	if TRUE, sample variances are substituted with their pooled estimate
cl	optional control list, passed to <code>optim()</code>
meth	optimization algorithm, passed to <code>optim()</code>
hess	if TRUE, return standard errors of parameter estimates from the hessian matrix

### Details

These functions use the joint distribution of population means to fit models using a full maximum-likelihood approach. This approach was found to have somewhat better performance than the "AD" approach, especially for noisy trends (Hunt, 2008).

### Value

a paleoTSfit object with the model fitting results

### Functions

- `opt.joint.URW()`: fit the URW model by the Joint parameterization
- `opt.joint.Stasis()`: fit the Stasis model by the Joint parameterization
- `opt.joint.StrictStasis()`: fit the Strict Stasis model by the Joint parameterization

### Note

It is easier to use the convenience function `fitSimple`.

### References

Hunt, G., M. J. Hopkins and S. Lidgard. 2015. Simple versus complex models of trait evolution and stasis as a response to environmental change. *Proc. Natl. Acad. Sci. USA* 112(16): 4885-4890.

### See Also

[fitSimple](#), [opt.GRW](#)

### Examples

```

x <- sim.GRW(ns = 20, ms = 1) # strong trend
plot(x)
w.grw <- opt.joint.GRW(x)
w.urw <- opt.joint.URW(x)
compareModels(w.grw, w.urw)

```

---

 opt.joint.OU

 Fit Ornstein-Uhlenbeck model using the "Joint" parameterization
 

---

**Description**

Fit Ornstein-Uhlenbeck model using the "Joint" parameterization

**Usage**

```
opt.joint.OU(
  y,
  pool = TRUE,
  cl = list(fnscale = -1),
  meth = "L-BFGS-B",
  hess = FALSE
)
```

**Arguments**

y	a paleoTS object
pool	if TRUE, sample variances are substituted with their pooled estimate
cl	optional control list, passed to <code>optim()</code>
meth	optimization algorithm, passed to <code>optim()</code>
hess	if TRUE, return standard errors of parameter estimates from the hessian matrix

**Details**

This function fits an Ornstein-Uhlenbeck (OU) model to time-series data. The OU model has four generating parameters: an ancestral trait value (`anc`), an optimum value (`theta`), the strength of attraction to the optimum (`alpha`), and a parameter that reflects the tendency of traits to diffuse (`vstep`). In a microevolutionary context, these parameters can be related to natural selection and genetic drift; see Hunt et al. (2008).

**Value**

a `paleoTSfit` object with the model fitting results

**Note**

It is easier to use the convenience function `fitSimple`. Note also that preliminary work found that the "AD" parameterization did not perform as well for the OU model and thus it is not implemented here.

**References**

Hunt, G., M. A. Bell and M. P. Travis. 2008. Evolution toward a new adaptive optimum: phenotypic evolution in a fossil stickleback lineage. *Evolution* 62(3): 700-710.

**See Also**

[fitSimple](#), [opt.joint.GRW](#)

**Examples**

```
x <- sim.OU(vs = 0.5) # most defaults OK
w <- opt.joint.OU(x)
plot(x, modelFit = w)
```

---

opt.punc

*Fit a model of trait evolution with specified punctuation(s)*

---

**Description**

Fit a model of trait evolution with specified punctuation(s)

**Usage**

```
opt.punc(
  y,
  gg,
  pool = TRUE,
  cl = list(fnscale = -1),
  meth = "L-BFGS-B",
  hess = FALSE,
  oshare
)
```

```
opt.joint.punc(
  y,
  gg,
  pool = TRUE,
  cl = list(fnscale = -1),
  meth = "L-BFGS-B",
  hess = FALSE,
  oshare
)
```

**Arguments**

y	a paleoTS object
gg	vector of indices indicating different segments
pool	if TRUE, sample variances are substituted with their pooled estimate
cl	optional control list, passed to <code>optim()</code>
meth	optimization algorithm, passed to <code>optim()</code>
hess	if TRUE, return standard errors of parameter estimates from the
oshare	logical, if TRUE, variance assumed to be shared (equal) across segments

**Details**

The sequence is divided into segments, which are separated by punctuations. Means for each segment are given by the vector `theta` with variances given by the vector `omega` (or a single value if `oshare = TRUE`). This function calls `optim` to numerically fit this model to a time-series, `y`.

**Value**

a `paleoTSfit` object with the results of the model fitting

**Functions**

- `opt.joint.punc()`: fits the punctuation model using the joint parameterization

**Note**

These functions would be used in the uncommon situation in which there is a prior hypothesis as to where the punctuation(s) take place. Normally users will instead use the function `fitGpunc`, which uses these functions to fit a range of possible timings for the punctuations.

**See Also**

[fitGpunc](#)

**Examples**

```
x <- sim.punc(ns = c(15, 15), theta = c(0,3), omega = c(0.1, 0.1))
w.sta <- fitSimple(x, model = "Stasis", method = "Joint")
w.punc <- opt.joint.punc(x, gg = rep(1:2, each = 15), oshare = TRUE)
compareModels(w.sta, w.punc)
```

---

opt.ssm.GRW

*Fit evolutionary models using state-space models (SSM)*

---

**Description**

Fit evolutionary models using state-space models (SSM)

**Usage**

```
opt.ssm.GRW(y, pool = TRUE, cl = list(fnscale = -1), hess = FALSE)
opt.ssm.URW(y, pool = TRUE, cl = list(fnscale = -1), hess = FALSE)
opt.ssm.Stasis(y, pool = TRUE, cl = list(fnscale = -1), hess = FALSE)
opt.ssm.StrictStasis(y, pool = TRUE, cl = list(fnscale = -1), hess = FALSE)
opt.ssm.OU(y, pool = TRUE, cl = list(fnscale = -1), hess = FALSE)
```

```

opt.ssm.ACDC(y, pool = TRUE, cl = list(fnscale = -1), hess = FALSE)

opt.ssm.covOU(y, z, pool = TRUE, cl = list(fnscale = -1), hess = FALSE)

opt.ssm.URWshift(y, gg, pool = TRUE, cl = list(fnscale = -1), hess = FALSE)

opt.ssm.covOU_vshift(
  y,
  z,
  gg,
  pool = TRUE,
  cl = list(fnscale = -1),
  hess = FALSE
)

```

### Arguments

<code>y</code>	a paleoTS object
<code>pool</code>	if TRUE, sample variances are substituted with their pooled estimate
<code>cl</code>	optional control list, passed to <code>optim()</code>
<code>hess</code>	if TRUE, return standard errors of parameter estimates from the hessian matrix
<code>z</code>	a covariate vector, used only for the covOU models
<code>gg</code>	a grouping vector, used only for the URWshift and covOU_vshift models

### Details

These functions use a state space model formulation to compute likelihoods and fit models.

Functions to fit the OU covariate tracking models (`covOU`, `covOU_vshift`) require a covariate argument, `z`. At present, only the OU covariate tracking with a shift in the step variance (`covOU_vshift`) requires the grouping vector argument (`gg`).

### Value

a `paleoTSfit` object with the model fitting results

### Note

For GRW, URW, Stasis, StrictStasis, ACDC and OU models, it will likely be easier to use the convenience function `fitSimple` with argument `method = "SSM"`.

The grouping vector, `gg`, is a vector of length equal to the number of samples. It has one element for each sample and takes integer value from 1 to the number of sample groups separated by shiftpoints. See the example below.

### See Also

[fitSimple](#), [opt.joint.GRW](#)

**Examples**

```

y <- sim.GRW(ns = 30, vs = 2)
w1 <- opt.ssm.URW(y)
gg <- rep(1:2, each = 15) # shift occurs immediately after sample 15
w2 <- opt.ssm.URWshift(y, gg = gg) # test model in which the step variance shifts
compareModels(w1, w2)

```

---

plot.paleoTS

*Plot a paleoTS object*


---

**Description**

Plot a paleoTS object

**Usage**

```

## S3 method for class 'paleoTS'
plot(
  x,
  nse = 1,
  pool = FALSE,
  add = FALSE,
  modelFit = NULL,
  pch = 21,
  pt.bg = "white",
  lwd = 1.5,
  ylim = NULL,
  xlab = NULL,
  ylab = NULL,
  add.label = TRUE,
  ...
)

```

**Arguments**

x	a paleoTS object
nse	the number of standard errors represented by the error bars on the plot; defaults to 1
pool	logical indicating if variances should be pooled across samples for the purposes of displaying error bars; defaults to FALSE
add	logical, if TRUE, adds to existing plot
modelFit	optional model fit from fitting functions
pch	plotting symbol, defaults to 19
pt.bg	color fill for points, defaults to white
lwd	line width, defaults to 1.5

ylim	optional, y-limits of the plot
xlab	optional, label for x-axis
ylab	optional, label for y-axis
add.label	logical, if TRUE, writes label from x at the top of the plot
...	other arguments passed to plotting functions

**Value**

none.

**Examples**

```
x <- sim.GRW(ns = 30)
w <- fitSimple(x, model = "GRW", method = "Joint")
plot(x, modelFit = w)
plot(x, xlab = "Time [Myr]", ylab = "Body length [mm]", pch = 22, pt.bg = "blue")
```

---

pool.var	<i>Compute a pooled variance</i>
----------	----------------------------------

---

**Description**

Computes a pooled variance from samples in a paleontological time-series

**Usage**

```
pool.var(y, nn = NULL, minN = NULL, ret.paleoTS = FALSE)
```

**Arguments**

y	either a paleoTS object, or a vector of sample variances
nn	a vector of sample sizes
minN	sample size below which variances are replaced with pooled variances. See Details.
ret.paleoTS	if TRUE, a paleoTS object is returned. If FALSE, the value of the pooled variance is returned.

**Details**

A pooled variance of a set of populations is the weighted average of the individual variances of the populations, with the weight for each population equal to its sample size minus one.

For many kinds of traits, variation levels tend to be similar among closely related populations. When this is true and sample sizes are low, much of the observed differences in variance among samples will be due to the high noise of estimated the variances. Replacing the observed variances of all populations (or only those with  $nn < minN$ ) with the estimated pooled variance can reduce this noise.

**Value**

if `ret.paleoTS = TRUE` a paleoTS object with all (or some) variances replaced with the pooled variance; otherwise the pooled variance

**Examples**

```
data(cantius_L)
cant_all <- pool.var(cantius_L, ret.paleoTS = TRUE) # replace all variances with pooled variance
cant_n5 <- pool.var(cantius_L, minN = 5, ret.paleoTS = TRUE) # replace only pops with n < 5
```

---

```
print.paleoTSfit      Print a paleoTSfit object
```

---

**Description**

Print a paleoTSfit object

**Usage**

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

**Arguments**

<code>x</code>	a paleoTSfit object
<code>...</code>	other arguments for other print methods

**Value**

None; this function is used only to print

**Examples**

```
y <- sim.punc(theta = c(0, 2), omega = c(0.1, 0.1))
wg <- fitGpunc(y)
print(wg)
```

---

read.paleoTS	<i>Read a text-file with data from a paleontological time-series</i>
--------------	--

---

**Description**

Read a text-file with data from a paleontological time-series

**Usage**

```
read.paleoTS(file = NULL, oldest = "first", reset.time = TRUE, ...)
```

**Arguments**

file	file name; if not supplied, an interactive window prompts the user to navigate to the text file
oldest	"first" if samples are in order from oldest to youngest, "last" if the opposite
reset.time	logical; see <a href="#">as.paleoTS</a>
...	other arguments, passed to read.table

**Details**

This function reads a text file with a specified format and converts it into a paleoTS object. It will often be the easiest way for users to import their own data. The text file should have four columns without headers, in this order: sample size, sample means, sample variances, sample ages.

**Value**

a paleoTS object

**See Also**

[as.paleoTS](#)

---

sim.covTrack	<i>Simulate trait evolution that tracks a covariate</i>
--------------	---

---

**Description**

Simulate trait evolution that tracks a covariate

**Usage**

```
sim.covTrack(
  ns = 20,
  b = 1,
  evar = 0.1,
  z,
  nn = rep(20, times = ns),
  tt = 0:(ns - 1),
  vp = 1
)
```

**Arguments**

ns	number of populations in a sequence
b	slope of the relationship between the change in the covariate and the change in the trait
evar	residual variance of the same relationship
z	vector of covariate that the trait tracks
nn	vector of sample sizes for populations
tt	vector of times (ages) for populations
vp	phenotypic trait variance within each population

**Details**

In this model, changes in a trait are linearly related to changes in a covariate with a slope of  $b$  and residual variance  $evar$ :  $dx = b * dz + eps$ , where  $eps \sim N(0, evar)$ . This model was described, and applied to an example in which body size changes tracked changes in temperature, by Hunt et al. (2010).

**Value**

a paleoTS object

**Note**

For a trait sequence of length  $ns$ , the covariate,  $z$ , can be of length  $ns - 1$ , in which case it is interpreted as the vector of *changes*,  $dz$ . If  $z$  is of length  $ns$ , differences are taken and these are used as the  $dz$ 's.

**References**

Hunt, G, S. Wicaksono, J. E. Brown, and K. G. Macleod. 2010. Climate-driven body size trends in the ostracod fauna of the deep Indian Ocean. *Palaeontology* 53(6): 1255-1268.

**Examples**

```

set.seed(13)
z <- c(1, 2, 2, 4, 0, 8, 2, 3, 1, 9, 4, 3)
x <- sim.covTrack(ns = 12, z = z, b = 0.5, evar = 0.2)
plot(x, ylim = c(-1, 10))
lines(x$tt, z, col = "blue")

```

sim.GRW

*Simulate random walk or directional time-series for trait evolution***Description**

Simulate random walk or directional time-series for trait evolution

**Usage**

```
sim.GRW(ns = 20, ms = 0, vs = 0.1, vp = 1, nn = rep(20, ns), tt = 0:(ns - 1))
```

**Arguments**

ns	number of populations in the sequence
ms	mean of evolutionary "steps"
vs	variance of evolutionary "steps"
vp	phenotypic variance within populations
nn	vector of population sample sizes
tt	vector of population times (ages)

**Details**

The general random walk model considers time in discrete steps. At each time step, an evolutionary change is drawn at random from a distribution of possible evolutionary "steps." It turns out that the long-term dynamics of an evolving lineage depend only on the mean and variance of this step distribution. The former, *mstep*, determined the directionality in a sequence and the latter, *vstep*, determines its volatility.

**Value**

a paleoTS object

**Note**

This function simulates an unbiased random walk if *ms* is equal to zero and a general (or biased) random walk otherwise.

**See Also**

[sim.Stasis](#), [sim.OU](#), [as.paleoTS](#)

**Examples**

```
x.grw <- sim.GRW(ms = 0.5)
x.urw <- sim.GRW(ms = 0)
plot(x.grw, ylim = range(c(x.grw$mm, x.urw$mm)))
plot(x.urw, add = TRUE, col = "blue")
legend(x = "topleft", c("GRW", "URW"), col = c("black", "blue"), lty = 1)
```

---

sim.GRW.shift	<i>Simulate (general) random walk with shift(s) in generating parameters</i>
---------------	--

---

**Description**

Simulate (general) random walk with shift(s) in generating parameters

**Usage**

```
sim.GRW.shift(
  ns = c(10, 10),
  ms = c(0, 1),
  vs = c(0.5, 0.5),
  nn = rep(30, sum(ns)),
  tt = 0:(sum(ns) - 1),
  vp = 1
)
```

**Arguments**

ns	vector of the number of samples in each segment
ms	vector of mean step parameter in each segment
vs	vector of step variance parameter in each segment
nn	vector of sample sizes, one for each population
tt	vector of samples times (ages)
vp	phenotypic variance in each sample

**Details**

Simulates under a model in which a sequence is divided into two or more segments. Trait evolution proceeds as a general random walk, with each segment getting its own generating parameters (mstep, vstep).

**Value**

a paleoTS object with the simulated time-series

**See Also**

[sim.GRW](#), [sim.sgs](#), [opt.GRW.shift](#)

**Examples**

```
x <- sim.GRW.shift(ns = c(10,10,10), ms = c(0, 1, 0), vs = c(0.1,0.1,0.1))
plot(x)
abline(v = c(9.5, 19.5), lty = 3, lwd = 2, col = "blue") # shows where dynamics shift
text(c(5, 15, 25), c(2,2,2), paste("segment", 1:3, sep = " "), col = "blue")
```

---

sim.OU

---

*Simulate an Ornstein-Uhlenbeck time-series*


---

**Description**

Simulate an Ornstein-Uhlenbeck time-series

**Usage**

```
sim.OU(
  ns = 20,
  anc = 0,
  theta = 10,
  alpha = 0.3,
  vstep = 0.1,
  vp = 1,
  nn = rep(20, ns),
  tt = 0:(ns - 1)
)
```

**Arguments**

ns	number of populations in the sequence
anc	ancestral phenotype
theta	OU optimum (long-term mean)
alpha	strength of attraction to the optimum
vstep	step variance
vp	phenotypic variance of each sample
nn	vector of sample sizes
tt	vector of sample times (ages)

**Details**

This function simulates an Ornstein-Uhlenbeck (OU) process. In microevolutionary terms, this models a population ascending a nearby peak in the adaptive landscape. The optimal trait value is `theta`, `alpha` indicates the strength of attraction to that peak (= strength of stabilizing selection around `theta`), `vstep` measures the random walk component (from genetic drift) and `anc` is the trait value at the start of the sequence.

**Value**

a paleoTS object

**References**

Hunt, G., M. A. Bell and M. P. Travis. 2008. Evolution toward a new adaptive optimum: phenotypic evolution in a fossil stickleback lineage. *Evolution* 62(3): 700-710.

**See Also**

[opt.joint.OU](#)

**Examples**

```
x1 <- sim.OU(alpha = 0.8) # strong alpha
x2 <- sim.OU(alpha = 0.1) # weaker alpha
plot(x1)
plot(x2, add = TRUE, col = "blue")
```

---

sim.punc

*Simulate a punctuated time-series*

---

**Description**

Simulates punctuated trait evolution with punctuations that are rapid relative to the spacing of samples. In practice, the time-series is divided into two or more segments, each of which has its own mean and variance.

**Usage**

```
sim.punc(
  ns = c(10, 10),
  theta = c(0, 1),
  omega = rep(0, length(theta)),
  nn = rep(30, sum(ns)),
  tt = 0:(sum(ns) - 1),
  vp = 1
)
```

**Arguments**

ns	vector of the number of samples in each segment
theta	vector of means, one for each segment
omega	vector of variances, one for each segment.
nn	vector of sample sizes, one for each population
tt	vector of times (ages), one for each population
vp	phenotypic variance within each population

**Details**

Segments are separated by punctuations. Population means in the  $i$ th segment are drawn randomly from a normal distribution with a mean equal to  $i$ th element of  $\theta$  and variance equal to the  $i$ th element of  $\omega$ . The magnitudes of punctuations are determined by the differences in adjacent  $\theta$  values.

**Value**

a paleoTS object with the simulated time-series.

**See Also**

[fitGpunc](#)

**Examples**

```
x <- sim.punc(ns = c(15, 15), theta = c(0,3), omega = c(0.1, 0.1))
plot(x)
```

---

sim.sgs

*Simulate protracted punctuation*

---

**Description**

This function simulates a punctuated change that is protracted enough that it is captured by multiple transitional populations. Trait evolution starts in stasis, shifts to a general random walk, and then shifts back into stasis.

**Usage**

```
sim.sgs(
  ns = c(20, 20, 20),
  theta = 0,
  omega = 1,
  ms = 1,
  vs = 0.1,
  nn = rep(30, sum(ns)),
  tt = 0:(sum(ns) - 1),
  vp = 1
)
```

**Arguments**

ns	vector with the number of samples in each segment
theta	trait mean for initial stasis segment
omega	trait variance for stasis segments

ms	step mean during random walk segment
vs	step variance during random walk segment
nn	vector of sample sizes for each population
tt	vector of times (ages) for each population
vp	phenotypic trait variance for each population

### Details

Trait evolution proceeds in three segments: Stasis, General random walk, stasis (sgs). The initial stasis segment has a mean of  $\theta$  and variance  $\omega$  before shifting in the second segment to a general random walk with parameters  $ms$  and  $vs$ . Finally, the third segment is a return to stasis, centered around the trait value of the last population of the random walk.

### Value

a paleoTS object

### Examples

```
x <- sim.sgs() # default values OK
plot(x)
```

---

sim.Stasis

*Simulate Stasis time-series for trait evolution*

---

### Description

Simulate Stasis time-series for trait evolution

### Usage

```
sim.Stasis(
  ns = 20,
  theta = 0,
  omega = 0,
  vp = 1,
  nn = rep(20, ns),
  tt = 0:(ns - 1)
)
```

### Arguments

ns	number of populations in the sequence
theta	mean of populations
omega	variance among populations
vp	phenotypic variance within populations
nn	vector of population sample sizes
tt	vector of population times (ages)

**Value**

a paleoTS object

**See Also**

[sim.GRW](#), [sim.OU](#), [as.paleoTS](#)

**Examples**

```
x <- sim.Stasis(omega = 0.5, vp = 0.1)
w.sta <- fitSimple(x, model = "Stasis")
w.ss <- fitSimple(x, model = "StrictStasis")
compareModels(w.sta, w.ss)
```

---

sim.Stasis.RW

*Simulate trait evolution with a mode shift*

---

**Description**

Trait evolution is modeled as a shift from a random walk (general or unbiased) to stasis, or vice versa.

**Usage**

```
sim.Stasis.RW(
  ns = c(20, 20),
  order = c("Stasis-RW", "RW-Stasis"),
  anc = 0,
  omega = 1,
  ms = 0,
  vs = 1,
  vp = 1,
  nn = 30,
  tt = NULL
)
```

**Arguments**

ns	vector of the number of samples in each segment
order	whether stasis or random walk come first, one of "Stasis-RW" or "RW-Stasis"
anc	starting trait value
omega	variance of stasis segment
ms	step mean during random walk segment
vs	step variance during random walk segment
vp	phenotypic trait variance for each population
nn	vector of sample sizes for each population
tt	vector of times (ages) for each population

**Details**

The anc argument is the starting trait value, and if the first segment is stasis, this is also the value of the stasis mean. When the first segment is a random walk, the stasis mean in the second segment is equal to the true trait mean at the end of the initial random walk.

**Value**

a paleoTSfit object

**See Also**

[fitModeShift](#)

**Examples**

```
x1 <- sim.Stasis.RW(omega = 0.1, ms = 5, order = "Stasis-RW")
x2 <- sim.Stasis.RW(omega = 0.1, ms = 5, order = "RW-Stasis")
plot(x1)
plot(x2, add = TRUE, col = "blue")
abline(v = 19, lty=3)
```

---

std.paleoTS

---

*Convert time-series to standard deviation units*


---

**Description**

Convert time-series to standard deviation units

**Usage**

```
std.paleoTS(y, center = c("mean", "start"))
```

**Arguments**

y	a paleoTS object
center	optional translation of time-series according to "mean" or "start"; see Details

**Details**

The standardization expresses each sample mean as the deviation from the overall mean, divided by the pooled within-sample standard deviation. Sample variances are also divided by the pooled sample variance.

Essentially, this converts paleontological time-series data into standard deviation units, similar to the computation of evolutionary rates in haldanes. This operation *does not* change the relative fit of models, but it does facilitate the comparison of parameter estimates across time-series of traits measured in different units.

If argument center = "start" the time-series is translated such that the trait mean of the first sample is zero.

### Value

the converted paleoTS object

### Examples

```
x <- sim.Stasis(ns = 8, theta = 1, omega = 4, vp = 2)
xs <- std.paleoTS(x, center = "start")
plot(x, ylim = range(c(x$mm, xs$mm)))
plot(xs, col = "red", add = TRUE)
legend(x = "topright", c("unstandardized", "standardized"), lty=1, col=c("black", "red"), cex=0.7)
```

---

sub.paleoTS

*Subsample a paleontological time-series*


---

### Description

Subsampling is done according to the supplied logical vector or, if none is supplied, as a proportion of samples, randomly chosen.

### Usage

```
sub.paleoTS(y, ok = NULL, k = 0.1, reset.time = TRUE)
```

### Arguments

y	a paleoTS object
ok	a logical vector, TRUE for populations to retain
k	proportion of samples to retain, with the samples chosen randomly
reset.time	if TRUE, resets the time so that the first population time is zero

### Value

the sub-sampled paleoTS object

### Examples

```
x <- sim.GRW(ns=20)
plot(x)
xs1 <- sub.paleoTS(x, k = 0.5)
plot(xs1, add = TRUE, col="green")
keep <- rep(c(TRUE, FALSE), 10)
xs2 <- sub.paleoTS(x, ok = keep)
plot(xs2, add = TRUE, col = "red")
```

---

test.var.het	<i>Test for heterogeneity of variances among samples in a time-series</i>
--------------	---

---

**Description**

Test for heterogeneity of variances among samples in a time-series

**Usage**

```
test.var.het(y, method = "Bartlett")
```

**Arguments**

y	a paleoTS object
method	test to use; currently only "Bartlett" is implemented

**Value**

a list with the test statistic, degrees of freedom, and p-value

**Note**

Most often, this function will be used to assess if it is reasonable to pool variances across samples using `pool.var`. A significant result means that the null hypothesis of equal variances across samples is rejected. Even in this case, however, it may still be preferable to pool variances, at least for some populations, if sample sizes are quite low.

**References**

Sokal, R. R and F. J. Rohlf. 1995. *Biometry* 3rd Ed.

**See Also**

[pool.var](#)

**Examples**

```
data(cantius_L)
test.var.het(cantius_L) # significant, but still may want to pool variances
```

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