

# Package ‘LDATS’

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**Title** Latent Dirichlet Allocation Coupled with Time Series Analyses

**Version** 0.3.0

**Description** Combines Latent Dirichlet Allocation (LDA) and Bayesian multinomial time series methods in a two-stage analysis to quantify dynamics in high-dimensional temporal data. LDA decomposes multivariate data into lower-dimension latent groupings, whose relative proportions are modeled using generalized Bayesian time series models that include abrupt change-points and smooth dynamics. The methods are described in Blei et al. (2003) <[doi:10.1162/jmlr.2003.3.4-5.993](https://doi.org/10.1162/jmlr.2003.3.4-5.993)>, Western and Kleykamp (2004) <[doi:10.1093/pan/023](https://doi.org/10.1093/pan/023)>, Venables and Ripley (2002, ISBN-13:978-0387954578), and Christensen et al. (2018) <[doi:10.1002/ecy.2373](https://doi.org/10.1002/ecy.2373)>.

**URL** <https://weecology.github.io/LDATS/>,  
<https://github.com/weecology/LDATS>

**BugReports** <https://github.com/weecology/LDATS/issues>

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**License** MIT + file LICENSE

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---

AICc	<i>Calculate AICc</i>
------	-----------------------

---

### Description

Calculate the small sample size correction of [AIC](#) for the input object.

### Usage

```
AICc(object)
```

### Arguments

object            An object for which [AIC](#) and [logLik](#) have defined methods.

### Value

numeric value of AICc.

### Examples

```
dat <- data.frame(y = rnorm(50), x = rnorm(50))
mod <- lm(dat)
AICc(mod)
```

---

autocorr_plot	<i>Produce the autocorrelation panel for the TS diagnostic plot of a parameter</i>
---------------	--

---

### Description

Produce a vanilla ACF plot using [acf](#) for the parameter of interest (rho or eta) as part of [TS\\_diagnostics\\_plot](#).

### Usage

```
autocorr_plot(x)
```

### Arguments

x                Vector of parameter values drawn from the posterior distribution, indexed to the iteration by the order of the vector.

**Value**

NULL.

**Examples**

```
autocorr_plot(rnorm(100, 0, 1))
```

---

check\_changepoints      *Check that a set of change point locations is proper*

---

**Description**

Check that the change point locations are numeric and conformable to interger values.

**Usage**

```
check_changepoints(changepoints = NULL)
```

**Arguments**

changepoints      Change point locations to evaluate.

**Value**

An error message is thrown if changepoints are not proper, else NULL.

**Examples**

```
check_changepoints(100)
```

---

check\_control      *Check that a control list is proper*

---

**Description**

Check that a list of controls is of the right class.

**Usage**

```
check_control(control, eclass = "list")
```

**Arguments**

control      Control list to evaluate.  
eclass      Expected class of the list to be evaluated.

**Value**

an error message is thrown if the input is improper, otherwise NULL.

**Examples**

```
check_control(list())
```

---

```
check_document_covariate_table
```

*Check that the document covariate table is proper*

---

**Description**

Check that the table of document-level covariates is conformable to a data frame and of the right size (correct number of documents) for the document-topic output from the LDA models.

**Usage**

```
check_document_covariate_table(  
  document_covariate_table,  
  LDA_models = NULL,  
  document_term_table = NULL  
)
```

**Arguments**

document_covariate_table	Document covariate table to evaluate.
LDA_models	Reference LDA model list (class LDA_set) that includes as its first element a properly fitted LDA model with a gamma slot with the document-topic distribution.
document_term_table	Optional input for checking when LDA_models is NULL

**Value**

An error message is thrown if document\_covariate\_table is not proper, else NULL.

**Examples**

```
data(rodents)  
check_document_covariate_table(rodents$document_covariate_table)
```

---

`check_document_term_table`*Check that document term table is proper*

---

**Description**

Check that the table of observations is conformable to a matrix of integers.

**Usage**

```
check_document_term_table(document_term_table)
```

**Arguments**

`document_term_table`

Table of observation count data (rows: documents, columns: terms. May be a class `matrix` or `data.frame` but must be conformable to a matrix of integers, as verified by [check\\_document\\_term\\_table](#).

**Value**

an error message is thrown if the input is improper, otherwise `NULL`.

**Examples**

```
data(rodents)
check_document_term_table(rodents$document_term_table)
```

---

`check_formula`*Check that a formula is proper*

---

**Description**

Check that formula is actually a [formula](#) and that the response and predictor variables are all included in data.

**Usage**

```
check_formula(data, formula)
```

**Arguments**

data	data.frame including [1] the time variable (indicated in timename), [2] the predictor variables (required by formula) and [3], the multinomial response variable (indicated in formula) as verified by <a href="#">check_timename</a> and <a href="#">check_formula</a> . Note that the response variables should be formatted as a data.frame object named as indicated by the response entry in the control list, such as gamma for a standard TS analysis on LDA output.
formula	formula to evaluate.

**Value**

An error message is thrown if formula is not proper, else NULL.

**Examples**

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
check_formula(data, gamma ~ 1)
```

---

check_formulas	<i>Check that formulas vector is proper and append the response variable</i>
----------------	--

---

**Description**

Check that the vector of formulas is actually formatted as a vector of [formula](#) objects and that the predictor variables are all included in the document covariate table.

**Usage**

```
check_formulas(formulas, document_covariate_table, control = list())
```

**Arguments**

formulas	Vector of the formulas to evaluate.
document_covariate_table	Document covariate table used to evaluate the availability of the data required by the formula inputs.
control	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <a href="#">TS_control</a> .



**Value**

An error message is thrown if formulas is not proper, else NULL.

**Examples**

```
data(rodents)
check_formulas(~ 1, rodents$document_covariate_table)
```

---

check_LDA_models	<i>Check that LDA model input is proper</i>
------------------	---

---

**Description**

Check that the LDA\_models input is either a set of LDA models (class LDA\_set, produced by [LDA\\_set](#)) or a singular LDA model (class LDA, produced by [LDA](#)).

**Usage**

```
check_LDA_models(LDA_models)
```

**Arguments**

LDA\_models      List of LDA models or singular LDA model to evaluate.

**Value**

An error message is thrown if LDA\_models is not proper, else NULL.

**Examples**

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDAs <- LDA_set(document_term_table, topics = 2, nseeds = 1)
LDA_models <- select_LDA(LDAs)
check_LDA_models(LDA_models)
```

---

check\_nchangepts    *Check that nchangepts vector is proper*

---

**Description**

Check that the vector of numbers of changepts is conformable to integers greater than 1.

**Usage**

```
check_nchangepts(nchangepts)
```

**Arguments**

nchangepts    Vector of the number of changepts to evaluate.

**Value**

An error message is thrown if nchangepts is not proper, else NULL.

**Examples**

```
check_nchangepts(0)  
check_nchangepts(2)
```

---

check\_seeds            *Check that nseeds value or seeds vector is proper*

---

**Description**

Check that the vector of numbers of seeds is conformable to integers greater than 0.

**Usage**

```
check_seeds(nseeds)
```

**Arguments**

nseeds            integer number of seeds (replicate starts) to use for each value of topics in the LDAs. Must be conformable to a positive integer value.

**Value**

an error message is thrown if the input is improper, otherwise NULL.

**Examples**

```
check_seeds(1)
check_seeds(2)
```

---

check_timename	<i>Check that the time vector is proper</i>
----------------	---

---

**Description**

Check that the vector of time values is included in the document covariate table and that it is either a integer-conformable or a date. If it is a date, the input is converted to an integer, resulting in the timestep being 1 day, which is often not desired behavior.

**Usage**

```
check_timename(document_covariate_table, timename)
```

**Arguments**

document_covariate_table	Document covariate table used to query for the time column.
timename	Column name for the time variable to evaluate.

**Value**

An error message is thrown if timename is not proper, else NULL.

**Examples**

```
data(rodents)
check_timename(rodents$document_covariate_table, "newmoon")
```

---

check_topics	<i>Check that topics vector is proper</i>
--------------	---

---

**Description**

Check that the vector of numbers of topics is conformable to integers greater than 1.

**Usage**

```
check_topics(topics)
```

**Arguments**

topics            Vector of the number of topics to evaluate for each model. Must be conformable to integer values.

**Value**

an error message is thrown if the input is improper, otherwise NULL.

**Examples**

```
check_topics(2)
```

---

check\_weights

*Check that weights vector is proper*

---

**Description**

Check that the vector of document weights is numeric and positive and inform the user if the average weight isn't 1.

**Usage**

```
check_weights(weights)
```

**Arguments**

weights            Vector of the document weights to evaluate, or TRUE for triggering internal weighting by document sizes.

**Value**

An error message is thrown if weights is not proper, else NULL.

**Examples**

```
check_weights(1)
wts <- runif(100, 0.1, 100)
check_weights(wts)
wts2 <- wts / mean(wts)
check_weights(wts2)
check_weights(TRUE)
```

---

count_trips	<i>Count trips of the ptMCMC particles</i>
-------------	--

---

### Description

Count the full trips (from one extreme temperature chain to the other and back again; Katzgraber *et al.* 2006) for each of the ptMCMC particles, as identified by their id on initialization.

This function was designed to work within `TS` and process the output of `est_changepoints` as a component of `diagnose_ptMCMC`, but has been generalized and would work with any output from a ptMCMC as long as `ids` is formatted properly.

### Usage

```
count_trips(ids)
```

### Arguments

`ids` matrix of identifiers of the particles in each chain for each iteration of the ptMCMC algorithm (rows: chains, columns: iterations).

### Value

list of [1] vector of within particle trip counts (`$trip_counts`), and [2] vector of within-particle average trip rates (`$trip_rates`).

### References

Katzgraber, H. G., S. Trebst, D. A. Huse. And M. Troyer. 2006. Feedback-optimized parallel tempering Monte Carlo. *Journal of Statistical Mechanics: Theory and Experiment* **3**:P03018 [link](#).

### Examples

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
data <- data[order(data[,"newmoon"]), ]
rho_dist <- est_changepoints(data, gamma ~ 1, 1, "newmoon", weights,
                           TS_control())
count_trips(rho_dist$ids)
```

---

diagnose\_ptMCMC

*Calculate ptMCMC summary diagnostics*


---

### Description

Summarize the step and swap acceptance rates as well as trip metrics from the saved output of a ptMCMC estimation.

### Usage

```
diagnose_ptMCMC(ptMCMCout)
```

### Arguments

ptMCMCout	Named list of saved data objects from a ptMCMC estimation including elements named <code>step_accepts</code> (matrix of logical outcomes of each step; rows: chains, columns: iterations), <code>swap_accepts</code> (matrix of logical outcomes of each swap; rows: chain pairs, columns: iterations), and <code>ids</code> (matrix of particle identifiers; rows: chains, columns: iterations). <code>ptMCMCout = NULL</code> indicates no use of ptMCMC and so the function returns NULL.
-----------	--

### Details

Within-chain step acceptance rates are averaged for each of the chains from the raw step acceptance histories (`ptMCMCout$step_accepts`) and between-chain swap acceptance rates are similarly averaged for each of the neighboring pairs of chains from the raw swap acceptance histories (`ptMCMCout$swap_accepts`). Trips are defined as movement from one extreme chain to the other and back again (Katzgraber *et al.* 2006). Trips are counted and turned to per-iteration rates using [count\\_trips](#).

This function was first designed to work within [TS](#) and process the output of [est\\_changepoints](#), but has been generalized and would work with any output from a ptMCMC as long as `ptMCMCout` is formatted properly.

### Value

list of [1] within-chain average step acceptance rates (`$step_acceptance_rate`), [2] average between-chain swap acceptance rates (`$swap_acceptance_rate`), [3] within particle trip counts (`$trip_counts`), and [4] within-particle average trip rates (`$trip_rates`).

### References

Katzgraber, H. G., S. Trebst, D. A. Huse. And M. Troyer. 2006. Feedback-optimized parallel tempering Monte Carlo. *Journal of Statistical Mechanics: Theory and Experiment* **3**:P03018 [link](#).

**Examples**

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
data <- data[order(data[, "newmoon"]), ]
rho_dist <- est_changepoints(data, gamma ~ 1, 1, "newmoon",
                             weights, TS_control())
diagnose_ptMCMC(rho_dist)
```

---

document_weights	<i>Calculate document weights for a corpus</i>
------------------	--

---

**Description**

Simple calculation of document weights based on the average number of words in a document within the corpus (mean value = 1).

**Usage**

```
document_weights(document_term_table)
```

**Arguments**

document\_term\_table

Table of observation count data (rows: documents, columns: terms. May be a class matrix or data.frame but must be conformable to a matrix of integers, as verified by [check\\_document\\_term\\_table](#).

**Value**

Vector of weights, one for each document, with the average sample receiving a weight of 1.0.

**Examples**

```
data(rodents)
document_weights(rodents$document_term_table)
```

---

ecdf_plot	<i>Produce the posterior distribution ECDF panel for the TS diagnostic plot of a parameter</i>
-----------	--

---

### Description

Produce a vanilla ECDF (empirical cumulative distribution function) plot using `ecdf` for the parameter of interest (`rho` or `eta`) as part of `TS_diagnostics_plot`. A horizontal line is added to show the median of the posterior.

### Usage

```
ecdf_plot(x, xlab = "parameter value")
```

### Arguments

<code>x</code>	Vector of parameter values drawn from the posterior distribution, indexed to the iteration by the order of the vector.
<code>xlab</code>	character value used to label the x axis.

### Value

NULL.

### Examples

```
ecdf_plot(rnorm(100, 0, 1))
```

---

est_changepoints	<i>Use ptMCMC to estimate the distribution of change point locations</i>
------------------	--

---

### Description

This function executes ptMCMC-based estimation of the change point location distributions for multinomial Time Series analyses.

### Usage

```
est_changepoints(  
  data,  
  formula,  
  nchangepoints,  
  timename,  
  weights,  
  control = list()  
)
```



**Arguments**

data	data.frame including [1] the time variable (indicated in <code>timename</code> ), [2] the predictor variables (required by <code>formula</code> ) and [3], the multinomial response variable (indicated in <code>formula</code> ) as verified by <code>check_timename</code> and <code>check_formula</code> . Note that the response variables should be formatted as a data.frame object named as indicated by the response entry in the control list, such as <code>gamma</code> for a standard TS analysis on LDA output.
formula	<code>formula</code> defining the regression between relationship the change points. Any predictor variable included must also be a column in <code>data</code> and any (multinomial) response variable must be a set of columns in <code>data</code> , as verified by <code>check_formula</code> .
nchangepoints	integer corresponding to the number of change points to include in the model. 0 is a valid input (corresponding to no change points, so a singular time series model), and the current implementation can reasonably include up to 6 change points. The number of change points is used to dictate the segmentation of the time series into chunks fit with separate models dictated by <code>formula</code> .
timename	character element indicating the time variable used in the time series.
weights	Optional class <code>numeric</code> vector of weights for each document. Defaults to <code>NULL</code> , translating to an equal weight for each document. When using <code>multinom_TS</code> in a standard LDATS analysis, it is advisable to weight the documents by their total size, as the result of LDA is a matrix of proportions, which does not account for size differences among documents. For most models, a scaling of the weights (so that the average is 1) is most appropriate, and this is accomplished using <code>document_weights</code> .
control	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <code>TS_control</code> .

**Value**

List of saved data objects from the ptMCMC estimation of change point locations (unless `nchangepoints` is 0, then `NULL` is returned).

**Examples**

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
formula <- gamma ~ 1
nchangepoints <- 1
control <- TS_control()
data <- data[order(data[, "newmoon"]), ]
rho_dist <- est_changepoints(data, formula, nchangepoints, "newmoon",
                             weights, control)
```

---

est_regressors	<i>Estimate the distribution of regressors, unconditional on the change point locations</i>
----------------	---

---

### Description

This function uses the marginal posterior distributions of the change point locations (estimated by [est\\_changepoints](#)) in combination with the conditional (on the change point locations) posterior distributions of the regressors (estimated by [multinom\\_TS](#)) to estimate the marginal posterior distribution of the regressors, unconditional on the change point locations.

### Usage

```
est_regressors(rho_dist, data, formula, timename, weights, control = list())
```

### Arguments

rho_dist	List of saved data objects from the ptMCMC estimation of change point locations (unless nchangepoints is 0, then NULL) returned from <a href="#">est_changepoints</a> .
data	data.frame including [1] the time variable (indicated in timename), [2] the predictor variables (required by formula) and [3], the multinomial response variable (indicated in formula) as verified by <a href="#">check_timename</a> and <a href="#">check_formula</a> . Note that the response variables should be formatted as a data.frame object named as indicated by the response entry in the control list, such as gamma for a standard TS analysis on LDA output.
formula	<a href="#">formula</a> defining the regression between relationship the change points. Any predictor variable included must also be a column in data and any (multinomial) response variable must be a set of columns in data, as verified by <a href="#">check_formula</a> .
timename	character element indicating the time variable used in the time series.
weights	Optional class numeric vector of weights for each document. Defaults to NULL, translating to an equal weight for each document. When using <a href="#">multinom_TS</a> in a standard LDATS analysis, it is advisable to weight the documents by their total size, as the result of LDA is a matrix of proportions, which does not account for size differences among documents. For most models, a scaling of the weights (so that the average is 1) is most appropriate, and this is accomplished using <a href="#">document_weights</a> .
control	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <a href="#">TS_control</a> .

## Details

The general approach follows that of Western and Kleykamp (2004), although we note some important differences. Our regression models are fit independently for each chunk (segment of time), and therefore the variance-covariance matrix for the full model has 0 entries for covariances between regressors in different chunks of the time series. Further, because the regression model here is a standard (non-hierarchical) softmax (Ripley 1996, Venables and Ripley 2002, Bishop 2006), there is no error term in the regression (as there is in the normal model used by Western and Kleykamp 2004), and so the posterior distribution used here is a multivariate normal, as opposed to a multivariate t, as used by Western and Kleykamp (2004).

## Value

matrix of draws (rows) from the marginal posteriors of the coefficients across the segments (columns).

## References

- Bishop, C. M. 2006. *Pattern Recognition and Machine Learning*. Springer, New York, NY, USA.
- Ripley, B. D. 1996. *Pattern Recognition and Neural Networks*. Cambridge University Press, Cambridge, UK.
- Venables, W. N. and B. D. Ripley. 2002. *Modern and Applied Statistics with S*. Fourth Edition. Springer, New York, NY, USA.
- Western, B. and M. Kleykamp. 2004. A Bayesian change point model for historical time series analysis. *Political Analysis* **12**:354-374. [link](#).

## Examples

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
formula <- gamma ~ 1
nchangepoints <- 1
control <- TS_control()
data <- data[order(data[, "newmoon"]), ]
rho_dist <- est_changepoints(data, formula, nchangepoints, "newmoon",
                             weights, control)
eta_dist <- est_regressors(rho_dist, data, formula, "newmoon", weights,
                           control)
```

---

expand_TS	<i>Expand the TS models across the factorial combination of LDA models, formulas, and number of change points</i>
-----------	---

---

### Description

Expand the completely crossed combination of model inputs: LDA model results, formulas, and number of change points.

### Usage

```
expand_TS(LDA_models, formulas, nchangepts)
```

### Arguments

LDA_models	List of LDA models (class LDA_set, produced by <a href="#">LDA_set</a> ) or a singular LDA model (class LDA, produced by <a href="#">LDA</a> ).
formulas	Vector of <a href="#">formula</a> (s) for the continuous (non-change point) component of the time series models. Any predictor variable included in a formula must also be a column in the document_covariate_table. Each element (formula) in the vector is evaluated for each number of change points and each LDA model.
nchangepts	Vector of integers corresponding to the number of change points to include in the time series models. 0 is a valid input corresponding to no change points ( <i>i.e.</i> , a singular time series model), and the current implementation can reasonably include up to 6 change points. Each element in the vector is the number of change points used to segment the data for each formula (entry in formulas) component of the TS model, for each selected LDA model.

### Value

Expanded data.frame table of the three values (columns) for each unique model run (rows): [1] the LDA model (indicated as a numeric element reference to the LDA\_models object), [2] the regressor formula, and [3] the number of changepts.

### Examples

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDAs <- LDA_set(document_term_table, topics = 2:3, nseeds = 2)
LDA_models <- select_LDA(LDAs)
weights <- document_weights(document_term_table)
formulas <- c(~ 1, ~ newmoon)
nchangepts <- 0:1
expand_TS(LDA_models, formulas, nchangepts)
```

---

iftrue	<i>Replace if TRUE</i>
--------	------------------------

---

**Description**

If the focal input is TRUE, replace it with alternative.

**Usage**

```
iftrue(x = TRUE, alt = NULL)
```

**Arguments**

x	Focal input.
alt	Alternative value.

**Value**

x if not TRUE, alt otherwise.

**Examples**

```
iftrue()
iftrue(TRUE, 1)
iftrue(2, 1)
iftrue(FALSE, 1)
```

---

jornada	<i>Jornada rodent data</i>
---------	----------------------------

---

**Description**

Counts of 17 rodent species across 24 sampling events, with the count being the total number observed across three trapping webs (146 traps in total) (Lightfoot *et al.* 2012).

**Usage**

```
jornada
```

**Format**

A list of two data.frame-class objects with rows corresponding to documents (sampling events). One element is the document term table (called `document_term_table`), which contains counts of the species (terms) in each sample (document), and the other is the document covariate table (called `document_covariate_table`) with columns of covariates (time step, year, season).

**Source**

<https://lter.jornada.nmsu.edu/data-catalog/>

**References**

Lightfoot, D. C., A. D. Davidson, D. G. Parker, L. Hernandez, and J. W. Laundre. 2012. Bottom-up regulation of desert grassland and shrubland rodent communities: implications of species-specific reproductive potentials. *Journal of Mammalogy* **93**:1017-1028. [link](#).

---

LDATS

*Package to conduct two-stage analyses combining Latent Dirichlet Allocation with Bayesian Time Series models*

---

**Description**

Performs two-stage analysis of multivariate temporal data using a combination of Latent Dirichlet Allocation (Blei *et al.* 2003) and Bayesian Time Series models (Western and Kleykamp 2004) that we extend for multinomial data using softmax regression (Venables and Ripley 2002) following Christensen *et al.* (2018).

**Documentation**

[Technical mathematical manuscript](#)

[End-user-focused vignette worked example](#)

[Computational pipeline vignette](#)

[Comparison to Christensen \*et al.\*](#)

**References**

Blei, D. M., A. Y. Ng, and M. I. Jordan. 2003. Latent Dirichlet Allocation. *Journal of Machine Learning Research* **3**:993-1022. [link](#).

Christensen, E., D. J. Harris, and S. K. M. Ernest. 2018. Long-term community change through multiple rapid transitions in a desert rodent community. *Ecology* **99**:1523-1529. [link](#).

Venables, W. N. and B. D. Ripley. 2002. *Modern and Applied Statistics with S*. Fourth Edition. Springer, New York, NY, USA.

Western, B. and M. Kleykamp. 2004. A Bayesian change point model for historical time series analysis. *Political Analysis* **12**:354-374. [link](#).

---

`LDA_msg`*Create the model-running-message for an LDA*

---

**Description**

Produce and print the message for a given LDA model.

**Usage**

```
LDA_msg(mod_topics, mod_seeds, control = list())
```

**Arguments**

<code>mod_topics</code>	integer value corresponding to the number of topics in the model.
<code>mod_seeds</code>	integer value corresponding to the seed used for the model.
<code>control</code>	Class <code>LDA_controls</code> list of control parameters to be used in LDA (note that "seed" will be overwritten).

**Examples**

```
LDA_msg(mod_topics = 4, mod_seeds = 2)
```

---

`LDA_set`*Run a set of Latent Dirichlet Allocation models*

---

**Description**

For a given dataset consisting of counts of words across multiple documents in a corpus, conduct multiple Latent Dirichlet Allocation (LDA) models (using the Variational Expectation Maximization (VEM) algorithm; Blei *et al.* 2003) to account for [1] uncertainty in the number of latent topics and [2] the impact of initial values in the estimation procedure.

`LDA_set` is a list wrapper of [LDA](#) in the `topicmodels` package (Grun and Hornik 2011).

`check_LDA_set_inputs` checks that all of the inputs are proper for `LDA_set` (that the table of observations is conformable to a matrix of integers, the number of topics is an integer, the number of seeds is an integer and the controls list is proper).

**Usage**

```
LDA_set(document_term_table, topics = 2, nseeds = 1, control = list())
```

```
check_LDA_set_inputs(document_term_table, topics, nseeds, control)
```

**Arguments**

document_term_table	Table of observation count data (rows: documents, columns: terms. May be a class matrix or data.frame but must be conformable to a matrix of integers, as verified by <a href="#">check_document_term_table</a> ).
topics	Vector of the number of topics to evaluate for each model. Must be conformable to integer values.
nseeds	Number of seeds (replicate starts) to use for each value of topics. Must be conformable to integer value.
control	A list of parameters to control the running and selecting of LDA models. Values not input assume default values set by <a href="#">LDA_set_control</a> . Values for running the LDAs replace defaults in (LDAcontrol, see <a href="#">LDA</a> (but if seed is given, it will be overwritten; use <code>iseed</code> instead).

**Value**

LDA\_set: list (class: LDA\_set) of LDA models (class: LDA\_VEM). check\_LDA\_set\_inputs: an error message is thrown if any input is improper, otherwise NULL.

**References**

- Blei, D. M., A. Y. Ng, and M. I. Jordan. 2003. Latent Dirichlet Allocation. *Journal of Machine Learning Research* **3**:993-1022. [link](#).
- Grun B. and K. Hornik. 2011. topicmodels: An R Package for Fitting Topic Models. *Journal of Statistical Software* **40**:13. [link](#).

**Examples**

```
data(rodents)
lda_data <- rodents$document_term_table
r_LDA <- LDA_set(lda_data, topics = 2, nseeds = 2)
```

---

LDA\_set\_control

*Create control list for set of LDA models*


---

**Description**

This function provides a simple creation and definition of the list used to control the set of LDA models. It is set up to be easy to work with the existing control capacity of [LDA](#).

**Usage**

```
LDA_set_control(quiet = FALSE, measurer = AIC, selector = min, izeed = 2, ...)
```



**Arguments**

quiet	logical indicator of whether the model should run quietly.
measurer, selector	Function names for use in evaluation of the LDA models. measurer is used to create a value for each model and selector operates on the values to choose the model(s) to pass on.
iseed	integer initial seed for the model set.
...	Additional arguments to be passed to <a href="#">LDA</a> as a control input.

**Value**

list for controlling the LDA model fit.

**Examples**

```
LDA_set_control()
```

---

LDA\_TS

---

*Run a full set of Latent Dirichlet Allocations and Time Series models*


---

**Description**

Conduct a complete LDATS analysis (Christensen *et al.* 2018), including running a suite of Latent Dirichlet Allocation (LDA) models (Blei *et al.* 2003, Grun and Hornik 2011) via [LDA\\_set](#), selecting LDA model(s) via [select\\_LDA](#), running a complete set of Bayesian Time Series (TS) models (Western and Kleykamp 2004) via [TS\\_on\\_LDA](#) on the chosen LDA model(s), and selecting the best TS model via [select\\_TS](#).

`conform_LDA_TS_data` converts the data input to match internal and sub-function specifications.

`check_LDA_TS_inputs` checks that the inputs to LDA\_TS are of proper classes for a full analysis.

**Usage**

```
LDA_TS(  
  data,  
  topics = 2,  
  nseeds = 1,  
  formulas = ~1,  
  nchangepts = 0,  
  timename = "time",  
  weights = TRUE,  
  control = list()  
)
```

```

conform_LDA_TS_data(data, quiet = FALSE)

check_LDA_TS_inputs(
  data = NULL,
  topics = 2,
  nseeds = 1,
  formulas = ~1,
  nchangepoints = 0,
  timename = "time",
  weights = TRUE,
  control = list()
)

```

## Arguments

data	<p>Either a document term table or a list including at least a document term table (with the word "term" in the name of the element) and optionally also a document covariate table (with the word "covariate" in the name of the element).</p> <p>The document term table is a table of observation count data (rows: documents, columns: terms) that may be a <code>matrix</code> or <code>data.frame</code>, but must be conformable to a matrix of integers, as verified by <a href="#">check_document_term_table</a>.</p> <p>The document covariate table is a table of associated data (rows: documents, columns: time index and covariate options) that may be a <code>matrix</code> or <code>data.frame</code>, but must be conformable to a data table, as verified by <a href="#">check_document_covariate_table</a>. Every model needs a covariate to describe the time value for each document (in whatever units and whose name in the table is input in <code>timename</code>) that dictates the application of the change points. <b><i>If a covariate table is not provided, the model assumes the observations were equi-spaced in time.</i></b> All covariates named within specific models in <code>formulas</code> must be included.</p>
topics	Vector of the number of topics to evaluate for each model. Must be conformable to integer values.
nseeds	integer number of seeds (replicate starts) to use for each value of topics in the LDAs. Must be conformable to integer value.
formulas	Vector of <a href="#">formula</a> (s) for the continuous (non-change point) component of the time series models. Any predictor variable included in a formula must also be a column in the <code>document_covariate_table</code> . Each element (formula) in the vector is evaluated for each number of change points and each LDA model.
nchangepoints	Vector of integers corresponding to the number of change points to include in the time series models. 0 is a valid input corresponding to no change points ( <i>i.e.</i> , a singular time series model), and the current implementation can reasonably include up to 6 change points. Each element in the vector is the number of change points used to segment the data for each formula (entry in <code>formulas</code> ) component of the TS model, for each selected LDA model.
timename	character element indicating the time variable used in the time series. Defaults to "time". The variable must be integer-conformable or a Date. If the variable

	named is a Date, the input is converted to an integer, resulting in the timestep being 1 day, which is often not desired behavior.
weights	Optional input for overriding standard weighting for documents in the time series. Defaults to TRUE, translating to an appropriate weighting of the documents based on the size (number of words) each document (the result of LDA is a matrix of proportions, which does not account for size differences among documents. Alternatively can be NULL for an equal weighting among documents or a numeric vector.
control	A list of parameters to control the running and selecting of LDA and TS models. Values not input assume default values set by <code>LDA_TS_control</code> .
quiet	logical indicator for <code>conform_LDA_TS_data</code> to indicate if messages should be printed.

### Value

LDA\_TS: a class LDA\_TS list object including all fitted LDA and TS models and selected models specifically as elements "LDA models" (from `LDA_set`), "Selected LDA model" (from `select_LDA`), "TS models" (from `TS_on_LDA`), and "Selected TS model" (from `select_TS`).

`conform_LDA_TS_data`: a data list that is ready for analyses using the stage-specific functions.

`check_LDA_TS_inputs`: an error message is thrown if any input is improper, otherwise NULL.

### References

Blei, D. M., A. Y. Ng, and M. I. Jordan. 2003. Latent Dirichlet Allocation. *Journal of Machine Learning Research* **3**:993-1022. [link](#).

Christensen, E., D. J. Harris, and S. K. M. Ernest. 2018. Long-term community change through multiple rapid transitions in a desert rodent community. *Ecology* **99**:1523-1529. [link](#).

Grun B. and K. Hornik. 2011. topicmodels: An R Package for Fitting Topic Models. *Journal of Statistical Software* **40**:13. [link](#).

Western, B. and M. Kleykamp. 2004. A Bayesian change point model for historical time series analysis. *Political Analysis* **12**:354-374. [link](#).

### Examples

```
data(rodents)

mod <- LDA_TS(data = rodents, topics = 2, nseeds = 1, formulas = ~1,
              nchangepoints = 1, timename = "newmoon")

conform_LDA_TS_data(rodents)
check_LDA_TS_inputs(rodents, timename = "newmoon")
```

---

LDA\_TS\_control      *Create the controls list for the LDATS model*

---

### Description

Create and define a list of control options used to run the LDATS model, as implemented by [LDA\\_TS](#).

### Usage

```
LDA_TS_control(
  quiet = FALSE,
  measurer_LDA = AIC,
  selector_LDA = min,
  iseed = 2,
  memoise = TRUE,
  response = "gamma",
  lambda = 0,
  measurer_TS = AIC,
  selector_TS = min,
  ntemps = 6,
  penultimate_temp = 2^6,
  ultimate_temp = 1e+10,
  q = 0,
  nit = 10000,
  magnitude = 12,
  burnin = 0,
  thin_frac = 1,
  summary_prob = 0.95,
  seed = NULL,
  ...
)
```

### Arguments

quiet	logical indicator of whether the model should run quietly.
measurer_LDA, selector_LDA	Function names for use in evaluation of the LDA models. <code>measurer_LDA</code> is used to create a value for each model and <code>selector_LDA</code> operates on the values to choose the model.
iseed	integer initial seed for the LDA model set.
memoise	logical indicator of whether the multinomial functions should be memoised (via <code>memoise</code> ). Memoisation happens to both <code>multinom_TS</code> and <code>multinom_TS_chunk</code> .
response	character element indicating the response variable used in the time series. Should be set to "gamma" for LDATS.
lambda	numeric "weight" decay term used to set the prior on the regressors within each chunk-level model. Defaults to 0, corresponding to a fully vague prior.

measurer_TS, selector_TS	Function names for use in evaluation of the TS models. measurer_TS is used to create a value for each model and selector_TS operates on the values to choose the model.
ntemps	integer number of temperatures (chains) to use in the ptMCMC algorithm.
penultimate_temp	Penultimate temperature in the ptMCMC sequence.
ultimate_temp	Ultimate temperature in the ptMCMC sequence.
q	Exponent controlling the ptMCMC temperature sequence from the focal chain (reference with temperature = 1) to the penultimate chain. 0 (default) implies a geometric sequence. 1 implies squaring before exponentiating.
nit	integer number of iterations (steps) used in the ptMCMC algorithm.
magnitude	Average magnitude (defining a geometric distribution) for the proposed step size in the ptMCMC algorithm.
burnin	integer number of iterations to remove from the beginning of the ptMCMC algorithm.
thin_frac	Fraction of iterations to retain, from the ptMCMC. Must be (0, 1], and the default value of 1 represents no thinning.
summary_prob	Probability used for summarizing the posterior distributions (via the highest posterior density interval, see <a href="#">HPDinterval</a> ) of the TS model.
seed	Input to set.seed in the time series model for replication purposes.
...	Additional arguments to be passed to <a href="#">LDA</a> as a control input.

**Value**

list of control lists, with named elements LDAcontrol, TScontrol, and quiet.

**Examples**

```
LDA_TS_control()
```

---

logLik.LDA_VEM	<i>Calculate the log likelihood of a VEM LDA model fit</i>
----------------	--

---

**Description**

Imported but updated calculations from topicmodels package, as applied to Latent Dirichlet Allocation fit with Variational Expectation Maximization via [LDA](#).

**Usage**

```
## S3 method for class 'LDA_VEM'
logLik(object, ...)
```

**Arguments**

object            A LDA\_VEM-class object.  
 ...                Not used, simply included to maintain method compatibility.

**Details**

The number of degrees of freedom is 1 (for alpha) plus the number of entries in the document-topic matrix. The number of observations is the number of entries in the document-term matrix.

**Value**

Log likelihood of the model logLik, also with df (degrees of freedom) and nobs (number of observations) values.

**References**

Buntine, W. 2002. Variational extensions to EM and multinomial PCA. *European Conference on Machine Learning, Lecture Notes in Computer Science* **2430**:23-34. [link](#).  
 Grun B. and K. Hornik. 2011. topicmodels: An R Package for Fitting Topic Models. *Journal of Statistical Software* **40**:13. [link](#).  
 Hoffman, M. D., D. M. Blei, and F. Bach. 2010. Online learning for latent Dirichlet allocation. *Advances in Neural Information Processing Systems* **23**:856-864. [link](#).

**Examples**

```
data(rodents)
lda_data <- rodents$document_term_table
r_LDA <- LDA_set(lda_data, topics = 2)
logLik(r_LDA[[1]])
```

---

```
logLik.multinom_TS_fit
```

*Log likelihood of a multinomial TS model*

---

**Description**

Convenience function to simply extract the logLik element (and df and nobs) from a multinom\_TS\_fit object fit by multinom\_TS. Extends logLik from multinom to multinom\_TS\_fit objects.

**Usage**

```
## S3 method for class 'multinom_TS_fit'
logLik(object, ...)
```

**Arguments**

object            A multinom\_TS\_fit-class object.  
 ...              Not used, simply included to maintain method compatibility.

**Value**

Log likelihood of the model, as class logLik, with attributes df (degrees of freedom) and nobs (the number of weighted observations, accounting for size differences among documents).

**Examples**

```
data(rodents)
dtt <- rodents$document_term_table
lda <- LDA_set(dtt, 2, 1, list(quiet = TRUE))
dct <- rodents$document_covariate_table
dct$gamma <- lda[[1]]@gamma
weights <- document_weights(dtt)
mts <- multinom_TS(dct, formula = gamma ~ 1, changepoints = c(20,50),
                  timename = "newmoon", weights = weights)
logLik(mts)
```

---

logLik.TS\_fit

*Determine the log likelihood of a Time Series model*


---

**Description**

Convenience function to extract and format the log likelihood of a TS\_fit-class object fit by [multinom\\_TS](#).

**Usage**

```
## S3 method for class 'TS_fit'
logLik(object, ...)
```

**Arguments**

object            Class TS\_fit object to be evaluated.  
 ...              Not used, simply included to maintain method compatibility.

**Value**

Log likelihood of the model logLik, also with df (degrees of freedom) and nobs (number of observations) values.

**Examples**

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
TSmod <- TS(data, gamma ~ 1, nchangepts = 1, "newmoon", weights)
logLik(TSmod)
```

---

logsumexp

*Calculate the log-sum-exponential (LSE) of a vector*

---

**Description**

Calculate the exponent of a vector (offset by the max), sum the elements, calculate the log, remove the offset.

**Usage**

```
logsumexp(x)
```

**Arguments**

x                    numeric vector

**Value**

The LSE.

**Examples**

```
logsumexp(1:10)
```



---

memoise_fun	<i>Logical control on whether or not to memoise</i>
-------------	---

---

**Description**

This function provides a simple, logical toggle control on whether the function fun should be memoised via `memoise` or not.

**Usage**

```
memoise_fun(fun, memoise_tf = TRUE)
```

**Arguments**

fun	Function name to (potentially) be memoised.
memoise_tf	logical value indicating if fun should be memoised.

**Value**

fun, memoised if desired.

**Examples**

```
sum_memo <- memoise_fun(sum)
```

---

messageq	<i>Optionally generate a message based on a logical input</i>
----------	---

---

**Description**

Given the input to quiet, generate the message(s) in msg or not.

**Usage**

```
messageq(msg = NULL, quiet = FALSE)
```

**Arguments**

msg	character vector of the message(s) to generate or NULL. If more than one element is contained in msg, they are concatenated with a newline between.
quiet	logical indicator controlling if the message is generated.

**Examples**

```
messageq("hello")  
messageq("hello", TRUE)
```

---

`mirror_vcov`*Create a properly symmetric variance covariance matrix*

---

**Description**

A wrapper on `vcov` to produce a symmetric matrix. If the default matrix returned by `vcov` is symmetric it is returned simply. If it is not, in fact, symmetric (as occurs occasionally with `multinom` applied to proportions), the matrix is made symmetric by averaging the lower and upper triangles. If the relative difference between the upper and lower triangles for any entry is more than 0.1

**Usage**

```
mirror_vcov(x)
```

**Arguments**

`x` Model object that has a defined method for `vcov`.

**Value**

Properly symmetric variance covariance matrix.

**Examples**

```
dat <- data.frame(y = rnorm(50), x = rnorm(50))
mod <- lm(dat)
mirror_vcov(mod)
```

---

`modalvalue`*Determine the mode of a distribution*

---

**Description**

Find the most common entry in a vector. Ties are not allowed, the first value encountered within the modal set if there are ties is deemed the mode.

**Usage**

```
modalvalue(x)
```

**Arguments**

`x` numeric vector.

**Value**

Numeric value of the mode.

**Examples**

```
d1 <- c(1, 1, 1, 2, 2, 3)
modalvalue(d1)
```

---

multinom\_TS

*Fit a multinomial change point Time Series model*

---

**Description**

Fit a set of multinomial regression models (via [multinom](#), Venables and Ripley 2002) to a time series of data divided into multiple segments (a.k.a. chunks) based on given locations for a set of change points.

`check_multinom_TS_inputs` checks that the inputs to `multinom_TS` are of proper classes for an analysis.

**Usage**

```
multinom_TS(
  data,
  formula,
  changepoints = NULL,
  timename = "time",
  weights = NULL,
  control = list()
)

check_multinom_TS_inputs(
  data,
  formula = gamma ~ 1,
  changepoints = NULL,
  timename = "time",
  weights = NULL,
  control = list()
)
```

**Arguments**

`data` `data.frame` including [1] the time variable (indicated in `timename`), [2] the predictor variables (required by `formula`) and [3], the multinomial response variable (indicated in `formula`) as verified by [check\\_timename](#) and [check\\_formula](#). Note that the response variables should be formatted as a `data.frame` object

	named as indicated by the response entry in the control list, such as gamma for a standard TS analysis on LDA output. See Examples.
formula	<a href="#">formula</a> defining the regression between relationship the change points. Any predictor variable included must also be a column in data and any (multinomial) response variable must be a set of columns in data, as verified by <a href="#">check_formula</a> .
changepoints	Numeric vector indicating locations of the change points. Must be conformable to integer values. Validity checked by <a href="#">check_changepoints</a> and <a href="#">verify_changepoint_locations</a> .
timename	character element indicating the time variable used in the time series. Defaults to "time". The variable must be integer-conformable or a Date. If the variable named is a Date, the input is converted to an integer, resulting in the timestep being 1 day, which is often not desired behavior.
weights	Optional class numeric vector of weights for each document. Defaults to NULL, translating to an equal weight for each document. When using <code>multinom_TS</code> in a standard LDATS analysis, it is advisable to weight the documents by their total size, as the result of LDA is a matrix of proportions, which does not account for size differences among documents. For most models, a scaling of the weights (so that the average is 1) is most appropriate, and this is accomplished using <a href="#">document_weights</a> .
control	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <a href="#">TS_control</a> .

### Value

`multinom_TS`: Object of class `multinom_TS_fit`, which is a list of [1] chunk-level model fits ("chunk models"), [2] the total log likelihood combined across all chunks ("logLik"), and [3] a data.frame of chunk beginning and ending times ("logLik" with columns "start" and "end").

`check_multinom_TS_inputs`: an error message is thrown if any input is improper, otherwise NULL.

### References

Venables, W. N. and B. D. Ripley. 2002. *Modern and Applied Statistics with S*. Fourth Edition. Springer, New York, NY, USA.

### Examples

```
data(rodents)
dtt <- rodents$document_term_table
lda <- LDA_set(dtt, 2, 1, list(quiet = TRUE))
dct <- rodents$document_covariate_table
dct$gamma <- lda[[1]]@gamma
weights <- document_weights(dtt)
check_multinom_TS_inputs(dct, timename = "newmoon")
mts <- multinom_TS(dct, formula = gamma ~ 1, changepoints = c(20,50),
  timename = "newmoon", weights = weights)
```

---

multinom\_TS\_chunk      *Fit a multinomial Time Series model chunk*

---

## Description

Fit a multinomial regression model (via `multinom`, Ripley 1996, Venables and Ripley 2002) to a defined chunk of time (a.k.a. segment) [`chunk$start`, `chunk$end`] within a time series.

## Usage

```
multinom_TS_chunk(
  data,
  formula,
  chunk,
  timename = "time",
  weights = NULL,
  control = list()
)
```

## Arguments

<code>data</code>	Class <code>data.frame</code> object including the predictor and response variables.
<code>formula</code>	Formula as a <code>formula</code> or <code>character</code> object describing the chunk.
<code>chunk</code>	Length-2 vector of times: [1] <code>start</code> , the start time for the chunk and [2] <code>end</code> , the end time for the chunk.
<code>timename</code>	character element indicating the time variable used in the time series. Defaults to <code>"time"</code> . The variable must be integer-conformable or a <code>Date</code> . If the variable named is a <code>Date</code> , the input is converted to an integer, resulting in the timestep being 1 day, which is often not desired behavior.
<code>weights</code>	Optional class <code>numeric</code> vector of weights for each document. Defaults to <code>NULL</code> , translating to an equal weight for each document. When using <code>multinom_TS</code> in a standard LDATS analysis, it is advisable to weight the documents by their total size, as the result of <code>LDA</code> is a matrix of proportions, which does not account for size differences among documents. For most models, a scaling of the weights (so that the average is 1) is most appropriate, and this is accomplished using <code>document_weights</code> .
<code>control</code>	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <code>TS_control</code> .

## Value

Fitted model object for the chunk, of classes `multinom` and `nnet`.

## References

- Ripley, B. D. 1996. Pattern Recognition and Neural Networks. Cambridge.
- Venables, W. N. and B. D. Ripley. 2002. Modern Applied Statistics with S. Fourth edition. Springer.

## Examples

```
data(rodents)
dtt <- rodents$document_term_table
lda <- LDA_set(dtt, 2, 1, list(quiet = TRUE))
dct <- rodents$document_covariate_table
dct$gamma <- lda[[1]]@gamma
weights <- document_weights(dtt)
chunk <- c(start = 0, end = 100)
mtsc <- multinom_TS_chunk(dct, formula = gamma ~ 1, chunk = chunk,
                          timename = "newmoon", weights = weights)
```

---

normalize

*Normalize a vector*

---

## Description

Normalize a numeric vector to be on the scale of [0,1].

## Usage

```
normalize(x)
```

## Arguments

x                    numeric vector.

## Value

Normalized x.

## Examples

```
normalize(1:10)
```

---

package_chunk_fits	<i>Package the output of the chunk-level multinomial models into a multinom_TS_fit list</i>
--------------------	---

---

### Description

Takes the list of fitted chunk-level models returned from `TS_chunk_memo` (the memoised version of `multinom_TS_chunk` and packages it as a `multinom_TS_fit` object. This involves naming the model fits based on the chunk time windows, combining the log likelihood values across the chunks, and setting the class of the output object.

### Usage

```
package_chunk_fits(chunks, fits)
```

### Arguments

<code>chunks</code>	Data frame of start and end times for each chunk (row).
<code>fits</code>	List of chunk-level fits returned by <code>TS_chunk_memo</code> , the memoised version of <code>multinom_TS_chunk</code> .

### Value

Object of class `multinom_TS_fit`, which is a list of [1] chunk-level model fits, [2] the total log likelihood combined across all chunks, and [3] the chunk time data table.

### Examples

```
data(rodents)
dtt <- rodents$document_term_table
lda <- LDA_set(dtt, 2, 1, list(quiet = TRUE))
dct <- rodents$document_covariate_table
dct$gamma <- lda[[1]]@gamma
weights <- document_weights(dtt)
formula <- gamma ~ 1
changepoints <- c(20,50)
timename <- "newmoon"
TS_chunk_memo <- memoise_fun(multinom_TS_chunk, TRUE)
chunks <- prep_chunks(dct, changepoints, timename)
nchunks <- nrow(chunks)
fits <- vector("list", length = nchunks)
for (i in 1:nchunks){
  fits[[i]] <- TS_chunk_memo(dct, formula, chunks[i, ], timename,
                           weights, TS_control())
}
package_chunk_fits(chunks, fits)
```

---

package\_LDA\_set      *Package the output from LDA\_set*

---

## Description

Name the elements (LDA models) and set the class (LDA\_set) of the models returned by [LDA\\_set](#).

## Usage

```
package_LDA_set(mods, mod_topics, mod_seeds)
```

## Arguments

`mods`                Fitted models returned from [LDA](#).  
`mod_topics`        Vector of integer values corresponding to the number of topics in each model.  
`mod_seeds`         Vector of integer values corresponding to the seed used for each model.

## Value

`lis` (class: LDA\_set) of LDA models (class: LDA\_VEM).

## Examples

```
data(rodents)
document_term_table <- rodents$document_term_table
topics <- 2
nseeds <- 2
control <- LDA_set_control()
mod_topics <- rep(topics, each = length(seq(2, nseeds * 2, 2)))
iseed <- control$iseed
mod_seeds <- rep(seq(iseed, iseed + (nseeds - 1)* 2, 2), length(topics))
nmods <- length(mod_topics)
mods <- vector("list", length = nmods)
for (i in 1:nmods){
  LDA_msg(mod_topics[i], mod_seeds[i], control)
  control_i <- prep_LDA_control(seed = mod_seeds[i], control = control)
  mods[[i]] <- topicmodels::LDA(document_term_table, k = mod_topics[i],
                              control = control_i)
}
package_LDA_set(mods, mod_topics, mod_seeds)
```



---

package\_LDA\_TS      *Package the output of LDA\_TS*

---

## Description

Combine the objects returned by [LDA\\_set](#), [select\\_LDA](#), [TS\\_on\\_LDA](#), and [select\\_TS](#), name them as elements of the list, and set the class of the list as LDA\_TS, for the return from [LDA\\_TS](#).

## Usage

```
package_LDA_TS(LDAs, sel_LDA, TSs, sel_TSs)
```

## Arguments

LDAs	List (class: LDA_set) of LDA models (class: LDA), as returned by <a href="#">LDA_set</a> .
sel_LDA	A reduced version of LDAs that only includes the LDA model(s) selected by <a href="#">select_LDA</a> . Still should be of class LDA_set.
TSs	Class TS_on_LDA list of results from <a href="#">TS</a> applied for each model on each LDA model input, as returned by <a href="#">TS_on_LDA</a> .
sel_TSs	A reduced version of TSs (of class TS_fit) that only includes the TS model chosen via <a href="#">select_TS</a> .

## Value

Class LDA\_TS-class object including all fitted models and selected models specifically, ready to be returned from [LDA\\_TS](#).

## Examples

```
data(rodents)
data <- rodents
control <- LDA_TS_control()
dtt <- data$document_term_table
dct <- data$document_covariate_table
weights <- document_weights(dtt)
LDAs <- LDA_set(dtt, 2, 1, control$LDA_set_control)
sel_LDA <- select_LDA(LDAs, control$LDA_set_control)
TSs <- TS_on_LDA(sel_LDA, dct, ~1, 1, "newmoon", weights,
                 control$TS_control)
sel_TSs <- select_TS(TSs, control$TS_control)
package_LDA_TS(LDAs, sel_LDA, TSs, sel_TSs)
```

---

package_TS	<i>Summarize the Time Series model</i>
------------	--

---

### Description

Calculate relevant summaries for the run of a Time Series model within `TS` and package the output as a `TS_fit`-class object.

### Usage

```
package_TS(data, formula, timename, weights, control, rho_dist, eta_dist)
```

### Arguments

data	data.frame including [1] the time variable (indicated in <code>timename</code> ), [2] the predictor variables (required by <code>formula</code> ) and [3], the multinomial response variable (indicated in <code>formula</code> ) as verified by <code>check_timename</code> and <code>check_formula</code> . Note that the response variables should be formatted as a data.frame object named as indicated by the response entry in the control list, such as <code>gamma</code> for a standard TS analysis on LDA output.
formula	<code>formula</code> defining the regression between relationship the change points. Any predictor variable included must also be a column in <code>data</code> and any (multinomial) response variable must be a set of columns in <code>data</code> , as verified by <code>check_formula</code> .
timename	character element indicating the time variable used in the time series.
weights	Optional class numeric vector of weights for each document. Defaults to NULL, translating to an equal weight for each document. When using <code>multinom_TS</code> in a standard LDATS analysis, it is advisable to weight the documents by their total size, as the result of LDA is a matrix of proportions, which does not account for size differences among documents. For most models, a scaling of the weights (so that the average is 1) is most appropriate, and this is accomplished using <code>document_weights</code> .
control	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <code>TS_control</code> .
rho_dist	List of saved data objects from the ptMCMC estimation of change point locations returned by <code>est_changepoints</code> (unless <code>nchangepoints</code> is 0, then NULL).
eta_dist	Matrix of draws (rows) from the marginal posteriors of the coefficients across the segments (columns), as estimated by <code>est_regressors</code> .

### Value

`TS_fit`-class list containing the following elements, many of which are hidden for printing, but are accessible:

**data** data input to the function.

- formula** [formula](#) input to the function.
- nchangepoints** `nchangepoints` input to the function.
- weights** `weights` input to the function.
- timename** `timename` input to the function.
- control** `control` input to the function.
- lls** Iteration-by-iteration [logLik](#) values for the full time series fit by [multinom\\_TS](#).
- rhos** Iteration-by-iteration change point estimates from [est\\_changepoints](#).
- etas** Iteration-by-iteration marginal regressor estimates from [est\\_regressors](#), which have been unconditioned with respect to the change point locations.
- ptMCMC\_diagnostics** ptMCMC diagnostics, see [diagnose\\_ptMCMC](#)
- rho\_summary** Summary table describing rhos (the change point locations), see [summarize\\_rhos](#).
- rho\_vcov** Variance-covariance matrix for the estimates of rhos (the change point locations), see [measure\\_rho\\_vcov](#).
- eta\_summary** Summary table describing ets (the regressors), see [summarize\\_etas](#).
- eta\_vcov** Variance-covariance matrix for the estimates of etas (the regressors), see [measure\\_eta\\_vcov](#).
- logLik** Across-iteration average of log-likelihoods (`lls`).
- nparams** Total number of parameters in the full model, including the change point locations and regressors.
- AIC** Penalized negative log-likelihood, based on `logLik` and `nparams`.

## Examples

```

data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
formula <- gamma ~ 1
nchangepoints <- 1
control <- TS_control()
data <- data[order(data[, "newmoon"]), ]
rho_dist <- est_changepoints(data, formula, nchangepoints, "newmoon",
                             weights, control)
eta_dist <- est_regressors(rho_dist, data, formula, "newmoon", weights,
                           control)
package_TS(data, formula, "newmoon", weights, control, rho_dist,
            eta_dist)

```

---

package\_TS\_on\_LDA      *Package the output of TS\_on\_LDA*

---

### Description

Set the class and name the elements of the results list returned from applying [TS](#) to the combination of TS models requested for the LDA model(s) input.

### Usage

```
package_TS_on_LDA(TSmods, LDA_models, models)
```

### Arguments

TSmods	list of results from <a href="#">TS</a> applied for each model on each LDA model input.
LDA_models	List of LDA models (class <code>LDA_set</code> , produced by <a href="#">LDA_set</a> ) or a singular LDA model (class <code>LDA</code> , produced by <a href="#">LDA</a> ).
models	<code>data.frame</code> object returned from <a href="#">expand_TS</a> that contains the combinations of LDA models, and formulas and nchangepts used in the TS models.

### Value

Class `TS_on_LDA` list of results from [TS](#) applied for each model on each LDA model input.

### Examples

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDAs <- LDA_set(document_term_table, topics = 2:3, nseeds = 2)
LDA_models <- select_LDA(LDAs)
weights <- document_weights(document_term_table)
mods <- expand_TS(LDA_models, c(~ 1, ~ newmoon), 0:1)
nmods <- nrow(mods)
TSmods <- vector("list", nmods)
for(i in 1:nmods){
  formula_i <- mods$formula[[i]]
  nchangepts_i <- mods$nchangepts[i]
  data_i <- prep_TS_data(document_covariate_table, LDA_models, mods, i)
  TSmods[[i]] <- TS(data_i, formula_i, nchangepts_i, "newmoon",
                    weights, TS_control())
}
package_TS_on_LDA(TSmods, LDA_models, mods)
```

---

plot.LDA_set	<i>Plot a set of LDATS LDA models</i>
--------------	---------------------------------------

---

**Description**

Generalization of the [plot](#) function to work on a list of LDA topic models (class LDA\_set).

**Usage**

```
## S3 method for class 'LDA_set'  
plot(x, ...)
```

**Arguments**

x	An LDA_set object of LDA topic models.
...	Additional arguments to be passed to subfunctions.

**Value**

NULL.

**Examples**

```
data(rodents)  
lda_data <- rodents$document_term_table  
r_LDA <- LDA_set(lda_data, topics = 2, nseeds = 2)  
plot(r_LDA)
```

---

plot.LDA_TS	<i>Plot the key results from a full LDATS analysis</i>
-------------	--

---

**Description**

Generalization of the [plot](#) function to work on fitted LDA\_TS model objects (class LDA\_TS) returned by [LDA\\_TS](#).

**Usage**

```
## S3 method for class 'LDA_TS'  
plot(  
  x,  
  ...,  
  cols = set_LDA_TS_plot_cols(),  
  bin_width = 1,  
)
```

```

  xname = NULL,
  border = NA,
  selection = "median"
)

```

### Arguments

x	A LDA_TS object of a full LDATS model fit by <a href="#">LDA_TS</a> .
...	Additional arguments to be passed to subfunctions. Not currently used, just retained for alignment with plot.
cols	list of elements used to define the colors for the two panels of the summary plot, as generated simply using <a href="#">set_LDA_TS_plot_cols</a> . cols has two elements: LDA and TS, each corresponding the set of plots for its stage in the full model. LDA contains entries cols and option (see <a href="#">set_LDA_plot_colors</a> ). TS contains two entries, rho and gamma, each corresponding to the related panel, and each containing default values for entries named cols, option, and alpha (see <a href="#">set_TS_summary_plot_cols</a> , <a href="#">set_gamma_colors</a> , and <a href="#">set_rho_hist_colors</a> ).
bin_width	Width of the bins used in the histograms of the summary time series plot, in units of the time variable used to fit the model (the x-axis).
xname	Label for the x-axis in the summary time series plot. Defaults to NULL, which results in usage of the timename element of the control list (held in control\$TS_control\$timename). To have no label printed, set xname = "".
border	Border for the histogram, default is NA.
selection	Indicator of the change points to use in the time series summary plot. Currently only defined for "median" and "mode".

### Value

NULL.

### Examples

```

data(rodents)
mod <- LDA_TS(data = rodents, topics = 2, nseeds = 1, formulas = ~1,
              nchangepts = 1, timename = "newmoon")
plot(mod, binwidth = 5, xlab = "New moon")

```

## Description

Create an LDATS LDA summary plot, with a top panel showing the topic proportions for each word and a bottom panel showing the topic proportions of each document/over time. The plot function is defined for class LDA\_VEM specifically (see [LDA](#)).

LDA\_plot\_top\_panel creates an LDATS LDA summary plot top panel showing the topic proportions word-by-word.

LDA\_plot\_bottom\_panel creates an LDATS LDA summary plot bottom panel showing the topic proportions over time/documents.

## Usage

```
## S3 method for class 'LDA_VEM'
plot(
  x,
  ...,
  xtime = NULL,
  xname = NULL,
  cols = NULL,
  option = "C",
  alpha = 0.8,
  LDATS = FALSE
)

LDA_plot_top_panel(
  x,
  cols = NULL,
  option = "C",
  alpha = 0.8,
  together = FALSE,
  LDATS = FALSE
)

LDA_plot_bottom_panel(
  x,
  xtime = NULL,
  xname = NULL,
  cols = NULL,
  option = "C",
  alpha = 0.8,
  together = FALSE,
  LDATS = FALSE
)
```

## Arguments

x                    Object of class LDA\_VEM.

...	Not used, retained for alignment with base function.
xtime	Optional x values used to plot the topic proportions according to a specific time value (rather than simply the order of observations).
xname	Optional name for the x values used in plotting the topic proportions (otherwise defaults to "Document").
cols	Colors to be used to plot the topics. Any valid color values ( <i>e.g.</i> , see <a href="#">colors</a> , <a href="#">rgb</a> ) can be input as with a standard plot. The default (cols = NULL) triggers use of <a href="#">viridis</a> color options (see option).
option	A character string indicating the color option from <a href="#">viridis</a> to use if 'cols == NULL'. Four options are available: "magma" (or "A"), "inferno" (or "B"), "plasma" (or "C", the default option), "viridis" (or "D") and "cividis" (or "E").
alpha	Numeric value [0,1] that indicates the transparency of the colors used. Supported only on some devices, see <a href="#">rgb</a> .
LDATS	logical indicating if the LDA plot is part of a larger LDATS plot output.
together	logical indicating if the subplots are part of a larger LDA plot output.

### Value

NULL.

### Examples

```
data(rodents)
lda_data <- rodents$document_term_table
r_LDA <- LDA_set(lda_data, topics = 4, nseeds = 10)
best_lda <- select_LDA(r_LDA)[[1]]
plot(best_lda, option = "cividis")
LDA_plot_top_panel(best_lda, option = "cividis")
LDA_plot_bottom_panel(best_lda, option = "cividis")
```

---

plot.TS\_fit

*Plot an LDATS TS model*

---

### Description

Generalization of the [plot](#) function to work on fitted TS model objects (class `TS_fit`) returned from [TS](#).



**Usage**

```
## S3 method for class 'TS_fit'
plot(
  x,
  ...,
  plot_type = "summary",
  interactive = FALSE,
  cols = set_TS_summary_plot_cols(),
  bin_width = 1,
  xname = NULL,
  border = NA,
  selection = "median",
  LDATS = FALSE
)
```

**Arguments**

x	A TS_fit object of a multinomial time series model fit by <a href="#">TS</a> .
...	Additional arguments to be passed to subfunctions. Not currently used, just retained for alignment with <a href="#">plot</a> .
plot_type	"diagnostic" or "summary".
interactive	logical input, should be codeTRUE unless testing.
cols	list of elements used to define the colors for the two panels of the summary plot, as generated simply using <a href="#">set_TS_summary_plot_cols</a> . cols has two elements rho and gamma, each corresponding to the related panel, and each containing default values for entries named cols, option, and alpha. See <a href="#">set_gamma_colors</a> and <a href="#">set_rho_hist_colors</a> for details on usage.
bin_width	Width of the bins used in the histograms of the summary time series plot, in units of the x-axis (the time variable used to fit the model).
xname	Label for the x-axis in the summary time series plot. Defaults to NULL, which results in usage of the <code>timename</code> element of the control list (held in <code>control\$TS_control\$timename</code> ). To have no label printed, set <code>xname = ""</code> .
border	Border for the histogram, default is NA.
selection	Indicator of the change points to use in the time series summary plot. Currently only defined for "median" and "mode".
LDATS	logical indicating if the plot is part of a larger LDATS plot output.

**Value**

NULL.

**Examples**

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
```

```
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
TSmod <- TS(data, gamma ~ 1, nchangepts = 1, "newmoon", weights)
plot(TSmod)
```

---

posterior_plot	<i>Produce the posterior distribution histogram panel for the TS diagnostic plot of a parameter</i>
----------------	---

---

### Description

Produce a vanilla histogram plot using `hist` for the parameter of interest (rho or eta) as part of [TS\\_diagnostics\\_plot](#). A vertical line is added to show the median of the posterior.

### Usage

```
posterior_plot(x, xlab = "parameter value")
```

### Arguments

x	Vector of parameter values drawn from the posterior distribution, indexed to the iteration by the order of the vector.
xlab	character value used to label the x axis.

### Value

NULL.

### Examples

```
posterior_plot(rnorm(100, 0, 1))
```

---

prep_chunks	<i>Prepare the time chunk table for a multinomial change point Time Series model</i>
-------------	--

---

## Description

Creates the table containing the start and end times for each chunk within a time series, based on the change points (used to break up the time series) and the range of the time series. If there are no change points (i.e. `changepoints` is `NULL`), there is still a single chunk defined by the start and end of the time series.

## Usage

```
prep_chunks(data, changepoints = NULL, timename = "time")
```

## Arguments

<code>data</code>	Class <code>data.frame</code> object including the predictor and response variables, but specifically here containing the column indicated by the <code>timename</code> input.
<code>changepoints</code>	Numeric vector indicating locations of the change points. Must be conformable to integer values.
<code>timename</code>	character element indicating the time variable used in the time series. Defaults to "time". The variable must be integer-conformable or a Date. If the variable named is a Date, the input is converted to an integer, resulting in the timestep being 1 day, which is often not desired behavior.

## Value

`data.frame` of start and end times (columns) for each chunk (rows).

## Examples

```
data(rodents)
dtt <- rodents$document_term_table
lda <- LDA_set(dtt, 2, 1, list(quiet = TRUE))
dct <- rodents$document_covariate_table
dct$gamma <- lda[[1]]@gamma
chunks <- prep_chunks(dct, changepoints = 100, timename = "newmoon")
```

---

prep_cpts	<i>Initialize and update the change point matrix used in the ptMCMC algorithm</i>
-----------	---

---

### Description

Each of the chains is initialized by prep\_cpts using a draw from the available times (i.e. assuming a uniform prior), the best fit (by likelihood) draw is put in the focal chain with each subsequently worse fit placed into the subsequently hotter chain. update\_cpts updates the change points after every iteration in the ptMCMC algorithm.

### Usage

```
prep_cpts(data, formula, nchangepts, timename, weights, control = list())
```

```
update_cpts(cpts, swaps)
```

### Arguments

data	data.frame including [1] the time variable (indicated in timename), [2] the predictor variables (required by formula) and [3], the multinomial response variable (indicated in formula) as verified by <a href="#">check_timename</a> and <a href="#">check_formula</a> . Note that the response variables should be formatted as a data.frame object named as indicated by the response entry in the control list, such as gamma for a standard TS analysis on LDA output.
formula	formula defining the regression relationship between the change points, see <a href="#">formula</a> . Any predictor variable included must also be a column in data and any (multinomial) response variable must be a set of columns in data, as verified by <a href="#">check_formula</a> .
nchangepts	integer corresponding to the number of change points to include in the model. 0 is a valid input (corresponding to no change points, so a singular time series model), and the current implementation can reasonably include up to 6 change points. The number of change points is used to dictate the segmentation of the data for each continuous model and each LDA model.
timename	character element indicating the time variable used in the time series. Defaults to "time". The variable must be integer-conformable or a Date. If the variable named is a Date, the input is converted to an integer, resulting in the timestep being 1 day, which is often not desired behavior.
weights	Optional class numeric vector of weights for each document. Defaults to NULL, translating to an equal weight for each document. When using multinom_TS in a standard LDATS analysis, it is advisable to weight the documents by their total size, as the result of LDA is a matrix of proportions, which does not account for size differences among documents. For most models, a scaling of the weights (so that the average is 1) is most appropriate, and this is accomplished using <a href="#">document_weights</a> .

control	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <a href="#">TS_control</a> .
cpts	The existing matrix of change points.
swaps	Chain configuration after among-temperature swaps.

### Value

list of [1] matrix of change points (rows) for each temperature (columns) and [2] vector of log-likelihood values for each of the chains.

### Examples

```

data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
data <- data[order(data[, "newmoon"]), ]
saves <- prep_saves(1, TS_control())
inputs <- prep_ptMCMC_inputs(data, gamma ~ 1, 1, "newmoon", weights,
                             TS_control())
cpts <- prep_cpts(data, gamma ~ 1, 1, "newmoon", weights, TS_control())
ids <- prep_ids(TS_control())
for(i in 1:TS_control()$nit){
  steps <- step_chains(i, cpts, inputs)
  swaps <- swap_chains(steps, inputs, ids)
  saves <- update_saves(i, saves, steps, swaps)
  cpts <- update_cpts(cpts, swaps)
  ids <- update_ids(ids, swaps)
}

```

---

```
prep_ids
```

*Initialize and update the chain ids throughout the ptMCMC algorithm*

---

### Description

`prep_ids` creates and `update_ids` updates the active vector of identities (`ids`) for each of the chains in the ptMCMC algorithm. These `ids` are used to track trips of the particles among chains.

These functions were designed to work within `TS` and specifically `est_changepoints`, but have been generalized and would work within any general ptMCMC as long as `control`, `ids`, and `swaps` are formatted properly.

**Usage**

```
prep_ids(control = list())

update_ids(ids, swaps)
```

**Arguments**

control	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <a href="#">TS_control</a> .
ids	The existing vector of chain ids.
swaps	Chain configuration after among-temperature swaps.

**Value**

The vector of chain ids.

**Examples**

```
prep_ids()

data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
data <- data[order(data[, "newmoon"]), ]
saves <- prep_saves(1, TS_control())
inputs <- prep_ptMCMC_inputs(data, gamma ~ 1, 1, "newmoon", weights,
                           TS_control())
cpts <- prep_cpts(data, gamma ~ 1, 1, "newmoon", weights, TS_control())
ids <- prep_ids(TS_control())
for(i in 1:TS_control()$nit){
  steps <- step_chains(i, cpts, inputs)
  swaps <- swap_chains(steps, inputs, ids)
  saves <- update_saves(i, saves, steps, swaps)
  cpts <- update_cpts(cpts, swaps)
  ids <- update_ids(ids, swaps)
}
```

---

prep_LDA_control	<i>Set the control inputs to include the seed</i>
------------------	---

---

**Description**

Update the control list for the LDA model with the specific seed as indicated. And remove controls not used within the LDA itself.

**Usage**

```
prep_LDA_control(seed, control = list())
```

**Arguments**

seed	integer used to set the seed of the specific model.
control	Named list of control parameters to be used in <a href="#">LDA</a> Note that if control has an element named seed it will be overwritten by the seed argument of prep_LDA_control.

**Value**

list of controls to be used in the LDA.

**Examples**

```
prep_LDA_control(seed = 1)
```

---

prep_pbar	<i>Initialize and tick through the progress bar</i>
-----------	---

---

**Description**

prep\_pbar creates and update\_pbar steps through the progress bars (if desired) in [TS](#)

**Usage**

```
prep_pbar(control = list(), bar_type = "rho", nr = NULL)
```

```
update_pbar(pbar, control = list())
```

**Arguments**

control	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <code>TS_control</code> . Of use here is <code>quiet</code> which is a logical indicator of whether there should be information (i.e. the progress bar) printed during the run or not. Default is TRUE.
bar_type	"rho" (for change point locations) or "eta" (for regressors).
nr	integer number of unique realizations, needed when <code>bar_type = "eta"</code> .
pbar	The progress bar object returned from <code>prep_pbar</code> .

**Value**

`prep_pbar`: the initialized progress bar object.

`update_pbar`: the ticked-forward pbar.

**Examples**

```
pb <- prep_pbar(control = list(nit = 2)); pb
pb <- update_pbar(pb); pb
pb <- update_pbar(pb); pb
```

---

<code>prep_proposal_dist</code>	<i>Pre-calculate the change point proposal distribution for the ptMCMC algorithm</i>
---------------------------------	--

---

**Description**

Calculate the proposal distribution in advance of actually running the ptMCMC algorithm in order to decrease computation time. The proposal distribution is a joint of three distributions: [1] a multinomial distribution selecting among the change points within the chain, [2] a binomial distribution selecting the direction of the step of the change point (earlier or later in the time series), and [3] a geometric distribution selecting the magnitude of the step.

**Usage**

```
prep_proposal_dist(nchangepts, control = list())
```

**Arguments**

nchangepts	Integer corresponding to the number of change points to include in the model. 0 is a valid input (corresponding to no change points, so a singular time series model), and the current implementation can reasonably include up to 6 change points. The number of change points is used to dictate the segmentation of the data for each continuous model and each LDA model.
------------	---



**control** A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by `TS_control`. Currently relevant here is `magnitude`, which controls the magnitude of the step size (is the average of the geometric distribution).

### Value

list of two matrix elements: [1] the size of the proposed step for each iteration of each chain and [2] the identity of the change point location to be shifted by the step for each iteration of each chain.

### Examples

```
prep_proposal_dist(nchangepoints = 2)
```

---

prep\_ptMCMC\_inputs      *Prepare the inputs for the ptMCMC algorithm estimation of change points*

---

### Description

Package the static inputs (controls and data structures) used by the ptMCMC algorithm in the context of estimating change points.

This function was designed to work within `TS` and specifically `est_changepoints`. It is still hard-coded to do so, but has the capacity to be generalized to work with any estimation via ptMCMC with additional coding work.

### Usage

```
prep_ptMCMC_inputs(  
  data,  
  formula,  
  nchangepoints,  
  timename,  
  weights = NULL,  
  control = list()  
)
```

### Arguments

**data** Class data.frame object including [1] the time variable (indicated in `control`), [2] the predictor variables (required by `formula`) and [3], the multinomial response variable (indicated in `formula`).

**formula** formula describing the continuous change. Any predictor variable included must also be a column in the data. Any (multinomial) response variable must also be a set of columns in data.

nchangepts	Integer corresponding to the number of change points to include in the model. 0 is a valid input (corresponding to no change points, so a singular time series model), and the current implementation can reasonably include up to 6 change points. The number of change points is used to dictate the segmentation of the data for each continuous model and each LDA model.
timename	character element indicating the time variable used in the time series. Defaults to "time". The variable must be integer-conformable or a Date. If the variable named is a Date, the input is converted to an integer, resulting in the timestep being 1 day, which is often not desired behavior.
weights	Optional class numeric vector of weights for each document. Defaults to NULL, translating to an equal weight for each document. When using <code>multinom_TS</code> in a standard LDATS analysis, it is advisable to weight the documents by their total size, as the result of LDA is a matrix of proportions, which does not account for size differences among documents. For most models, a scaling of the weights (so that the average is 1) is most appropriate, and this is accomplished using <code>document_weights</code> .
control	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <code>TS_control</code> .

### Value

Class `ptMCMC_inputs` list, containing the static inputs for use within the `ptMCMC` algorithm for estimating change points.

### Examples

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
data <- data[order(data[, "newmoon"]), ]
saves <- prep_saves(1, TS_control())
inputs <- prep_ptMCMC_inputs(data, gamma ~ 1, 1, "newmoon", weights,
                           TS_control())
```

**Description**

prep\_saves creates the data structure used to save the output from each iteration of the ptMCMC algorithm, which is added via update\_saves. Once the ptMCMC is complete, the saved data objects are then processed (burn-in iterations are dropped and the remaining iterations are thinned) via process\_saves.

This set of functions was designed to work within [TS](#) and specifically [est\\_changepoints](#). They are still hardcoded to do so, but have the capacity to be generalized to work with any estimation via ptMCMC with additional coding work.

**Usage**

```
prep_saves(nchangepoints, control = list())
```

```
update_saves(i, saves, steps, swaps)
```

```
process_saves(saves, control = list())
```

**Arguments**

- |               |   |
|---------------|---|
| nchangepoints | integer corresponding to the number of change points to include in the model. 0 is a valid input (corresponding to no change points, so a singular time series model), and the current implementation can reasonably include up to 6 change points. The number of change points is used to dictate the segmentation of the data for each continuous model and each LDA model. |
| control       | A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <a href="#">TS_control</a> .  |
| i             | integer iteration index.  |
| saves         | The existing list of saved data objects.  |
| steps         | Chain configuration after within-temperature steps.   |
| swaps         | Chain configuration after among-temperature swaps.  |

**Value**

list of ptMCMC objects: change points (`$cpts`), log-likelihoods (`$lls`), chain ids (`$ids`), step acceptances (`$step_accepts`), and swap acceptances (`$swap_accepts`).

**Examples**

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
data <- data[order(data[, "newmoon"]), ]
```

```

saves <- prep_saves(1, TS_control())
inputs <- prep_ptMCMC_inputs(data, gamma ~ 1, 1, "newmoon", weights,
                           TS_control())
cpts <- prep_cpts(data, gamma ~ 1, 1, "newmoon", weights, TS_control())
ids <- prep_ids(TS_control())
for(i in 1:TS_control()$nit){
  steps <- step_chains(i, cpts, inputs)
  swaps <- swap_chains(steps, inputs, ids)
  saves <- update_saves(i, saves, steps, swaps)
  cpts <- update_cpts(cpts, swaps)
  ids <- update_ids(ids, swaps)
}
process_saves(saves, TS_control())

```

---

```
prep_temp_sequence      Prepare the ptMCMC temperature sequence
```

---

### Description

Create the series of temperatures used in the ptMCMC algorithm.

This function was designed to work within [TS](#) and [est\\_changepoints](#) specifically, but has been generalized and would work with any ptMCMC model as long as `control` includes the relevant control parameters (and provided that the [check\\_control](#) function and its use here are generalized).

### Usage

```
prep_temp_sequence(control = list())
```

### Arguments

<code>control</code>	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <a href="#">TS_control</a> .
----------------------	--

### Value

vector of temperatures.

### Examples

```
prep_temp_sequence()
```

---

prep_TS_data	<i>Prepare the model-specific data to be used in the TS analysis of LDA output</i>
--------------	--

---

### Description

Append the estimated topic proportions from a fitted LDA model to the document covariate table to create the data structure needed for [TS](#).

### Usage

```
prep_TS_data(document_covariate_table, LDA_models, mods, i = 1)
```

### Arguments

document_covariate_table	Document covariate table (rows: documents, columns: time index and covariate options). Every model needs a covariate to describe the time value for each document (in whatever units and whose name in the table is input in <code>timename</code> ) that dictates the application of the change points. In addition, all covariates named within specific models in <code>formula</code> must be included. Must be a conformable to a data table, as verified by <a href="#">check_document_covariate_table</a> .
LDA_models	List of LDA models (class <code>LDA_set</code> , produced by <a href="#">LDA_set</a> ) or a singular LDA model (class <code>LDA</code> , produced by <a href="#">LDA</a> ).
mods	The <code>data.table</code> created by <a href="#">expand_TS</a> that contains each of the models (defined by the LDA model to use and the <code>and</code> formula number of changepoints for the TS model). Indexed here by <code>i</code> .
i	integer index referencing the row in <code>mods</code> to use.

### Value

Class `data.frame` object including [1] the time variable (indicated in `control`), [2] the predictor variables (required by `formula`) and [3], the multinomial response variable (indicated in `formula`), ready for input into [TS](#).

### Examples

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDAs <- LDA_set(document_term_table, topics = 2:3, nseeds = 2)
LDA_models <- select_LDA(LDAs)
weights <- document_weights(document_term_table)
formulas <- c(~ 1, ~ newmoon)
mods <- expand_TS(LDA_models, formulas = ~1, nchangepoints = 0)
data1 <- prep_TS_data(document_covariate_table, LDA_models, mods)
```

---

print.LDA_TS	<i>Print the selected LDA and TS models of LDA_TS object</i>
--------------	--

---

**Description**

Convenience function to print only the selected elements of a LDA\_TS-class object returned by [LDA\\_TS](#)

**Usage**

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

**Arguments**

x	Class LDA_TS object to be printed.
...	Not used, simply included to maintain method compatibility.

**Value**

The selected models in x as a two-element list with the TS component only returning the non-hidden components.

**Examples**

```
data(rodents)  
mod <- LDA_TS(data = rodents, topics = 2, nseeds = 1, formulas = ~1,  
              nchangepts = 1, timename = "newmoon")  
print(mod)
```

---

print.TS_fit	<i>Print a Time Series model fit</i>
--------------	--------------------------------------

---

**Description**

Convenience function to print only the most important components of a TS\_fit-class object fit by [TS](#).

**Usage**

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

**Arguments**

x                    Class TS\_fit object to be printed.  
 ...                  Not used, simply included to maintain method compatibility.

**Value**

The non-hidden parts of x as a list.

**Examples**

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
TSmod <- TS(data, gamma ~ 1, nchangepts = 1, "newmoon", weights)
print(TSmod)
```

---

print.TS\_on\_LDA            *Print a set of Time Series models fit to LDAs*

---

**Description**

Convenience function to print only the names of a TS\_on\_LDA-class object generated by [TS\\_on\\_LDA](#).

**Usage**

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

**Arguments**

x                    Class TS\_on\_LDA object to be printed.  
 ...                  Not used, simply included to maintain method compatibility.

**Value**

character vector of the names of x's models.

**Examples**

```

data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDAs <- LDA_set(document_term_table, topics = 2:3, nseeds = 2)
LDA_models <- select_LDA(LDAs)
weights <- document_weights(document_term_table)
formulas <- c(~ 1, ~ newmoon)
mods <- TS_on_LDA(LDA_models, document_covariate_table, formulas,
                 nchangepts = 0:1, timename = "newmoon", weights)
print(mods)

```

---

```
print_model_run_message
```

*Print the message to the console about which combination of the Time Series and LDA models is being run*

---

**Description**

If desired, print a message at the beginning of every model combination stating the TS model and the LDA model being evaluated.

**Usage**

```
print_model_run_message(models, i, LDA_models, control)
```

**Arguments**

models	data.frame object returned from <a href="#">expand_TS</a> that contains the combinations of LDA models, and formulas and nchangepts used in the TS models.
i	integer index of the row to use from models.
LDA_models	List of LDA models (class LDA_set, produced by <a href="#">LDA_set</a> ) or a singular LDA model (class LDA, produced by <a href="#">LDA</a> ).
control	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <a href="#">TS_control</a> . Of particular importance here is the logical-class element named quiet.

**Value**

NULL.



**Examples**

```

data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDAs <- LDA_set(document_term_table, topics = 2:3, nseeds = 2)
LDA_models <- select_LDA(LDAs)
weights <- document_weights(document_term_table)
formulas <- c(~ 1, ~ newmoon)
nchangepts <- 0:1
mods <- expand_TS(LDA_models, formulas, nchangepts)
print_model_run_message(mods, 1, LDA_models, TS_control())

```

---

proposed_step_mods	<i>Fit the chunk-level models to a time series, given a set of proposed change points within the ptMCMC algorithm</i>
--------------------	---

---

**Description**

This function wraps around `TS_memo` (optionally memoised `multinom_TS`) to provide a simpler interface within the `ptMCMC` algorithm and is implemented within `propose_step`.

**Usage**

```
proposed_step_mods(prop_changepts, inputs)
```

**Arguments**

`prop_changepts` matrix of proposed change points across chains.

`inputs` Class `ptMCMC_inputs` list, containing the static inputs for use within the `ptMCMC` algorithm.

**Value**

List of models associated with the proposed step, with an element for each chain.

**Examples**

```

data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
data <- data[order(data[, "newmoon"]), ]
saves <- prep_saves(1, TS_control())
inputs <- prep_ptMCMC_inputs(data, gamma ~ 1, 1, "newmoon", weights,

```

```

                                TS_control())
cpts <- prep_cpts(data, gamma ~ 1, 1, "newmoon", weights, TS_control())
i <- 1
pdist <- inputs$pdist
ntemps <- length(inputs$temps)
selection <- cbind(pdist$which_steps[i, ], 1:ntemps)
prop_changepts <- cpts$changepts
curr_changepts_s <- cpts$changepts[selection]
prop_changepts_s <- curr_changepts_s + pdist$steps[i, ]
if(all(is.na(prop_changepts_s))){
  prop_changepts_s <- NULL
}
prop_changepts[selection] <- prop_changepts_s
mods <- proposed_step_mods(prop_changepts, inputs)

```

---

rho\_lines

*Add change point location lines to the time series plot*


---

### Description

Adds vertical lines to the plot of the time series of fitted proportions associated with the change points of interest.

### Usage

```
rho_lines(spec_rhos)
```

### Arguments

spec\_rhos      numeric vector indicating the locations along the x axis where the specific change points being used are located.

### Examples

```

data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
TSmod <- TS(data, gamma ~ 1, nchangepts = 1, "newmoon", weights)
pred_gamma_TS_plot(TSmod)
rho_lines(200)

```

---

rodents	<i>Portal rodent data</i>
---------	---------------------------

---

**Description**

An example LDATS dataset, functionally that used in Christensen *et al.* (2018). The data are counts of 21 rodent species across 436 sampling events, with the count being the total number observed across 8 50 m x 50 m plots, each sampled using 49 live traps (Brown 1998, Ernest *et al.* 2016).

**Usage**

```
rodents
```

**Format**

A list of two data.frame-class objects with rows corresponding to documents (sampling events). One element is the document term table (called `document_term_table`), which contains counts of the species (terms) in each sample (document), and the other is the document covariate table (called `document_covariate_table`) with columns of covariates (newmoon number, sin and cos of the fraction of the year).

**Source**

<https://github.com/weecology/PortalData/tree/master/Rodents>

**References**

- Brown, J. H. 1998. The desert granivory experiments at Portal. Pages 71-95 in W. J. Resetarits Jr. and J. Bernardo, *editors*, *Experimental Ecology*. Oxford University Press, New York, New York, USA.
- Christensen, E., D. J. Harris, and S. K. M. Ernest. 2018. Long-term community change through multiple rapid transitions in a desert rodent community. *Ecology* **99**:1523-1529. [link](#).
- Ernest, S. K. M., *et al.* 2016. Long-term monitoring and experimental manipulation of a Chihuahuan desert ecosystem near Portal, Arizona (1977-2013). *Ecology* **97**:1082. [link](#).

---

select_LDA	<i>Select the best LDA model(s) for use in time series</i>
------------	--

---

**Description**

Select the best model(s) of interest from an `LDA_set` object, based on a set of user-provided functions. The functions default to choosing the model with the lowest AIC value.

**Usage**

```
select_LDA(LDA_models = NULL, control = list())
```

**Arguments**

LDA_models	An object of class LDA_set produced by <a href="#">LDA_set</a> .
control	A list of parameters to control the running and selecting of LDA models. Values not input assume default values set by <a href="#">LDA_set_control</a> . Values for running the LDAs replace defaults in (LDAcontrol, see <a href="#">LDA</a> (but if seed is given, it will be overwritten; use i seed instead).

**Value**

A reduced version of LDA\_models that only includes the selected LDA model(s). The returned object is still an object of class LDA\_set.

**Examples**

```
data(rodents)
lda_data <- rodents$document_term_table
r_LDA <- LDA_set(lda_data, topics = 2, nseeds = 2)
select_LDA(r_LDA)
```

---

select\_TS

*Select the best Time Series model*


---

**Description**

Select the best model of interest from an TS\_on\_LDA object generated by [TS\\_on\\_LDA](#), based on a set of user-provided functions. The functions default to choosing the model with the lowest AIC value.

Presently, the set of functions should result in a singular selected model. If multiple models are chosen via the selection, only the first is returned.

**Usage**

```
select_TS(TS_models, control = list())
```

**Arguments**

TS_models	An object of class TS_on_LDA produced by <a href="#">TS_on_LDA</a> .
control	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <a href="#">TS_control</a> .

**Value**

A reduced version of TS\_models that only includes the selected TS model. The returned object is a single TS model object of class TS\_fit.

**Examples**

```

data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDAs <- LDA_set(document_term_table, topics = 2:3, nseeds = 2)
LDA_models <- select_LDA(LDAs)
weights <- document_weights(document_term_table)
formulas <- c(~ 1, ~ newmoon)
mods <- TS_on_LDA(LDA_models, document_covariate_table, formulas,
                 nchangepoints = 0:1, timename = "newmoon", weights)
select_TS(mods)

```

---

set_gamma_colors	<i>Prepare the colors to be used in the gamma time series</i>
------------------	---

---

**Description**

Based on the inputs, create the set of colors to be used in the time series of the fitted gamma (topic proportion) values.

**Usage**

```
set_gamma_colors(x, cols = NULL, option = "D", alpha = 1)
```

**Arguments**

x	Object of class <code>TS_fit</code> , fit by <a href="#">TS</a> .
cols	Colors to be used to plot the time series of fitted topic proportions.
option	A character string indicating the color option from <a href="#">viridis</a> to use if "cols == NULL". Four options are available: "magma" (or "A"), "inferno" (or "B"), "plasma" (or "C"), "viridis" (or "D", the default option) and "cividis" (or "E").
alpha	Numeric value [0,1] that indicates the transparency of the colors used. Supported only on some devices, see <a href="#">rgb</a> .

**Value**

Vector of character hex codes indicating colors to use.

**Examples**

```

data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma

```

```
weights <- document_weights(document_term_table)
TSmod <- TS(data, gamma ~ 1, nchangepts = 1, "newmoon", weights)
set_gamma_colors(TSmod)
```

---

set\_LDA\_plot\_colors     *Prepare the colors to be used in the LDA plots*

---

### Description

Based on the inputs, create the set of colors to be used in the LDA plots made by `plot.LDA_TS`.

### Usage

```
set_LDA_plot_colors(x, cols = NULL, option = "C", alpha = 0.8)
```

### Arguments

x	Object of class LDA.
cols	Colors to be used to plot the topics. Any valid color values ( <i>e.g.</i> , see <a href="#">colors</a> , <a href="#">rgb</a> ) can be input as with a standard plot. The default (cols = NULL) triggers use of <a href="#">viridis</a> color options (see option).
option	A character string indicating the color option from <a href="#">viridis</a> to use if 'cols == NULL'. Four options are available: "magma" (or "A"), "inferno" (or "B"), "plasma" (or "C", the default option), "viridis" (or "D") and "cividis" (or "E").
alpha	Numeric value [0,1] that indicates the transparency of the colors used. Supported only on some devices, see <a href="#">rgb</a> .

### Value

vector of character hex codes indicating colors to use.

### Examples

```
data(rodents)
lda_data <- rodents$document_term_table
r_LDA <- LDA_set(lda_data, topics = 4, nseeds = 10)
set_LDA_plot_colors(r_LDA[[1]])
```

---

set\_LDA\_TS\_plot\_cols *Create the list of colors for the LDATS summary plot*

---

### Description

A default list generator function that produces the options for the colors controlling the panels of the LDATS summary plots, needed because the change point histogram panel should be in a different color scheme than the LDA and fitted time series model panels, which should be in a matching color scheme. See [set\\_LDA\\_plot\\_colors](#), [set\\_TS\\_summary\\_plot\\_cols](#), [set\\_gamma\\_colors](#), and [set\\_rho\\_hist\\_colors](#) for specific details on usage.

### Usage

```
set_LDA_TS_plot_cols(
  rho_cols = NULL,
  rho_option = "D",
  rho_alpha = 0.4,
  gamma_cols = NULL,
  gamma_option = "C",
  gamma_alpha = 0.8
)
```

### Arguments

rho_cols	Colors to be used to plot the histograms of change points. Any valid color values ( <i>e.g.</i> , see <a href="#">colors</a> , <a href="#">rgb</a> ) can be input as with a standard plot. The default (rho_cols = NULL) triggers use of <a href="#">viridis</a> color options (see rho_option).
rho_option	A character string indicating the color option from <a href="#">viridis</a> to use if 'rho_cols == NULL'. Four options are available: "magma" (or "A"), "inferno" (or "B"), "plasma" (or "C"), "viridis" (or "D", the default option) and "cividis" (or "E").
rho_alpha	Numeric value [0,1] that indicates the transparency of the colors used. Supported only on some devices, see <a href="#">rgb</a> .
gamma_cols	Colors to be used to plot the LDA topic proportions, time series of observed topic proportions, and time series of fitted topic proportions. Any valid color values ( <i>e.g.</i> , see <a href="#">colors</a> , <a href="#">rgb</a> ) can be input as with a standard plot. The default (gamma_cols = NULL) triggers use of <a href="#">viridis</a> color options (see gamma_option).
gamma_option	A character string indicating the color option from <a href="#">viridis</a> to use if gamma_cols == NULL'. Four options are available: "magma" (or "A"), "inferno" (or "B"), "plasma" (or "C", the default option), "viridis" (or "D") and "cividis" (or "E").
gamma_alpha	Numeric value [0,1] that indicates the transparency of the colors used. Supported only on some devices, see <a href="#">rgb</a> .

**Value**

list of elements used to define the colors for the two panels of the summary plot, as generated simply using `set_LDA_TS_plot_cols`. `cols` has two elements: LDA and TS, each corresponding the set of plots for its stage in the full model. LDA contains entries `cols` and `options` (see `set_LDA_plot_colors`). TS contains two entries, `rho` and `gamma`, each corresponding to the related panel, and each containing default values for entries named `cols`, `option`, and `alpha` (see `set_TS_summary_plot_cols`, `set_gamma_colors`, and `set_rho_hist_colors`).

**Examples**

```
set_LDA_TS_plot_cols()
```

---

```
set_rho_hist_colors    Prepare the colors to be used in the change point histogram
```

---

**Description**

Based on the inputs, create the set of colors to be used in the change point histogram.

**Usage**

```
set_rho_hist_colors(x = NULL, cols = NULL, option = "D", alpha = 1)
```

**Arguments**

<code>x</code>	matrix of change point locations (element rhos) from an object of class <code>TS_fit</code> , fit by <code>TS</code> .
<code>cols</code>	Colors to be used to plot the histograms of change points. Any valid color values (e.g., see <code>colors</code> , <code>rgb</code> ) can be input as with a standard plot. The default ( <code>rho_cols = NULL</code> ) triggers use of <code>viridis</code> color options (see <code>rho_option</code> ).
<code>option</code>	A character string indicating the color option from <code>viridis</code> to use if <code>cols == NULL</code> . Four options are available: "magma" (or "A"), "inferno" (or "B"), "plasma" (or "C"), "viridis" (or "D", the default option) and "cividis" (or "E").
<code>alpha</code>	Numeric value [0,1] that indicates the transparency of the colors used. Supported only on some devices, see <code>rgb</code> .

**Value**

Vector of character hex codes indicating colors to use.



**Examples**

```

data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
TSmod <- TS(data, gamma ~ 1, nchangepts = 1, "newmoon", weights)
set_rho_hist_colors(TSmod$rhos)

```

---

```
set_TS_summary_plot_cols
```

*Create the list of colors for the TS summary plot*

---

**Description**

A default list generator function that produces the options for the colors controlling the panels of the TS summary plots, so needed because the panels should be in different color schemes. See [set\\_gamma\\_colors](#) and [set\\_rho\\_hist\\_colors](#) for specific details on usage.

**Usage**

```

set_TS_summary_plot_cols(
  rho_cols = NULL,
  rho_option = "D",
  rho_alpha = 0.4,
  gamma_cols = NULL,
  gamma_option = "C",
  gamma_alpha = 0.8
)

```

**Arguments**

<code>rho_cols</code>	Colors to be used to plot the histograms of change points. Any valid color values (e.g., see <a href="#">colors</a> , <a href="#">rgb</a> ) can be input as with a standard plot. The default ( <code>rho_cols = NULL</code> ) triggers use of <a href="#">viridis</a> color options (see <code>rho_option</code> ).
<code>rho_option</code>	A character string indicating the color option from <a href="#">viridis</a> to use if ' <code>rho_cols == NULL</code> '. Four options are available: "magma" (or "A"), "inferno" (or "B"), "plasma" (or "C"), "viridis" (or "D", the default option) and "cividis" (or "E").
<code>rho_alpha</code>	Numeric value [0,1] that indicates the transparency of the colors used. Supported only on some devices, see <a href="#">rgb</a> .
<code>gamma_cols</code>	Colors to be used to plot the LDA topic proportions, time series of observed topic proportions, and time series of fitted topic proportions. Any valid color values (e.g., see <a href="#">colors</a> , <a href="#">rgb</a> ) can be input as with a standard plot. The default ( <code>gamma_cols = NULL</code> ) triggers use of <a href="#">viridis</a> color options (see <code>gamma_option</code> ).

gamma_option	A character string indicating the color option from <a href="#">viridis</a> to use if gamma_cols == NULL'. Four options are available: "magma" (or "A"), "inferno" (or "B"), "plasma" (or "C"), "viridis" (or "D", the default option) and "cividis" (or "E").
gamma_alpha	Numeric value [0,1] that indicates the transparency of the colors used. Supported only on some devices, see <a href="#">rgb</a> .

### Value

list of elements used to define the colors for the two panels. Contains two elements rho and gamma, each corresponding to the related panel, and each containing default values for entries named cols, option, and alpha.

### Examples

```
set_TS_summary_plot_cols()
```

---

sim_LDA_data	<i>Simulate LDA data from an LDA structure given parameters</i>
--------------	---

---

### Description

For a given set of parameters alpha and Beta and document-specific total word counts, simulate a document-by-term matrix. Additional structuring variables (the numbers of topics (k), documents (M), terms (V)) are inferred from input objects.

### Usage

```
sim_LDA_data(N, Beta, alpha = NULL, Theta = NULL, seed = NULL)
```

### Arguments

N	A vector of document sizes (total word counts). Must be integer conformable. Is used to infer the total number of documents.
Beta	matrix of categorical distribution parameters defining terms within topics. Dimension: k x V (number of topics x number of terms). Used to infer both (k) and (V). Must be non-negative and sum to 1 within topics.
alpha	Single positive numeric value for the Dirichlet distribution parameter defining topics within documents. To specifically define document topic probabilities, use Theta.
Theta	matrix of probabilities defining topics within documents. Dimension: M x k (documents x topics). Must be non-negative and sum to 1 within documents. To generally define document topic probabilities, use alpha.
seed	Input to <a href="#">set.seed</a> .

**Value**

A document-by-term matrix of counts (dim:  $M \times V$ ).

**Examples**

```
N <- c(10, 22, 15, 31)
alpha <- 1.2
Beta <- matrix(c(0.1, 0.1, 0.8, 0.2, 0.6, 0.2), 2, 3, byrow = TRUE)
sim_LDA_data(N, Beta, alpha = alpha)
Theta <- matrix(c(0.2, 0.8, 0.8, 0.2, 0.5, 0.5, 0.9, 0.1), 4, 2,
               byrow = TRUE)
sim_LDA_data(N, Beta, Theta = Theta)
```

---

sim_LDA_TS_data	<i>Simulate LDA_TS data from LDA and TS model structures and parameters</i>
-----------------	---

---

**Description**

For a given set of covariates  $X$ ; parameters  $Beta$ ,  $Eta$ ,  $\rho$ , and  $err$ ; and document-specific time stamps  $tD$  and lengths  $N$ ), simulate a document-by-topic matrix. Additional structuring variables (the numbers of topics ( $k$ ), terms ( $V$ ), documents ( $M$ ), segments ( $S$ ), and covariates per segment ( $C$ )) are inferred from input objects.

**Usage**

```
sim_LDA_TS_data(N, Beta, X, Eta, rho, tD, err = 0, seed = NULL)
```

**Arguments**

$N$	A vector of document sizes (total word counts). Must be integer conformable. Is used to infer the total number of documents.
$Beta$	matrix of categorical distribution parameters defining terms within topics. Dimension: $k \times V$ (number of topics $\times$ number of terms). Used to infer both ( $k$ ) and ( $V$ ). Must be non-negative and sum to 1 within topics.
$X$	matrix of covariates, dimension $M$ (number of documents) $\times$ $C$ (number of covariates, including the intercept) (a.k.a the design matrix).
$Eta$	matrix of regression parameters across the segments, dimension: $SC$ (number of segments $\times$ number of covariates, including the intercept) $\times$ $k$ (number of topics).
$\rho$	Vector of integer-conformable time locations of changepoints or NULL if no changepoints. Used to determine the number of segments. Must exist within the bounds of the times of the documents, $tD$ .
$tD$	Vector of integer-conformable times of the documents. Must be of length $M$ (as determined by $X$ ).

err	Additive error on the link-scale. Must be a non-negative numeric value. Default value of 0 indicates no error.
seed	Input to <code>set.seed</code> .

**Value**

A document-by-term matrix of counts (dim: M x V).

**Examples**

```
N <- c(10, 22, 15, 31)
tD <- c(1, 3, 4, 6)
rho <- 3
X <- cbind(rep(1, 4), 1:4)
Eta <- cbind(c(0.5, 0.3, 0.9, 0.5), c(1.2, 1.1, 0.1, 0.5))
Beta <- matrix(c(0.1, 0.1, 0.8, 0.2, 0.6, 0.2), 2, 3, byrow = TRUE)
err <- 1
sim_LDA_TS_data(N, Beta, X, Eta, rho, tD, err)
```

---

sim\_TS\_data

---

*Simulate TS data from a TS model structure given parameters*


---

**Description**

For a given set of covariates  $X$ ; parameters  $\text{Eta}$ ,  $\text{rho}$ , and  $\text{err}$ ; and document-specific time stamps  $\text{tD}$ , simulate a document-by-topic matrix. Additional structuring variables (numbers of topics ( $k$ ), documents ( $M$ ), segments ( $S$ ), and covariates per segment ( $C$ )) are inferred from input objects.

**Usage**

```
sim_TS_data(X, Eta, rho, tD, err = 0, seed = NULL)
```

**Arguments**

$X$	matrix of covariates, dimension $M$ (number of documents) x $C$ (number of covariates, including the intercept) (a.k.a. the design matrix).
$\text{Eta}$	matrix of regression parameters across the segments, dimension: $SC$ (number of segments x number of covariates, including the intercept) x $k$ (number of topics).
$\text{rho}$	Vector of integer-conformable time locations of changepoints or NULL if no changepoints. Used to determine the number of segments. Must exist within the bounds of the times of the documents, $\text{tD}$ .
$\text{tD}$	Vector of integer-conformable times of the documents. Must be of length $M$ (as determined by $X$ ).
err	Additive error on the link-scale. Must be a non-negative numeric value. Default value of 0 indicates no error.
seed	Input to <code>set.seed</code> .

**Value**

A document-by-topic matrix of probabilities (dim: M x k).

**Examples**

```
tD <- c(1, 3, 4, 6)
rho <- 3
X <- cbind(rep(1, 4), 1:4)
Eta <- cbind(c(0.5, 0.3, 0.9, 0.5), c(1.2, 1.1, 0.1, 0.5))
sim_TS_data(X, Eta, rho, tD, err = 1)
```

---

softmax

*Calculate the softmax of a vector or matrix of values*

---

**Description**

Calculate the softmax (normalized exponential) of a vector of values or a set of vectors stacked rowwise.

**Usage**

```
softmax(x)
```

**Arguments**

x                    numeric vector or matrix

**Value**

The softmax of x.

**Examples**

```
dat <- matrix(runif(100, -1, 1), 25, 4)
softmax(dat)
softmax(dat[,1])
```

---

step\_chains

---

*Conduct a within-chain step of the ptMCMC algorithm*


---

### Description

This set of functions steps the chains forward one iteration of the within-chain component of the ptMCMC algorithm. `step_chains` is the main function, comprised of a proposal (made by `prop_step`), an evaluation of that proposal (made by `eval_step`), and then an update of the configuration (made by `take_step`).

This set of functions was designed to work within `TS` and specifically `est_changepoints`. They are still hardcoded to do so, but have the capacity to be generalized to work with any estimation via ptMCMC with additional coding work.

### Usage

```
step_chains(i, cpts, inputs)

propose_step(i, cpts, inputs)

eval_step(i, cpts, prop_step, inputs)

take_step(cpts, prop_step, accept_step)
```

### Arguments

<code>i</code>	integer iteration index.
<code>cpts</code>	matrix of change point locations across chains.
<code>inputs</code>	Class <code>ptMCMC_inputs</code> list, containing the static inputs for use within the ptMCMC algorithm.
<code>prop_step</code>	Proposed step output from <code>propose_step</code> .
<code>accept_step</code>	logical indicator of acceptance of each chain's proposed step.

### Details

For each iteration of the ptMCMC algorithm, all of the chains have the potential to take a step. The possible step is proposed under a proposal distribution (here for change points we use a symmetric geometric distribution), the proposed step is then evaluated and either accepted or not (following the Metropolis-Hastings rule; Metropolis, *et al.* 1953, Hasting 1960, Gupta *et al.* 2018), and then accordingly taken or not (the configurations are updated).

### Value

`step_chains`: list of change points, log-likelihoods, and logical indicators of acceptance for each chain.

propose\_step: list of change points and log-likelihood values for the proposal.

eval\_step: logical vector indicating if each chain's proposal was accepted.

take\_step: list of change points, log-likelihoods, and logical indicators of acceptance for each chain.

## References

Gupta, S., L. Hainsworth, J. S. Hogg, R. E. C. Lee, and J. R. Faeder. 2018. Evaluation of parallel tempering to accelerate Bayesian parameter estimation in systems biology. [link](#).

Hastings, W. K. 1970. Monte Carlo sampling methods using Markov Chains and their applications. *Biometrika* **57**:97-109. [link](#).

Metropolis, N., A. W. Rosenbluth, M. N. Rosenbluth, A. H. Teller, and E. Teller. 1953. Equations of state calculations by fast computing machines. *Journal of Chemical Physics* **21**: 1087-1092. [link](#).

## Examples

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
data <- data[order(data[, "newmoon"]), ]
saves <- prep_saves(1, TS_control())
inputs <- prep_ptMCMC_inputs(data, gamma ~ 1, 1, "newmoon", weights,
                             TS_control())
cpts <- prep_cpts(data, gamma ~ 1, 1, "newmoon", weights, TS_control())
ids <- prep_ids(TS_control())
for(i in 1:TS_control()$nit){
  steps <- step_chains(i, cpts, inputs)
  swaps <- swap_chains(steps, inputs, ids)
  saves <- update_saves(i, saves, steps, swaps)
  cpts <- update_cpts(cpts, swaps)
  ids <- update_ids(ids, swaps)
}
# within step_chains()
cpts <- prep_cpts(data, gamma ~ 1, 1, "newmoon", weights, TS_control())
i <- 1
prop_step <- propose_step(i, cpts, inputs)
accept_step <- eval_step(i, cpts, prop_step, inputs)
take_step(cpts, prop_step, accept_step)
```

---

summarize_etas	<i>Summarize the regressor (eta) distributions</i>
----------------	--

---

**Description**

summarize\_etas calculates summary statistics for each of the chunk-level regressors.

measure\_etas\_vcov generates the variance-covariance matrix for the regressors.

**Usage**

```
summarize_etas(etas, control = list())
```

```
measure_eta_vcov(etas)
```

**Arguments**

etas	Matrix of regressors (columns) across iterations of the ptMCMC (rows), as returned from <a href="#">est_regressors</a> .
control	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <a href="#">TS_control</a> .

**Value**

summarize\_etas: table of summary statistics for chunk-level regressors including mean, median, mode, posterior interval, standard deviation, MCMC error, autocorrelation, and effective sample size for each regressor.

measure\_eta\_vcov: variance-covariance matrix for chunk-level regressors.

**Examples**

```
etas <- matrix(rnorm(100), 50, 2)
summarize_etas(etas)
measure_eta_vcov(etas)
```

---

summarize_rhos	<i>Summarize the rho distributions</i>
----------------	--

---

**Description**

summarize\_rho calculates summary statistics for each of the change point locations.

measure\_rho\_vcov generates the variance-covariance matrix for the change point locations.



**Usage**

```
summarize_rhos(rhos, control = list())

measure_rho_vcov(rhos)
```

**Arguments**

**rhos** Matrix of change point locations (columns) across iterations of the ptMCMC (rows) or NULL if no change points are in the model, as returned from [est\\_changepoints](#).

**control** A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by [TS\\_control](#).

**Value**

**summarize\_rhos**: table of summary statistics for change point locations including mean, median, mode, posterior interval, standard deviation, MCMC error, autocorrelation, and effective sample size for each change point location.

**measure\_rho\_vcov**: variance-covariance matrix for change point locations.

**Examples**

```
rhos <- matrix(sample(80:100, 100, TRUE), 50, 2)
summarize_rhos(rhos)
measure_rho_vcov(rhos)
```

---

 swap\_chains

---

*Conduct a set of among-chain swaps for the ptMCMC algorithm*


---

**Description**

This function handles the among-chain swapping based on temperatures and likelihood differentials.

This function was designed to work within [TS](#) and specifically [est\\_changepoints](#). It is still hard-coded to do so, but has the capacity to be generalized to work with any estimation via ptMCMC with additional coding work.

**Usage**

```
swap_chains(chainsin, inputs, ids)
```

**Arguments**

**chainsin** Chain configuration to be evaluated for swapping.

**inputs** Class ptMCMC\_inputs list, containing the static inputs for use within the ptMCMC algorithm.

**ids** The vector of integer chain ids.

## Details

The ptMCMC algorithm couples the chains (which are taking their own walks on the distribution surface) through "swaps", where neighboring chains exchange configurations (Geyer 1991, Falcioni and Deem 1999) following the Metropolis criterion (Metropolis *et al.* 1953). This allows them to share information and search the surface in combination (Earl and Deem 2005).

## Value

list of updated change points, log-likelihoods, and chain ids, as well as a vector of acceptance indicators for each swap.

## References

Earl, D. J. and M. W. Deem. 2005. Parallel tempering: theory, applications, and new perspectives. *Physical Chemistry Chemical Physics* **7**: 3910-3916. [link](#).

Falcioni, M. and M. W. Deem. 1999. A biased Monte Carlo scheme for zeolite structure solution. *Journal of Chemical Physics* **110**: 1754-1766. [link](#).

Geyer, C. J. 1991. Markov Chain Monte Carlo maximum likelihood. *In Computing Science and Statistics: Proceedings of the 23rd Symposium on the Interface*. pp 156-163. American Statistical Association, New York, USA. [link](#).

Metropolis, N., A. W. Rosenbluth, M. N. Rosenbluth, A. H. Teller, and E. Teller. 1953. Equations of state calculations by fast computing machines. *Journal of Chemical Physics* **21**: 1087-1092. [link](#).

## Examples

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
data <- data[order(data[,"newmoon"]), ]
saves <- prep_saves(1, TS_control())
inputs <- prep_ptMCMC_inputs(data, gamma ~ 1, 1, "newmoon", weights,
                           TS_control())
cpts <- prep_cpts(data, gamma ~ 1, 1, "newmoon", weights, TS_control())
ids <- prep_ids(TS_control())
for(i in 1:TS_control()$nit){
  steps <- step_chains(i, cpts, inputs)
  swaps <- swap_chains(steps, inputs, ids)
  saves <- update_saves(i, saves, steps, swaps)
  cpts <- update_cpts(cpts, swaps)
  ids <- update_ids(ids, swaps)
}
```

---

trace_plot	<i>Produce the trace plot panel for the TS diagnostic plot of a parameter</i>
------------	---

---

### Description

Produce a trace plot for the parameter of interest (rho or eta) as part of `TS_diagnostics_plot`. A horizontal line is added to show the median of the posterior.

### Usage

```
trace_plot(x, ylab = "parameter value")
```

### Arguments

x	Vector of parameter values drawn from the posterior distribution, indexed to the iteration by the order of the vector.
ylab	character value used to label the y axis.

### Value

NULL.

### Examples

```
trace_plot(rnorm(100, 0, 1))
```

---

TS	<i>Conduct a single multinomial Bayesian Time Series analysis</i>
----	---

---

### Description

This is the main interface function for the LDATS application of Bayesian change point Time Series analyses (Christensen *et al.* 2018), which extends the model of Western and Kleykamp (2004; see also Ruggieri 2013) to multinomial (proportional) response data using softmax regression (Ripley 1996, Venables and Ripley 2002, Bishop 2006) using a generalized linear modeling approach (McCullagh and Nelder 1989). The models are fit using parallel tempering Markov Chain Monte Carlo (ptMCMC) methods (Earl and Deem 2005) to locate change points and neural networks (Ripley 1996, Venables and Ripley 2002, Bishop 2006) to estimate regressors.

`check_TS_inputs` checks that the inputs to TS are of proper classes for a full analysis.

**Usage**

```
TS(
  data,
  formula = gamma ~ 1,
  nchangepoints = 0,
  timename = "time",
  weights = NULL,
  control = list()
)
```

```
check_TS_inputs(
  data,
  formula = gamma ~ 1,
  nchangepoints = 0,
  timename = "time",
  weights = NULL,
  control = list()
)
```

**Arguments**

<code>data</code>	<code>data.frame</code> including [1] the time variable (indicated in <code>timename</code> ), [2] the predictor variables (required by <code>formula</code> ) and [3], the multinomial response variable (indicated in <code>formula</code> ) as verified by <a href="#">check_timename</a> and <a href="#">check_formula</a> . Note that the response variables should be formatted as a <code>data.frame</code> object named as indicated by the response entry in the <code>control</code> list, such as <code>gamma</code> for a standard TS analysis on LDA output. See Examples.
<code>formula</code>	<a href="#">formula</a> defining the regression between relationship the change points. Any predictor variable included must also be a column in <code>data</code> and any (multinomial) response variable must be a set of columns in <code>data</code> , as verified by <a href="#">check_formula</a> .
<code>nchangepoints</code>	integer corresponding to the number of change points to include in the model. 0 is a valid input (corresponding to no change points, so a singular time series model), and the current implementation can reasonably include up to 6 change points. The number of change points is used to dictate the segmentation of the time series into chunks fit with separate models dictated by <code>formula</code> .
<code>timename</code>	character element indicating the time variable used in the time series. Defaults to "time". The variable must be integer-conformable or a Date. If the variable named is a Date, the input is converted to an integer, resulting in the timestep being 1 day, which is often not desired behavior.
<code>weights</code>	Optional class <code>numeric</code> vector of weights for each document. Defaults to <code>NULL</code> , translating to an equal weight for each document. When using <code>multinom_TS</code> in a standard LDATS analysis, it is advisable to weight the documents by their total size, as the result of LDA is a matrix of proportions, which does not account for size differences among documents. For most models, a scaling of the weights (so that the average is 1) is most appropriate, and this is accomplished using <code>document_weights</code> .

**control** A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by [TS\\_control](#).

### Value

TS: TS\_fit-class list containing the following elements, many of which are hidden for printing, but are accessible:

**data** data input to the function.

**formula** [formula](#) input to the function.

**nchangepoints** nchangepoints input to the function.

**weights** weights input to the function.

**control** control input to the function.

**lls** Iteration-by-iteration [logLik](#) values for the full time series fit by [multinom\\_TS](#).

**rhos** Iteration-by-iteration change point estimates from [est\\_changepoints](#).

**etas** Iteration-by-iteration marginal regressor estimates from [est\\_regressors](#), which have been unconditioned with respect to the change point locations.

**ptMCMC\_diagnostics** ptMCMC diagnostics, see [diagnose\\_ptMCMC](#)

**rho\_summary** Summary table describing rhos (the change point locations), see [summarize\\_rhos](#).

**rho\_vcov** Variance-covariance matrix for the estimates of rhos (the change point locations), see [measure\\_rho\\_vcov](#).

**eta\_summary** Summary table describing etas (the regressors), see [summarize\\_etas](#).

**eta\_vcov** Variance-covariance matrix for the estimates of etas (the regressors), see [measure\\_eta\\_vcov](#).

**logLik** Across-iteration average of log-likelihoods (lls).

**nparams** Total number of parameters in the full model, including the change point locations and regressors.

**deviance** Penalized negative log-likelihood, based on [logLik](#) and [nparams](#).

**check\_TS\_inputs**: An error message is thrown if any input is not proper, else NULL.

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### Examples

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)

TSmod <- TS(data, gamma ~ 1, nchangepoints = 1, "newmoon", weights)

check_TS_inputs(data, timename = "newmoon")
```

---

TS\_control

*Create the controls list for the Time Series model*

---

### Description

This function provides a simple creation and definition of a list used to control the time series model fit occurring within [TS](#).

### Usage

```
TS_control(
  memoise = TRUE,
  response = "gamma",
  lambda = 0,
  measurer = AIC,
  selector = min,
  ntemps = 6,
  penultimate_temp = 2^6,
  ultimate_temp = 1e+10,
  q = 0,
  nit = 10000,
  magnitude = 12,
  quiet = FALSE,
  burnin = 0,
  thin_frac = 1,
  summary_prob = 0.95,
  seed = NULL
)
```

**Arguments**

memoise	logical indicator of whether the multinomial functions should be memoised (via <a href="#">memoise</a> ). Memoisation happens to both <a href="#">multinom_TS</a> and <a href="#">multinom_TS_chunk</a> .
response	character element indicating the response variable used in the time series.
lambda	numeric "weight" decay term used to set the prior on the regressors within each chunk-level model. Defaults to 0, corresponding to a fully vague prior.
measurer, selector	Function names for use in evaluation of the TS models. <code>measurer</code> is used to create a value for each model and <code>selector</code> operates on the values to choose the model.
ntemps	integer number of temperatures (chains) to use in the ptMCMC algorithm.
penultimate_temp	Penultimate temperature in the ptMCMC sequence.
ultimate_temp	Ultimate temperature in the ptMCMC sequence.
q	Exponent controlling the ptMCMC temperature sequence from the focal chain (reference with temperature = 1) to the penultimate chain. 0 (default) implies a geometric sequence. 1 implies squaring before exponentiating.
nit	integer number of iterations (steps) used in the ptMCMC algorithm.
magnitude	Average magnitude (defining a geometric distribution) for the proposed step size in the ptMCMC algorithm.
quiet	logical indicator of whether the model should run quietly (if FALSE, a progress bar and notifications are printed).
burnin	integer number of iterations to remove from the beginning of the ptMCMC algorithm.
thin_frac	Fraction of iterations to retain, must be $(0, 1]$ , and the default value of 1 represents no thinning.
summary_prob	Probability used for summarizing the posterior distributions (via the highest posterior density interval, see <a href="#">HPDinterval</a> ).
seed	Input to <code>set.seed</code> for replication purposes.

**Value**

list, with named elements corresponding to the arguments.

**Examples**

```
TS_control()
```

---

TS\_diagnostics\_plot *Plot the diagnostics of the parameters fit in a TS model*

---

### Description

Plot 4-panel figures (showing trace plots, posterior ECDF, posterior density, and iteration autocorrelation) for each of the parameters (change point locations and regressors) fitted within a multinomial time series model (fit by [TS](#)).

`eta_diagnostics_plots` creates the diagnostic plots for the regressors (etas) of a time series model.

`rho_diagnostics_plots` creates the diagnostic plots for the change point locations (rho) of a time series model.

### Usage

```
TS_diagnostics_plot(x, interactive = TRUE)
```

```
eta_diagnostics_plots(x, interactive)
```

```
rho_diagnostics_plots(x, interactive)
```

### Arguments

`x` Object of class `TS_fit`, generated by [TS](#) to have its diagnostics plotted.  
`interactive` logical input, should be code `TRUE` unless testing.

### Value

NULL.

### Examples

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
TSmod <- TS(data, gamma ~ 1, nchangepts = 1, "newmoon", weights)
TS_diagnostics_plot(TSmod)
```



---

 TS\_on\_LDA

---

*Conduct a set of Time Series analyses on a set of LDA models*


---

### Description

This is a wrapper function that expands the main Time Series analyses function ([TS](#)) across the LDA models (estimated using [LDA](#) or [LDA\\_set](#) and the Time Series models, with respect to both continuous time formulas and the number of discrete changepoints. This function allows direct passage of the control parameters for the parallel tempering MCMC through to the main Time Series function, [TS](#), via the `ptMCMC_controls` argument.

`check_TS_on_LDA_inputs` checks that the inputs to `TS_on_LDA` are of proper classes for a full analysis.

### Usage

```
TS_on_LDA(
  LDA_models,
  document_covariate_table,
  formulas = ~1,
  nchangepoints = 0,
  timename = "time",
  weights = NULL,
  control = list()
)

check_TS_on_LDA_inputs(
  LDA_models,
  document_covariate_table,
  formulas = ~1,
  nchangepoints = 0,
  timename = "time",
  weights = NULL,
  control = list()
)
```

### Arguments

`LDA_models` List of LDA models (class `LDA_set`, produced by [LDA\\_set](#)) or a singular LDA model (class `LDA`, produced by [LDA](#)).

`document_covariate_table` Document covariate table (rows: documents, columns: time index and covariate options). Every model needs a covariate to describe the time value for each document (in whatever units and whose name in the table is input in `timename`) that dictates the application of the change points. In addition, all covariates named within specific models in `formula` must be included. Must be a conformable to a data table, as verified by [check\\_document\\_covariate\\_table](#).

formulas	Vector of <code>formula</code> (s) for the continuous (non-change point) component of the time series models. Any predictor variable included in a formula must also be a column in the <code>document_covariate_table</code> . Each element (formula) in the vector is evaluated for each number of change points and each LDA model.
nchangepoints	Vector of integers corresponding to the number of change points to include in the time series models. 0 is a valid input corresponding to no change points ( <i>i.e.</i> , a singular time series model), and the current implementation can reasonably include up to 6 change points. Each element in the vector is the number of change points used to segment the data for each formula (entry in <code>formulas</code> ) component of the TS model, for each selected LDA model.
timename	character element indicating the time variable used in the time series. Defaults to "time". The variable must be integer-conformable or a Date. If the variable named is a Date, the input is converted to an integer, resulting in the timestep being 1 day, which is often not desired behavior.
weights	Optional class <code>numeric</code> vector of weights for each document. Defaults to <code>NULL</code> , translating to an equal weight for each document. When using <code>multinom_TS</code> in a standard LDATS analysis, it is advisable to weight the documents by their total size, as the result of <code>LDA</code> is a matrix of proportions, which does not account for size differences among documents. For most models, a scaling of the weights (so that the average is 1) is most appropriate, and this is accomplished using <code>document_weights</code> .
control	A list of parameters to control the fitting of the Time Series model including the parallel tempering Markov Chain Monte Carlo (ptMCMC) controls. Values not input assume defaults set by <code>TS_control</code> .

## Value

`TS_on_LDA`: `TS_on_LDA`-class list of results from `TS` applied for each model on each LDA model input.

`check_TS_inputs`: An error message is thrown if any input is not proper, else `NULL`.

## Examples

```
data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDAs <- LDA_set(document_term_table, topics = 2:3, nseeds = 2)
LDA_models <- select_LDA(LDAs)
weights <- document_weights(document_term_table)
formulas <- c(~ 1, ~ newmoon)
mods <- TS_on_LDA(LDA_models, document_covariate_table, formulas,
                 nchangepoints = 0:1, timename = "newmoon", weights)
```



cols	list of elements used to define the colors for the two panels, as generated simply using <code>set_TS_summary_plot_cols</code> . Has two elements rho and gamma, each corresponding to the related panel, and each containing default values for entries named cols, option, and alpha. See <code>set_gamma_colors</code> and <code>set_rho_hist_colors</code> for details on usage.
bin_width	Width of the bins used in the histograms, in units of the x-axis (the time variable used to fit the model).
xname	Label for the x-axis in the summary time series plot. Defaults to NULL, which results in usage of the <code>timename</code> element of the control list (held in <code>control\$TS_control\$timename</code> ). To have no label printed, set <code>xname = ""</code> .
border	Border for the histogram, default is NA.
selection	Indicator of the change points to use. Currently only defined for "median" and "mode".
LDATS	logical indicating if the plot is part of a larger LDATS plot output.
together	logical indicating if the subplots are part of a larger LDA plot output.

**Value**

NULL.

**Examples**

```

data(rodents)
document_term_table <- rodents$document_term_table
document_covariate_table <- rodents$document_covariate_table
LDA_models <- LDA_set(document_term_table, topics = 2)[[1]]
data <- document_covariate_table
data$gamma <- LDA_models@gamma
weights <- document_weights(document_term_table)
TSmod <- TS(data, gamma ~ 1, nchangepoints = 1, "newmoon", weights)
TS_summary_plot(TSmod)
pred_gamma_TS_plot(TSmod)
rho_hist(TSmod)

```

---

verify\_changepoint\_locations

*Verify the change points of a multinomial time series model*

---

**Description**

Verify that a time series can be broken into a set of chunks based on input change points.

**Usage**

```
verify_changepoint_locations(data, changepoints = NULL, timename = "time")
```

**Arguments**

data	Class data.frame object including the predictor and response variables.
changepoints	Numeric vector indicating locations of the change points. Must be conformable to integer values.
timename	character element indicating the time variable used in the time series. Defaults to "time". The variable must be integer-conformable or a Date. If the variable named is a Date, the input is converted to an integer, resulting in the timestep being 1 day, which is often not desired behavior.

**Value**

Logical indicator of the check passing TRUE or failing FALSE.

**Examples**

```
data(rodents)
dtt <- rodents$document_term_table
lda <- LDA_set(dtt, 2, 1, list(quiet = TRUE))
dct <- rodents$document_covariate_table
dct$gamma <- lda[[1]]@gamma
verify_changepoint_locations(dct, changepoints = 100,
                             timename = "newmoon")
```

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