

Package ‘mdgc’

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Type Package

Title Missing Data Imputation Using Gaussian Copulas

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Description Provides functions to impute missing values using Gaussian copulas for mixed data types as described by Christoffersen et al. (2021) <[doi:10.48550/arXiv.2102.02642](https://doi.org/10.48550/arXiv.2102.02642)>. The method is related to Hoff (2007) <[doi:10.1214/07-AOAS107](https://doi.org/10.1214/07-AOAS107)> and Zhao and Udell (2019) <[doi:10.48550/arXiv.1910.12845](https://doi.org/10.48550/arXiv.1910.12845)> but differs by making a direct approximation of the log marginal likelihood using an extended version of the Fortran code created by Genz and Bretz (2002) <[doi:10.1198/106186002394](https://doi.org/10.1198/106186002394)> in addition to also support multinomial variables.

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BugReports <https://github.com/boennecd/mdgc/issues>

URL <https://github.com/boennecd/mdgc>

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Contents

mdgc-package	2
get_mdgc	3
get_mdgc_log_ml	4
mdgc	6
mdgc_fit	9
mdgc_impute	11
mdgc_log_ml	13
mdgc_start_value	15

Index	17
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mdgc-package	<i>mdgc: Missing Data imputation using Gaussian Copulas</i>
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Description

The mdgc package is used to estimate Gaussian Copula models for mixed data types (continuous, binary, ordinal, and multinomial) that can be used for imputations. The main function is the `mdgc` function. The rest of the functions in the package give the user access to lower level functions.

Examples are provided at <https://github.com/boennecd/mdgc>. The package is still in a development stage and the API may change.

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References

Christoffersen, B., Clements, M., Humphreys, K., & Kjellström, H. (2021). *Asymptotically Exact and Fast Gaussian Copula Models for Imputation of Mixed Data Types*. <https://arxiv.org/abs/2102.02642>.

See Also

Useful links:

- <https://github.com/boennecd/mdgc>
- Report bugs at <https://github.com/boennecd/mdgc/issues>

`get_mdgc`*Get mdgc Object*

Description

Creates a mdgc object which is needed for estimation of the covariance matrix and the mean vector and to perform imputation.

Usage

```
get_mdgc(dat)
```

Arguments

`dat` [data.frame](#) with continuous, multinomial, ordinal, and binary variables.

Details

It is important to use appropriate classes for the [data.frame](#) columns:

- Continuous variables: should be [numerics](#).
- Binary variables: should be [logicals](#).
- Multinomial variables: should be [factors](#).
- Ordinal variables: should be [ordered](#).

Value

An object of class mdgc. It has the following elements:

`lower`, `upper`, `code`, `multinomial`, `idx_non_zero_mean`
arguments to pass to [get_mdgc_log_ml](#).

`margs` functions to get lower and upper bounds for each column of `dat`.

`reals`, `bins`, `ords`
indices of continuous, binary, and ordinal variables, respectively.

`truth` the numeric version of `dat`.

`means` starting values for the non-zero mean terms (see e.g. [mdgc_fit](#)).

See Also

[get_mdgc_log_ml](#), [mdgc_start_value](#)

Examples

```

# there is a bug on CRAN's check on Solaris which I have failed to reproduce.
# See https://github.com/r-hub/solarischeck/issues/8#issuecomment-796735501.
# Thus, this example is not run on Solaris
is_solaris <- tolower(Sys.info()[["sysname"]]) == "sunos"

if(!is_solaris){
  # randomly mask data
  set.seed(11)
  masked_data <- iris
  masked_data[matrix(runif(prod(dim(iris))) < .10, NROW(iris))] <- NA

  # use the functions in the package
  library(mdgc)
  obj <- get_mdgc(masked_data)
  print(class(obj))
}

```

get_mdgc_log_ml

Get Pointer to C++ Object to Approximate the Log Marginal Likelihood

Description

Creates a C++ object which is needed to approximate the log marginal likelihood. The object cannot be saved.

Usage

```

get_mdgc_log_ml(object, ...)

## S3 method for class 'mdgc'
get_mdgc_log_ml(object, ...)

## S3 method for class 'data.frame'
get_mdgc_log_ml(object, ...)

## Default S3 method:
get_mdgc_log_ml(
  object,
  lower,
  upper,
  code,
  multinomial,
  idx_non_zero_mean,
  ...
)

```

Arguments

object	mdgc object from <code>get_mdgc</code> or a <code>data.frame</code> to pass to <code>get_mdgc</code> . Ignored by the default method.
...	used to pass arguments to S3 methods.
lower	[# variables]x[# observations] matrix with lower bounds for each variable on the normal scale.
upper	[# variables]x[# observations] matrix with upper bounds for each variable on the normal scale.
code	[# variables]x[# observations] matrix integer code for the each variable on the normal scale. Zero implies an observed value (the value in upper), one implies a missing value, and two implies an interval.
multinomial	<code>list</code> where each element is 3x[# multinomial variables] <code>matrix</code> with multinomial outcomes. The first index is the outcome as an integer code, the second index is the number of categories, and the third index is the index of each multinomial variable (this is zero-based).
idx_non_zero_mean	indices for non-zero mean variables. Indices should be sorted.

Details

Indices are zero-based except the outcome index for multinomial variables.

`idx_non_zero_mean` indices with terms with code equal to zero (observed values) are ignored.

Value

A Rcpp::XPtr to pass to e.g. `mdgc_log_ml`.

See Also

[mdgc_fit](#), [mdgc_log_ml](#)

Examples

```
# there is a bug on CRAN's check on Solaris which I have failed to reproduce.
# See https://github.com/r-hub/solarischeck/issues/8#issuecomment-796735501.
# Thus, this example is not run on Solaris
is_solaris <- tolower(Sys.info()[["sysname"]]) == "sunos"

if(!is_solaris){
  # randomly mask data
  set.seed(11)
  masked_data <- iris
  masked_data[matrix(runif(prod(dim(iris))) < .10, NROW(iris))] <- NA

  # use the functions in the package
  library(mdgc)
  obj <- get_mdgc(masked_data)
  ptr <- get_mdgc_log_ml(obj)
```

```
}
```

mdgc

Perform Model Estimation and Imputation

Description

A convenience function to perform model estimation and imputation in one call. The learning rate is likely model specific and should be altered. See [mdgc_fit](#).

See the README at <https://github.com/boennecd/mdgc> for examples.

Usage

```
mdgc(  
  dat,  
  lr = 0.001,  
  maxit = 25L,  
  batch_size = NULL,  
  rel_eps = 0.001,  
  method = c("svrg", "adam", "aug_Lagran"),  
  seed = 1L,  
  epsilon = 1e-08,  
  beta_1 = 0.9,  
  beta_2 = 0.999,  
  n_threads = 1L,  
  do_reorder = TRUE,  
  abs_eps = -1,  
  maxpts = 10000L,  
  minvls = 100L,  
  verbose = FALSE,  
  irel_eps = rel_eps,  
  imaxit = maxpts,  
  iabs_eps = abs_eps,  
  iminvls = 1000L,  
  start_val = NULL,  
  decay = 0.98,  
  conv_crit = 1e-05,  
  use_aprx = FALSE  
)
```

Arguments

<code>dat</code>	<code>data.frame</code> with continuous, multinomial, ordinal, and binary variables.
<code>lr</code>	learning rate.
<code>maxit</code>	maximum number of iteration.

<code>batch_size</code>	number of observations in each batch.
<code>rel_eps</code>	relative error for each marginal likelihood factor.
<code>method</code>	estimation method to use. Can be "svrg", "adam", or "aug_Lagran".
<code>seed</code>	fixed seed to use. Use NULL if the seed should not be fixed.
<code>epsilon</code>	ADAM parameters.
<code>beta_1</code>	ADAM parameters.
<code>beta_2</code>	ADAM parameters.
<code>n_threads</code>	number of threads to use.
<code>do_reorder</code>	logical for whether to use a heuristic variable reordering. TRUE is likely the best option.
<code>abs_eps</code>	absolute convergence threshold for each marginal likelihood factor.
<code>maxpts</code>	maximum number of samples to draw for each marginal likelihood term.
<code>minvls</code>	minimum number of samples.
<code>verbose</code>	logical for whether to print output during the estimation.
<code>irel_eps</code>	relative error for each term in the imputation.
<code>imaxit</code>	maximum number of samples to draw in the imputation.
<code>iabs_eps</code>	absolute convergence threshold for each term in the imputation.
<code>iminvls</code>	minimum number of samples in the imputation.
<code>start_val</code>	starting value for the covariance matrix. Use NULL if unspecified.
<code>decay</code>	the learning rate used by SVRG is given by $lr * decay^{iteration_number}$.
<code>conv_crit</code>	relative convergence threshold.
<code>use_aprx</code>	logical for whether to use an approximation of <code>pnorm</code> and <code>qnorm</code> . This may yield a noticeable reduction in the computation time.

Details

It is important that the input for data has the appropriate types and classes. See [get_mdgc](#).

Value

A list with the following entries:

<code>ximp</code>	data.frame with the observed and imputed values.
<code>imputed</code>	output from mdgc_impute .
<code>vcov</code>	the estimated covariance matrix.
<code>mea</code>	the estimated non-zero mean terms.

Additional elements may be present depending on the chosen method. See [mdgc_fit](#).

References

- Kingma, D.P., & Ba, J. (2015). *Adam: A Method for Stochastic Optimization*. abs/1412.6980.
- Johnson, R., & Zhang, T. (2013). *Accelerating stochastic gradient descent using predictive variance reduction*. In Advances in neural information processing systems.

See Also

[get_mdgc](#), [mdgc_start_value](#), [get_mdgc_log_ml](#), [mdgc_fit](#), [mdgc_impute](#)

Examples

```
# there is a bug on CRAN's check on Solaris which I have failed to reproduce.
# See https://github.com/r-hub/solarischeck/issues/8#issuecomment-796735501.
# Thus, this example is not run on Solaris
is_solaris <- tolower(Sys.info()[["sysname"]]) == "sunos"

if(!is_solaris && require(catdata)){
  data(retinopathy)

  # prepare data and save true data set
  retinopathy$RET <- as.ordered(retinopathy$RET)
  retinopathy$SM <- as.logical(retinopathy$SM)

  # randomly mask data
  set.seed(28325145)
  truth <- retinopathy
  for(i in seq_along(retinopathy))
    retinopathy[[i]][runif(NROW(retinopathy)) < .3] <- NA

  cat("\nMasked data:\n")
  print(head(retinopathy, 10))
  cat("\n")

  # impute data
  impu <- mdgc(retinopathy, lr = 1e-3, maxit = 25L, batch_size = 25L,
              rel_eps = 1e-3, maxpts = 5000L, verbose = TRUE,
              n_threads = 1L, method = "svrg")

  # show correlation matrix
  cat("\nEstimated correlation matrix\n")
  print(impu$vcov)

  # compare imputed and true values
  cat("\nObserved;\n")
  print(head(retinopathy, 10))
  cat("\nImputed values:\n")
  print(head(impu$ximp, 10))
  cat("\nTruth:\n")
  print(head(truth, 10))

  # using augmented Lagrangian method
  cat("\n")
  impu_aug <- mdgc(retinopathy, maxit = 25L, rel_eps = 1e-3,
                  maxpts = 5000L, verbose = TRUE,
                  n_threads = 1L, method = "aug_Lagran")

  # compare the log-likelihood estimate
  obj <- get_mdgc_log_ml(retinopathy)
```

```

cat(sprintf(
  "Maximum log likelihood with SVRG vs. augmented Lagrangian:\n %.2f vs. %.2f\n",
  mdgc_log_ml(obj, vcov = impu $vcov, mea = impu $mea, rel_eps = 1e-3),
  mdgc_log_ml(obj, vcov = impu_aug$vcov, mea = impu_aug$mea, rel_eps = 1e-3))

# show correlation matrix
cat("\nEstimated correlation matrix (augmented Lagrangian)\n")
print(impu_aug$vcov)

cat("\nImputed values (augmented Lagrangian):\n")
print(head(impu_aug$ximp, 10))
}

```

mdgc_fit

Estimate the Model Parameters

Description

Estimates the covariance matrix and the non-zero mean terms. The `lr` parameter and the `batch_size` parameter are likely data dependent. Convergence should be monitored e.g. by using `verbose = TRUE` with `method = "svrg"`.

See the README at <https://github.com/boennecd/mdgc> for examples.

Usage

```

mdgc_fit(
  ptr,
  vcov,
  mea,
  lr = 0.001,
  rel_eps = 0.001,
  maxit = 25L,
  batch_size = NULL,
  method = c("svrg", "adam", "aug_Lagran"),
  seed = 1L,
  epsilon = 1e-08,
  beta_1 = 0.9,
  beta_2 = 0.999,
  n_threads = 1L,
  do_reorder = TRUE,
  abs_eps = -1,
  maxpts = 10000L,
  minvls = 100L,
  verbose = FALSE,
  decay = 0.98,
  conv_crit = 1e-06,

```

```

    use_aprx = FALSE,
    mu = 1,
    lambda = NULL
  )

```

Arguments

<code>ptr</code>	returned object from <code>get_mdgc_log_ml</code> .
<code>vcov, mea</code>	starting value for the covariance matrix and the non-zero mean entries.
<code>lr</code>	learning rate.
<code>rel_eps</code>	relative error for each marginal likelihood factor.
<code>maxit</code>	maximum number of iteration.
<code>batch_size</code>	number of observations in each batch.
<code>method</code>	estimation method to use. Can be "svrg", "adam", or "aug_Lagran".
<code>seed</code>	fixed seed to use. Use NULL if the seed should not be fixed.
<code>epsilon, beta_1, beta_2</code>	ADAM parameters.
<code>n_threads</code>	number of threads to use.
<code>do_reorder</code>	logical for whether to use a heuristic variable reordering. TRUE is likely the best option.
<code>abs_eps</code>	absolute convergence threshold for each marginal likelihood factor.
<code>maxpts</code>	maximum number of samples to draw for each marginal likelihood term.
<code>minvls</code>	minimum number of samples.
<code>verbose</code>	logical for whether to print output during the estimation.
<code>decay</code>	the learning rate used by SVRG is given by $lr * decay^{iteration_number}$.
<code>conv_crit</code>	relative convergence threshold.
<code>use_aprx</code>	logical for whether to use an approximation of <code>pnorm</code> and <code>qnorm</code> . This may yield a noticeable reduction in the computation time.
<code>mu</code>	starting value for the penalty in the augmented Lagrangian method.
<code>lambda</code>	starting values for the Lagrange multiplier estimates. NULL yields a default.

Value

An `list` with the following elements:

<code>result</code>	<code>list</code> with two elements: <code>vcov</code> is the estimated covariance matrix and <code>mea</code> is the estimated non-zero mean terms.
<code>estimates</code>	If present, the estimated parameters after each iteration.
<code>fun_vals</code>	If present, the output of <code>mdgc_log_ml</code> after each iteration.
<code>mu, lambda</code>	If present, the <code>mu</code> and <code>lambda</code> values at the end.

The elements that may be present depending on the chosen method.

References

Kingma, D.P., & Ba, J. (2015). *Adam: A Method for Stochastic Optimization*. abs/1412.6980.

Johnson, R., & Zhang, T. (2013). *Accelerating stochastic gradient descent using predictive variance reduction*. In Advances in neural information processing systems.

See Also

[mdgc_log_ml](#), [mdgc_start_value](#), [mdgc_impute](#).

Examples

```
# there is a bug on CRAN's check on Solaris which I have failed to reproduce.
# See https://github.com/r-hub/solarischeck/issues/8#issuecomment-796735501.
# Thus, this example is not run on Solaris
is_solaris <- tolower(Sys.info()[["sysname"]]) == "sunos"

if(!is_solaris){
  # randomly mask data
  set.seed(11)
  masked_data <- iris
  masked_data[matrix(runif(prod(dim(iris))) < .10, NROW(iris))] <- NA

  # use the functions in the package
  library(mdgc)
  obj <- get_mdgc(masked_data)
  ptr <- get_mdgc_log_ml(obj)
  start_vals <- mdgc_start_value(obj)

  fit <- mdgc_fit(ptr, start_vals, obj$means, rel_eps = 1e-2, maxpts = 10000L,
                 minvls = 1000L, use_aprx = TRUE, batch_size = 100L, lr = .001,
                 maxit = 100L, n_threads = 2L)
  print(fit$result$vcov)
  print(fit$result$mea)
}
```

mdgc_impute

Impute Missing Values

Description

Imputes missing values given a covariance matrix and mean vector using a similar quasi-random numbers method as [mdgc_log_ml](#).

Usage

```
mdgc_impute(
  object,
  vcov,
  mea,
  rel_eps = 0.001,
  maxit = 10000L,
  abs_eps = -1,
  n_threads = 1L,
  do_reorder = TRUE,
  minvls = 1000L,
  use_aprx = FALSE
)
```

Arguments

object	returned object from get_mdgc .
vcov	covariance matrix to condition on in the imputation.
mea	vector with non-zero mean entries to condition on.
rel_eps	relative convergence threshold for each term in the approximation.
maxit	maximum number of samples
abs_eps	absolute convergence threshold for each term in the approximation.
n_threads	number of threads to use.
do_reorder	logical for whether to use a heuristic variable reordering. TRUE is likely the best option.
minvls	minimum number of samples.
use_aprx	logical for whether to use an approximation of pnorm and qnorm . This may yield a noticeable reduction in the computation time.

Value

A list of lists with imputed values for the continuous variables and a vector with probabilities for each level for the ordinal, binary, and multinomial variables.

Examples

```
# there is a bug on CRAN's check on Solaris which I have failed to reproduce.
# See https://github.com/r-hub/solarischeck/issues/8#issuecomment-796735501.
# Thus, this example is not run on Solaris
is_solaris <- tolower(Sys.info()[["sysname"]]) == "sunos"

if(!is_solaris){
  # randomly mask data
  set.seed(11)
  masked_data <- iris
  masked_data[matrix(runif(prod(dim(iris))) < .10, NROW(iris))] <- NA
}
```

```

# use the functions in the package
library(mdgc)
obj <- get_mdgc(masked_data)
ptr <- get_mdgc_log_ml(obj)
start_vals <- mdgc_start_value(obj)

fit <- mdgc_fit(ptr, start_vals, obj$means, rel_eps = 1e-2, maxpts = 10000L,
               minvls = 1000L, use_aprx = TRUE, batch_size = 100L, lr = .001,
               maxit = 100L, n_threads = 2L)

# impute using the estimated values
imputed <- mdgc_impute(obj, fit$result$vcov, fit$result$mea, minvls = 1000L,
                      maxit = 10000L, n_threads = 2L, use_aprx = TRUE)
print(imputed[1:5]) # first 5 observations
print(head(masked_data, 5)) # observed
print(head(iris      , 5)) # truth
}

```

mdgc_log_ml

Evaluate the Log Marginal Likelihood and Its Derivatives

Description

Approximates the log marginal likelihood and the derivatives using randomized quasi-Monte Carlo. The method uses a generalization of the Fortran code by Genz and Bretz (2002).

Mean terms for observed continuous variables are always assumed to be zero.

The returned log marginal likelihood is not a proper log marginal likelihood if the ptr object is constructed from a mdgc object from `get_mdgc` as it does not include the log of the determinants of the Jacobians for the transformation of the continuous variables.

Usage

```

mdgc_log_ml(
  ptr,
  vcov,
  mea,
  rel_eps = 0.01,
  n_threads = 1L,
  comp_derivs = FALSE,
  indices = NULL,
  do_reorder = TRUE,
  maxpts = 100000L,
  abs_eps = -1,
  minvls = 100L,
  use_aprx = FALSE
)

```

Arguments

<code>ptr</code>	object returned by <code>get_mdgc_log_ml</code> .
<code>vcov</code>	covariance matrix.
<code>mea</code>	vector with non-zero mean entries.
<code>rel_eps</code>	relative error for each marginal likelihood factor.
<code>n_threads</code>	number of threads to use.
<code>comp_derivs</code>	logical for whether to approximate the gradient.
<code>indices</code>	integer vector with which terms (observations) to include. Must be zero-based. NULL yields all observations.
<code>do_reorder</code>	logical for whether to use a heuristic variable reordering. TRUE is likely the best option.
<code>maxpts</code>	maximum number of samples to draw for each marginal likelihood term.
<code>abs_eps</code>	absolute convergence threshold for each marginal likelihood factor.
<code>minvls</code>	minimum number of samples.
<code>use_aprx</code>	logical for whether to use an approximation of <code>pnorm</code> and <code>qnorm</code> . This may yield a noticeable reduction in the computation time.

Value

A numeric vector with a single element with the log marginal likelihood approximation. Two attributes are added if `comp_derivs` is TRUE: `"grad_vcov"` for the derivative approximation with respect to `vcov` and `"grad_mea"` for the derivative approximation with respect to `mea`.

References

Genz, A., & Bretz, F. (2002). *Comparison of Methods for the Computation of Multivariate t Probabilities*. Journal of Computational and Graphical Statistics.

Genz, A., & Bretz, F. (2008). *Computation of Multivariate Normal and t Probabilities*. Springer-Verlag, Heidelberg.

See Also

[mdgc_fit](#)

Examples

```
# there is a bug on CRAN's check on Solaris which I have failed to reproduce.
# See https://github.com/r-hub/solarischeck/issues/8#issuecomment-796735501.
# Thus, this example is not run on Solaris
is_solaris <- tolower(Sys.info()[["sysname"]]) == "sunos"

if(!is_solaris){
  # randomly mask data
  set.seed(11)
  masked_data <- iris
  masked_data[matrix(runif(prod(dim(iris))) < .10, NROW(iris))] <- NA
}
```

```

# use the functions in the package
library(mdgc)
obj <- get_mdgc(masked_data)
ptr <- get_mdgc_log_ml(obj)
start_vals <- mdgc_start_value(obj)
print(mdgc_log_ml(ptr, start_vals, obj$means))
print(mdgc_log_ml(ptr, start_vals, obj$means, use_aprx = TRUE))
print(mdgc_log_ml(ptr, start_vals, obj$means, use_aprx = TRUE,
                  comp_derivs = TRUE))
}

```

mdgc_start_value

Get Starting Value for the Covariance Matrix Using a Heuristic

Description

Uses a heuristic to get starting values for the covariance matrix. These can be passed e.g. to [mdgc_fit](#).

Usage

```
mdgc_start_value(object, ...)
```

```
## S3 method for class 'mdgc'
mdgc_start_value(object, ...)
```

```
## Default S3 method:
```

```
mdgc_start_value(
  object,
  lower,
  upper,
  code,
  multinomial,
  idx_non_zero_mean,
  mea,
  n_threads = 1L,
  ...
)
```

Arguments

object	mdgc object from get_mdgc . Ignored by the default method.
...	used to pass arguments to S3 methods.
lower	[# variables]x[# observations] matrix with lower bounds for each variable on the normal scale.

upper	[# variables]x[# observations] matrix with upper bounds for each variable on the normal scale.
code	[# variables]x[# observations] matrix integer code for the each variable on the normal scale. Zero implies an observed value (the value in upper), one implies a missing value, and two implies an interval.
multinomial	<code>list</code> where each element is 3x[# multinomial variables] <code>matrix</code> with multinomial outcomes. The first index is the outcome as an integer code, the second index is the number of categories, and the third index is the index of each multinomial variable (this is zero-based).
idx_non_zero_mean	indices for non-zero mean variables. Indices should be sorted.
mea	vector with non-zero mean entries.
n_threads	number of threads to use.

Value

The starting value for the covariance matrix.

Examples

```
# there is a bug on CRAN's check on Solaris which I have failed to reproduce.
# See https://github.com/r-hub/solarischeck/issues/8#issuecomment-796735501.
# Thus, this example is not run on Solaris
is_solaris <- tolower(Sys.info()[["sysname"]]) == "sunos"

if(!is_solaris){
  # randomly mask data
  set.seed(11)
  masked_data <- iris
  masked_data[matrix(runif(prod(dim(iris))) < .10, NROW(iris))] <- NA

  # use the functions in the package
  library(mdgc)
  obj <- get_mdgc(masked_data)
  ptr <- get_mdgc_log_ml(obj)
  start_vals <- mdgc_start_value(obj)
  print(start_vals) # starting value for the covariance matrix
}
```

Index

`_PACKAGE` (mdgc-package), 2

`data.frame`, 3, 5–7

`factor`, 3

`get_mdgc`, 3, 5, 7, 8, 12, 13, 15

`get_mdgc_log_ml`, 3, 4, 8, 10, 14

`list`, 5, 10, 16

`logical`, 3

`matrix`, 5, 16

`mdgc`, 2, 6

`mdgc-package`, 2

`mdgc_fit`, 3, 5–8, 9, 14, 15

`mdgc_impute`, 7, 8, 11, 11

`mdgc_log_ml`, 5, 10, 11, 13

`mdgc_start_value`, 3, 8, 11, 15

`numeric`, 3

`ordered`, 3

`pnorm`, 7, 10, 12, 14

`qnorm`, 7, 10, 12, 14