

# Package ‘mrgsolve’

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**Type** Package

**Title** Simulate from ODE-Based Models

**Version** 0.10.4

**Maintainer** Kyle T Baron <kyleb@metrumrg.com>

**Description** Fast simulation from ordinary differential equation (ODE) based models typically employed in quantitative pharmacology and systems biology.

**License** GPL (>= 2)

**URL** <https://github.com/metrumresearchgroup/mrgsolve>

**BugReports** <https://github.com/metrumresearchgroup/mrgsolve/issues>

**Depends** R (>= 3.1.2), methods

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**Suggests** lattice, testthat, xml2 (>= 1.2.0), rmarkdown, yaml, knitr, data.table

**LinkingTo** Rcpp (>= 0.12.12), RcppArmadillo (>= 0.7.900.2.0), BH (>= 1.62.0-1)

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'events.R' 'class\_rx.R' 'compile.R' 'data\_set.R' 'datasets.R'  
 'env.R' 'funset.R' 'handle\_spec\_block.R' 'idata\_set.R' 'init.R'  
 'inven.R' 'knobs.R' 'matlist.R' 'matrix.R' 'mcode.R'  
 'model\_include.R' 'modlib.R' 'modspec.R' 'mread.R'  
 'mrgindata.R' 'mrgsim\_q.R' 'mrgsims.R' 'mrgsolve.R' 'nmxml.R'  
 'param.R' 'print.R' 'r\_to\_cpp.R' 'realize\_addl.R' 'relabel.R'  
 'render.R' 'update.R' 'workflows.R'

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---

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| aboutsolver | <i>About the lsoda differential equation solver used by mrgsolve</i> |
|-------------|--|

---

## Description

The differential equation solver is a C++ translation of DLSODA from ODEPACK. The C++ translation was created by Dilawar Singh and hosted here <https://github.com/dilawar/lib soda>. As we understand the history of the code, Heng Li was also involved in early versions of the code written in C. There was a potentially-related project hosted here <https://github.com/sdwfrost/lib lsoda>.

## Details

The C++ translation by Dilawar Singh contains functions that appear to be based on BLAS and LAPACK routines. These functions have been renamed to be distinct from the respective BLAS and LAPACK function names. References are given in the section below.

## History

The following history was recorded in the source code published by Dilawar Singh:

```

/*
 * HISTORY:
 * This is a CPP version of the LSODA library for integration into MOOSE
 * somulator.
 * The original was aquired from
 * http://www.ccl.net/cca/software/SOURCES/C/kinetics2/index.shtml and modified
 * by
 * Heng Li <lh3lh3@gmail.com>. Heng merged several C files into one and added a
 * simpler interface. [Available
 * here](http://lh3lh3.users.sourceforge.net/download/lsoda.c)

```

```
* The original source code came with no license or copyright
* information. Heng Li released his modification under the MIT/X11 license. I
* maintain the same license. I have removed quite a lot of text/comments from
* this library. Please refer to the standard documentation.
*
* Contact: Dilawar Singh <dilawars@ncbs.res.in>
*/
```

## References

1. LAPACK: <https://netlib.org/lapack>
2. BLAS: <https://netlib.org/blas>

---

as.ev

*Coerce an object to class ev*

---

## Description

Coerce an object to class ev

## Usage

```
as.ev(x, ...)
```

```
## S4 method for signature 'data.frame'
as.ev(x, keep_id = TRUE, clean = FALSE, ...)
```

```
## S4 method for signature 'ev'
as.ev(x, ...)
```

## Arguments

|         |  |
|---------|--|
| x       | an object to coerce  |
| ...     | not used   |
| keep_id | if TRUE, ID column is retained if it exists                      |
| clean   | if TRUE, only dosing or ID information is retained in the result |

## Examples

```
data <- data.frame(amt = 100)

as.ev(data)
```

---

as.list,mrgmod-method *Coerce a model object to list*

---

### Description

Coerce a model object to list

### Usage

```
## S4 method for signature 'mrgmod'
as.list(x, deep = FALSE, ...)
```

### Arguments

|      |   |
|------|---|
| x    | mrgmod object   |
| deep | if TRUE, extra information is returned (see details). |
| ...  | not used  |

### Details

If deep is TRUE, then the values for trans,advan, and mindt are returned as well as a summary of internal model functions (with a call to mrgsolve:::funset).

### Slots

- npar: number of parameters
- neq: number of compartments or differential equations
- pars: names of model parameters
- covariates: names of parameters identified as covariates
- cmt: names of model compartments
- param: the parameter list
- init: initial condition list
- omega: \$OMEGA matrices, as a matlist object
- sigma: \$SIGMA matrices, as a matlist object
- fixed: named list of \$FIXED values
- model: model name
- project: model project directory
- soloc: directory where the model is being built
- sodll: complete path to the model shared object
- cfile: path for the model source code file
- shlib: list of compilation information
- start: simulation start time

- end: simulation end time
- delta: simulation time step
- add: additional simulation times
- capture: names of captured data items
- request: compartments requested upon simulation
- cmti: named indices for current output compartments
- capturei: named indices for current output capture
- random: names and labels of \$OMEGA and \$SIGMA
- code: model source code from cfile
- details: model details data frame
- atol: see [solversettings](#)
- rtol: see [solversettings](#)
- ss\_atol: absolute tolerance to use when advancing to PK steady state
- ss\_rtol: relative tolerance to use when advancing to PK steady state
- maxsteps: see [solversettings](#)
- hmin: see [solversettings](#)
- hmax: see [solversettings](#)
- envir: the model environment
- plugins: plugins invoked in the model
- digits: number of digits to request in simulated data
- tscale: multiplicative scalar for time in results only
- mindt: simulation output time below which there model will assume to have not advanced
- preclean: logical indicating to clean up compilation artifacts prior to compiling
- debug: print debugging information during simulation run
- verbose: print extra information during setup for model run

---

as.list,mrgsims-method

*Coerce an mrgsims object to list*

---

### Description

Coerce an mrgsims object to list

### Usage

```
## S4 method for signature 'mrgsims'
as.list(x, ...)
```

### Arguments

|     |                   |
|-----|-------------------|
| x   | an mrgsims object |
| ... | not used          |

**Description**

These are simple functions that may be helpful to create the matrix objects that mrgsolve expects. Functions are named based on whether they create a diagonal matrix (d), a block matrix (b), or a correlation matrix (c).

**Usage**

```
as_bmat(x, ...)  
  
## S4 method for signature 'list'  
as_bmat(x, ...)  
  
## S4 method for signature 'numeric'  
as_bmat(x, pat = "*", ...)  
  
## S4 method for signature 'data.frame'  
as_bmat(x, pat = "*", cols = NULL, ...)  
  
## S4 method for signature 'ANY'  
as_bmat(x, ...)  
  
as_dmat(x, ...)  
  
## S4 method for signature 'list'  
as_dmat(x, ...)  
  
## S4 method for signature 'ANY'  
as_dmat(x, ...)  
  
## S4 method for signature 'numeric'  
as_dmat(x, pat = "*", ...)  
  
## S4 method for signature 'data.frame'  
as_dmat(x, pat = "*", cols = NULL, ...)  
  
as_cmat(x, ...)
```

**Arguments**

|      |  |
|------|--|
| x    | data frame or list   |
| ...  | arguments passed to <a href="#">dmat</a> or <a href="#">bmat</a> |
| pat  | regular expression, character                                    |
| cols | column names to use instead of pat                               |



## Details

Use `as_dmat` to create a diagonal matrix, `as_bmat` to create a block matrix, and `as_cmat` to create a block matrix where diagonal elements are understood to be correlations rather than covariances. `as_cmat` uses `as_bmat` to form the matrix and then converts off-diagonal elements to covariances before returning.

The methods for `data.frame` will work down the rows of the data frame and make the appropriate matrix from the data in each row. The result is a list of matrices.

## Value

A numeric matrix for list and numeric methods. For `data.frames`, a list of matrices are returned.

## See Also

[bmat](#), [dmat](#), [cmat](#)

## Examples

```
df <- data.frame(
  OMEGA1.1 = c(1,2),
  OMEGA2.1 = c(11,22),
  OMEGA2.2 = c(3,4),
  SIGMA1.1 = 1,
  FOO=-1
)

as_bmat(df, "OMEGA")
as_dmat(df, "SIGMA")
as_dmat(df[1,], "OMEGA")
```

---

as\_data\_set

*Create a simulation data set from ev objects*

---

## Description

Create a simulation data set from ev objects

## Usage

```
as_data_set(x, ...)
```

## S4 method for signature 'ev'

```
as_data_set(x, ...)
```

## S4 method for signature 'data.frame'

```
as_data_set(x, ...)
```

**Arguments**

x                    ev objects  
 ...                 more ev objects

**Details**

The goal is to take a series of event objects and combine them into a single data set that can be passed to `data_set`. Each event object is added to the data frame as an ID or set of IDs that are distinct from the IDs in the other event objects. Note that including ID argument to the `ev` call where `length(ID)` is greater than one will render that set of events for all of IDs that are requested.

To get a data frame with one row (event) per ID look at `expand.ev`.

**Value**

a data frame suitable for passing into `data_set`

**Examples**

```
as_data_set(ev(amt=c(100,200), cmt=1, ID=1:3),
            ev(amt=300, time=24, ID=1:2),
            ev(amt=1000, ii=8, addl=10, ID=1:3))

# Instead of this, use expand.ev
as_data_set(ev(amt=100), ev(amt=200),ev(amt=300))
```

---

 as\_deslist

---

*Create a list of designs from a data frame*


---

**Description**

Create a list of designs from a data frame

**Usage**

```
as_deslist(data, descol = "ID")
```

**Arguments**

data                input data set; see details  
 descol             character column name to be used for design groups

**Details**

The input data set must have a column with the same name as the value of `descol`. Other column names should be `start` (the time of the first observation), `end` (the time of the last observation), `delta` (the time steps to take between `start` and `end`), and `add` (other, ad-hoc times). Note that `add` might be a `list-column` to get a vector of times for each time grid object.

**Value**

The function returns a list of `tgrid` objects, one for each unique value found in `descol`.

**Examples**

```
idata <- tibble::tibble(ID=1:4, end=seq(24,96,24), delta=6,
  add=list(c(122,124,135),c(111), c(99),c(88)))
```

```
idata <- dplyr::mutate(idata, GRP = ID %%2)
```

```
idata
```

```
l <- as_deslist(idata,"GRP")
```

```
l
```

```
lapply(l,stime)
```

```
lapply(as_deslist(idata, "ID"),stime)
```

---

blocks

*Return the code blocks from a model specification file*

---

**Description**

Return the code blocks from a model specification file

**Usage**

```
blocks(x, ...)
```

```
## S4 method for signature 'mrgmod'
blocks(x, ...)
```

```
## S4 method for signature 'character'
blocks(x, ...)
```

**Arguments**

```
x          model object or path to model specification file
...        passed along
```

**Examples**

```
mod <- mrgsolve::house()
mod %>% blocks
mod %>% blocks(PARAM, TABLE)
```

---

**BLOCK\_PARSE***Functions to parse code blocks*

---

**Description**

Most of the basic blocks are listed in this help topic. But see also [PKMODEL](#) which has more-involved options and is documented separately.

**Usage**

```
PARAM(  
  x,  
  env,  
  annotated = FALSE,  
  covariates = FALSE,  
  pos = 1,  
  as_object = FALSE,  
  ...  
)
```

```
FIXED(x, env, annotated = FALSE, pos = 1, ...)
```

```
THETA(x, env, annotated = FALSE, pos = 1, name = "THETA", fill = NULL, ...)
```

```
INIT(x, env, annotated = FALSE, pos = 1, as_object = FALSE, ...)
```

```
CMT(x, env, annotated = FALSE, pos = 1, as_object = FALSE, ...)
```

```
CAPTURE(x, env, annotated = FALSE, pos = 1, ...)
```

**Arguments**

|            |  |
|------------|--|
| x          | data   |
| env        | parse environment                            |
| annotated  | logical                                      |
| covariates | logical                                      |
| pos        | block position                               |
| as_object  | indicates that object code is being provided |
| ...        | passed                                       |
| name       | block name                                   |
| fill       | data to use for block contents               |

**See Also**

[PKMODEL](#)

---

c,matlist-method      *Operations with matlist objects*

---

**Description**

Operations with matlist objects

**Usage**

```
## S4 method for signature 'matlist'
c(x, ..., recursive = FALSE)
```

**Arguments**

|           |                       |
|-----------|-----------------------|
| x         | a matlist object      |
| ...       | other matlist objects |
| recursive | not used              |

---

c,tgrid-method      *Operations with tgrid objects*

---

**Description**

Operations with tgrid objects

**Usage**

```
## S4 method for signature 'tgrid'
c(x, ..., recursive = FALSE)

## S4 method for signature 'tgrids'
c(x, ..., recursive = FALSE)

## S4 method for signature 'tgrid,numeric'
e1 + e2

## S4 method for signature 'tgrid,numeric'
e1 * e2

## S4 method for signature 'tgrids,numeric'
e1 + e2

## S4 method for signature 'tgrids,numeric'
e1 * e2
```

**Arguments**

|           |                               |
|-----------|-------------------------------|
| x         | mrgmod object                 |
| ...       | passed along to other methods |
| recursive | not used                      |
| e1        | tgrid or tgrids object        |
| e2        | numeric value                 |

---

|           |  |
|-----------|--|
| carry_out | <i>Select items to carry into simulated output</i> |
|-----------|--|

---

**Description**

When items named in this function are found in the input data set (either [data\\_set](#) or [idata\\_set](#)), they are copied into the simulated output. Special items like `evid` or `amt` or the like are not copied from the data set per se, but they are copied from `datarecord` objects that are created during the simulation.

**Usage**

```
carry_out(x, ...)
```

```
carry.out(x, ...)
```

**Arguments**

|     |              |
|-----|--------------|
| x   | model object |
| ... | passed along |

**Details**

There is also a `carry.out` argument to [mrgsim](#) that can be set to accomplish the same thing as a call to `carry_out` in the pipeline.

`carry.out` and `carry_out`. Using the underscore version is now preferred.

---

|      |   |
|------|---|
| cmtn | <i>Get the compartment number from a compartment name</i> |
|------|---|

---

**Description**

Get the compartment number from a compartment name

**Usage**

```
cmtn(x, ...)
```

```
## S4 method for signature 'mrgmod'  
cmtn(x, tag, ...)
```

**Arguments**

|     |                  |
|-----|------------------|
| x   | model object     |
| ... | passed along     |
| tag | compartment name |

**Examples**

```
mod <- mrgsolve::house()  
mod %>% cmtn("CENT")
```

---

|      |                                      |
|------|--------------------------------------|
| code | <i>Extract the code from a model</i> |
|------|--------------------------------------|

---

**Description**

Extract the code from a model

**Usage**

```
code(x)
```

**Arguments**

|   |                          |
|---|--------------------------|
| x | an mrgsolve model object |
|---|--------------------------|

**Value**

a character vector of model code

---

 data\_set

*Select and modify a data set for simulation*


---

### Description

The input data set (`data_set`) is a data frame that specifies observations, model events, and / or parameter values for a population of individuals.

### Usage

```
data_set(x, data, ...)

## S4 method for signature 'mrgmod,data.frame'
data_set(
  x,
  data,
  .subset = TRUE,
  .select = TRUE,
  object = NULL,
  need = NULL,
  ...
)

## S4 method for signature 'mrgmod,ANY'
data_set(x, data, ...)

## S4 method for signature 'mrgmod,ev'
data_set(x, data, ...)

## S4 method for signature 'mrgmod,missing'
data_set(x, object, ...)
```

### Arguments

|                      |   |
|----------------------|---|
| <code>x</code>       | model object  |
| <code>data</code>    | data set  |
| <code>...</code>     | passed along  |
| <code>.subset</code> | an unquoted expression passed to <code>dplyr::filter</code> ; retain only certain rows in the data set  |
| <code>.select</code> | passed to <code>dplyr::select</code> ; retain only certain columns in the data set; this should be the result of a call to <code>dplyr::vars()</code> |
| <code>object</code>  | character name of an object existing in <code>\$ENV</code> to use for the data set  |
| <code>need</code>    | passed to <a href="#">inventory</a>   |



## Details

Input data sets are R data frames that can include columns with any valid name, however columns with selected names are treated specially by `mrgsolve` and incorporated into the simulation.

ID specifies the subject ID and is required for every input data set.

When columns have the same name as parameters (`$PARAM` in the model specification file), the values in those columns will be used to update the corresponding parameter as the simulation progresses.

Input data set may include the following columns related to PK dosing events: `time`, `cmt`, `amt`, `rate`, `ii`, `addl`, `ss`. Along with ID, `time` is a required column in the input data set unless `$PRED` is in use. Upper case PK dosing column names including `TIME`, `CMT`, `AMT`, `RATE`, `II`, `ADDL`, `SS` are also recognized. However, an error will be generated if a mix of upper case and lower case columns in this family are found.

`time` is the observation or event time, `cmt` is the compartment number (see [init](#)), `amt` is the dosing amount, `rate` is the infusion rate, `ii` is the dosing interval, `addl` specifies additional doses to administer, and `ss` is a flag for steady state dosing. These column names operate similarly to other non-linear mixed effects modeling software.

An error will be generated when `mrgsolve` detects that the data set is not sorted by `time` within an individual. Also, an error will be generated in case `mrgsolve` finds negative values for `time`, unless `$PRED` is in use.

Only numeric data can be brought in to the problem. Any non-numeric data columns will be dropped with warning. See [numerics\\_only](#), which is used to prepare the data set.

An error will be generated if any parameter columns in the input data set contain NA. Likewise, and error will be generated if missing values are found in the following columns: ID, `time`/`TIME`, `rate`/`RATE`.

See [exdatasets](#) for different example data sets.

## See Also

[idata\\_set](#), [ev](#), [valid\\_data\\_set](#), [valid\\_idata\\_set](#)

## Examples

```
mod <- mrgsolve::house()

data <- expand.ev(ID=1:3, amt=c(10,20))

mod %>% data_set(data, ID > 1) %>% mrgsim

data(extran1)
head(extran1)

mod %>% data_set(extran1) %>% mrgsim
mod %>% mrgsim(data=extran1)
```

---

design                      *Set observation designs for the simulation*

---

### Description

This function also allows you to assign different designs to different groups or individuals in a population.

### Usage

```
design(x, deslist = list(), descol = character(0), ...)
```

### Arguments

|         |  |
|---------|--|
| x       | model object   |
| deslist | a list of tgrid or tgrids objects or numeric vector to be used in place of ... |
| descol  | the idata column name (character) for design assignment                        |
| ...     | not used   |

### Details

This setup requires the use of an `idata_set`, with individual-level data passed in one ID per row. For each ID, specify a grouping variable in `idata` (`descol`). For each unique value of the grouping variable, make one `tgrid` object and pass them in order as ... or form them into a list and pass as `deslist`.

You must assign the `idata_set` before assigning the designs in the command chain (see the example below).

### Examples

```
peak <- tgrid(0,6,0.1)
sparse <- tgrid(0,24,6)

des1 <- c(peak,sparse)
des2 <- tgrid(0,72,4)

data <- expand.ev(ID = 1:10, amt=c(100,300))
data$GRP <- data$amt/100

idata <- data[,c("ID", "amt")]

mod <- mrgsolve::house()

mod %>%
  omat(dmat(1,1,1,1)) %>%
  carry_out(GRP) %>%
```

```
idata_set(idata) %>%  
design(list(des1, des2), "amt") %>%  
data_set(data) %>%  
mrgsim %>%  
plot(RES~time|GRP)
```

---

details

*Extract model details*

---

## Description

Extract model details

## Usage

```
details(x, complete = FALSE, values = TRUE, ...)
```

## Arguments

|          |  |
|----------|--|
| x        | a model object   |
| complete | logical; if TRUE, un-annotated parameters and compartments will be added to the output |
| values   | logical; if TRUE, a values column will be added to the output                          |
| ...      | not used   |

## Details

This function is not exported. You will have to call it with `mrgsolve:::details()`.

## Examples

```
mod <- mrgsolve:::house()  
mrgsolve:::details(mod)
```

---

|         |  |
|---------|--|
| drop_re | <i>Deprecated: drop random effect matrices from model object</i> |
|---------|--|

---

**Description**

Deprecated: drop random effect matrices from model object

**Usage**

```
drop_re(.x, ...)
```

```
drop.re(...)
```

**Arguments**

|    |          |
|----|----------|
| .x | not used |
|----|----------|

|     |          |
|-----|----------|
| ... | not used |
|-----|----------|

**Details**

Users are no longer allowed to remove random effect matrices from the model object. Use [zero\\_re](#) instead to convert the matrix to all zeros.

**See Also**

[zero\\_re](#)

---

|          |  |
|----------|--|
| env_eval | <i>Re-evaluate the code in the ENV block</i> |
|----------|--|

---

**Description**

The \$ENV block is a block of R code that can realize any sort of R object that might be used in running a model.

**Usage**

```
env_eval(x, seed = NULL)
```

**Arguments**

|   |              |
|---|--------------|
| x | model object |
|---|--------------|

|      |   |
|------|---|
| seed | passed to <a href="#">set.seed</a> if a numeric value is supplied |
|------|---|

**See Also**

[env\\_get](#), [env\\_ls](#)

---

|         |                                 |
|---------|---------------------------------|
| env_get | <i>Return model environment</i> |
|---------|---------------------------------|

---

**Description**

Return model environment

**Usage**

```
env_get(x, tolist = TRUE)
```

```
env_get_env(x)
```

**Arguments**

x                    model object

tolist                should the environment be coerced to list?

---

|        |  |
|--------|--|
| env_ls | <i>List objects in the model environment</i> |
|--------|--|

---

**Description**

Each model keeps an internal environment that allows the user to carry any R object along. Objects are coded in \$ENV.

**Usage**

```
env_ls(x, ...)
```

**Arguments**

x                    model object

...                  passed to `ls`

---

|            |  |
|------------|--|
| env_update | <i>Update objects in model environment</i> |
|------------|--|

---

**Description**

Update objects in model environment

**Usage**

```
env_update(.x, ..., .dots = list())
```

**Arguments**

|       |                            |
|-------|----------------------------|
| .x    | model object               |
| ...   | objects to update          |
| .dots | list of objects to updated |

---

|    |  |
|----|--|
| ev | <i>Event objects for simulating PK and other interventions</i> |
|----|--|

---

**Description**

An event object specifies dosing or other interventions that get implemented during simulation. Event objects do similar things as [data\\_set](#), but simpler and quicker.

**Usage**

```
ev(x, ...)

## S4 method for signature 'mrgmod'
ev(x, object = NULL, ...)

## S4 method for signature 'missing'
ev(
  time = 0,
  amt = 0,
  evid = 1,
  cmt = 1,
  ID = numeric(0),
  replicate = TRUE,
  until = NULL,
  tinf = NULL,
  realize_addl = FALSE,
  ...
)
```

```
## S4 method for signature 'ev'
ev(x, realize_addl = FALSE, ...)
```

### Arguments

|              |  |
|--------------|--|
| x            | a model object   |
| ...          | other items to be incorporated into the event object; see details  |
| object       | passed to show   |
| time         | event time   |
| amt          | dose amount  |
| evid         | event ID   |
| cmt          | compartment  |
| ID           | subject ID   |
| replicate    | logical; if TRUE, events will be replicated for each individual in ID  |
| until        | the expected maximum <b>observation</b> time for this regimen  |
| tinf         | infusion time; if greater than zero, then the rate item will be derived as amt/tinf                                  |
| realize_addl | if FALSE (default), no change to addl doses. If TRUE, addl doses are made explicit with <a href="#">realize_addl</a> |

### Details

- Required items in events objects include `time`, `amt`, `evid` and `cmt`.
- If not supplied, `evid` is assumed to be 1.
- If not supplied, `cmt` is assumed to be 1.
- If not supplied, `time` is assumed to be 0.
- If `amt` is not supplied, an error will be generated.
- If `total` is supplied, then `addl` will be set to `total - 1`.
- Other items can include `ii`, `ss`, and `addl` (see [data\\_set](#) for details on all of these items).
- ID may be specified as a vector.
- If `replicate` is TRUE (default), then the events regimen is replicated for each ID; otherwise, the number of event rows must match the number of IDs entered

### Value

events object

### See Also

[ev\\_rep](#), [ev\\_days](#), [ev\\_repeat](#), [ev\\_assign](#), [ev\\_seq](#), [mutate.ev](#), [as.ev](#), [ev\\_methods](#)

**Examples**

```

mod <- mrgsolve::house()

mod <- mod %>% ev(amt=1000, time=0, cmt=1)

loading <- ev(time=0, cmt=1, amt=1000)

maint <- ev(time=12, cmt=1, amt=500, ii=12, addl=10)

c(loading, maint)

loading$time

```

---

ev\_assign

*Replicate a list of events into a data set*


---

**Description**

Replicate a list of events into a data set

**Usage**

```

ev_assign(l, idata, evgroup, join = FALSE)

assign_ev(...)

```

**Arguments**

|         |  |
|---------|--|
| l       | list of event objects  |
| idata   | an idata set (one ID per row)  |
| evgroup | the character name of the column in idata that specifies event object to implement |
| join    | if TRUE, join idata to the data set before returning.                              |
| ...     | used to pass arguments from assign_ev to ev_assign                                 |

**Details**

ev\_assign connects events in a list passed in as the l argument to values in the data set identified in the evgroup argument. For making assignments, the unique values in the evgroup column are first sorted so that the first sorted unique value in evgroup is assigned to the first event in l, the second sorted value in evgroup column is assigned to the second event in l, and so on. This is a change from previous behavior, which did not sort the unique values in evgroup prior to making the assignments.



**Examples**

```

ev1 <- ev(amt=100)
ev2 <- ev(amt=300, rate=100, ii=12, addl=10)

idata <- data.frame(ID=1:10)
idata$arm <- 1+(idata$ID %%2)

ev_assign(list(ev1, ev2), idata, "arm", join=TRUE)

```

---

ev\_days

*Schedule dosing events on days of the week*


---

**Description**

This function lets you schedule doses on specific days of the week, allowing you to create dosing regimens on Monday/Wednesday/Friday, or Tuesday/Thursday, or every other day (however you want to define that) etc.

**Usage**

```

ev_days(
  ev = NULL,
  days = "",
  addl = 0,
  ii = 168,
  unit = c("hours", "days"),
  ...
)

```

**Arguments**

|      |  |
|------|--|
| ev   | an event object  |
| days | comma- or space-separated character string of valid days of the the week (see details) |
| addl | additional doses to administer   |
| ii   | inter-dose interval; intended use is to keep this at the default value                 |
| unit | time unit; the function can only currently handle hours or days                        |
| ...  | event objects named by one the valid days of the week (see details)                    |

**Details**

Valid names of the week are:

- m for Monday
- t for Tuesday
- w for Wednesday
- th for Thursday
- f for Friday
- sa for Saturday
- s for Sunday

The whole purpose of this function is to schedule doses on specific days of the week, in a repeating weekly schedule. Please do use caution when changing `ii` from its default value.

**Examples**

```
# Monday, Wednesday, Friday x 4 weeks
ev_days(ev(amt=100), days="m,w,f", addl=3)

# 50 mg Tuesdays, 100 mg Thursdays x 6 months
ev_days(t=ev(amt=50), th=ev(amt=100), addl=23)
```

---

 ev\_rep

*Replicate an event object*


---

**Description**

An event sequence can be replicated a certain number of times in a certain number of IDs.

**Usage**

```
ev_rep(x, ID = 1, n = NULL, wait = 0, as.ev = FALSE, id = NULL)
```

**Arguments**

|       |                                     |
|-------|-------------------------------------|
| x     | event object                        |
| ID    | numeric vector if IDs               |
| n     | passed to <a href="#">ev_repeat</a> |
| wait  | passed to <a href="#">ev_repeat</a> |
| as.ev | if TRUE an event object is returned |
| id    | deprecated; use ID instead          |

**Value**

A single data.frame or event object as determined by the value of `as.ev`.

**See Also**

[ev\\_repeat](#)

**Examples**

```
e1 <- c(ev(amt=100), ev(amt=200, ii=24, addl=2, time=72))
ev_rep(e1, 1:5)
```

---

|           |  |
|-----------|--|
| ev_repeat | <i>Repeat a block of dosing events</i> |
|-----------|--|

---

**Description**

Repeat a block of dosing events

**Usage**

```
ev_repeat(x, n, wait = 0, as.ev = FALSE)
```

**Arguments**

|       |  |
|-------|--|
| x     | event object or dosing data frame  |
| n     | number of times to repeat  |
| wait  | time to wait between repeats   |
| as.ev | if TRUE, an event object is returned; otherwise a data.frame is returned |

**Value**

See `as.ev` argument.

---

 ev\_rx

*Create intervention objects from Rx input*


---

### Description

See details below for Rx specification. Actual parsing is done by [parse\\_rx](#); this function can be used to debug Rx inputs.

### Usage

```
ev_rx(x, y, ...)
```

```
## S4 method for signature 'mrgmod,character'
```

```
ev_rx(x, y, ...)
```

```
## S4 method for signature 'character,missing'
```

```
ev_rx(x, df = FALSE, ...)
```

```
parse_rx(x)
```

### Arguments

|     |                                       |
|-----|---------------------------------------|
| x   | a model object or character Rx input  |
| y   | character Rx input; see details       |
| ... | not used at this time                 |
| df  | if TRUE then a data frame is returned |

### Value

The method dispatched on model object (mrgmod) returns another model object. The character method returns an event object. The `parse_rx` function return a list named with arguments for the event object constructor [ev](#).

### Rx specification

- The dose is found at the start of the string by sequential digits; this may be integer, decimal, or in scientific notation
- Use `in` to identify the dosing compartment number; must be integer
- Use `q` to identify the dosing interval; must be integer or decimal number (but not scientific notation)
- Use `over` to indicate an infusion and its duration; integer or decimal number
- Use `x` to indicate total number of doses; must be integer
- Use `then` or `,` to separate dosing periods
- User `af`ter to insert a lag in the start of a period; integer or decimal number (but not scientific notation)

**Examples**

```

# example("ev_rx")

ev_rx("100")

ev_rx("100 in 2")

ev_rx("100 q12 x 3")

ev_rx("100 over 2")

ev_rx("100 q 24 x 3 then 50 q12 x 2")

ev_rx("100 then 50 q 24 after 12")

ev_rx("100.2E-2 q4")

ev_rx("100 over 2.23")

ev_rx("100 q 12 x 3")

parse_rx("100 mg q 24 then 200 mg q12")

```

---

ev\_seq

*Schedule a series of event objects*


---

**Description**

Schedule a series of event objects

**Usage**

```

ev_seq(..., ID = NULL, .dots = NULL, id = NULL)

## S3 method for class 'ev'
seq(...)

```

**Arguments**

|       |   |
|-------|---|
| ...   | event objects or numeric arguments named wait |
| ID    | numeric vector of subject IDs                 |
| .dots | a list of event objects that replaces ...     |
| id    | deprecated; use ID                            |

## Details

The doses for the next event line start after all of the doses from the previous event line plus one dosing interval from the previous event line (see examples).

When numerics named `wait` are mixed in with the event objects, a period with no dosing activity is incorporated into the sequence, between the adjacent dosing event objects. Values for `wait` can be negative.

Values for `time` in any event object act like a prefix time spacer wherever that event occurs in the event sequence (see examples).

Use the generic `seq` when the first argument is an event object. If a waiting period is the first event, you will need to use `ev_seq`. When an event object has multiple rows, the end time for that sequence is taken to be one dosing interval after the event that takes place on the last row of the event object.

## Value

A single event object.

## Examples

```
e1 <- ev(amt=100, ii=12, addl=1)
e2 <- ev(amt=200)
seq(e1, e2)
seq(e1, .ii = 8, e2)
seq(e1, wait = 8, e2)
seq(e1, .ii = 8, e2, ID = 1:10)
ev_seq(.ii = 12, e1, .ii = 120, e2, .ii = 120, e1)
seq(ev(amt=100, ii=12), ev(time=8, amt=200))
```

---

exdatasets

*Example input data sets*

---

## Description

Example input data sets

**Usage**

```
data(exidata)
```

```
data(extran1)
```

```
data(extran2)
```

```
data(extran3)
```

```
data(exTheoph)
```

```
data(exBoot)
```

**Details**

- exidata holds individual-level parameters and other data items, one per row
- extran1 is a "condensed" data set
- extran2 is a full dataset
- extran3 is a full dataset with parameters
- exTheoph is the theophylline data set, ready for input into mrgsolve
- exBoot a set of bootstrap parameter estimates

**Examples**

```
mod <- mrgsolve::house() %>% update(end=240) %>% Req(CP)

## Full data set
data(exTheoph)
out <- mod %>% data_set(exTheoph) %>% mrgsim
out
plot(out)

## Condensed: mrgsolve fills in the observations
data(extran1)
out <- mod %>% data_set(extran1) %>% mrgsim
out
plot(out)

## Add a parameter to the data set
stopifnot(require(dplyr))
data <- extran1 %>% distinct(ID) %>% select(ID) %>%
  mutate(CL=exp(log(1.5) + rnorm(nrow(.), 0, sqrt(0.1)))) %>%
  left_join(extran1,.)

data

out <- mod %>% data_set(data) %>% carry_out(CL) %>% mrgsim
out
```

```
plot(out)

## idata
data(exidata)
out <- mod %>% idata_set(exidata) %>% ev(amt=100,ii=24,addl=10) %>% mrgsim
plot(out, CP~time|ID)
```

---

expand.idata

*Create template data sets for simulation*

---

## Description

Create template data sets for simulation

## Usage

```
expand.idata(...)
```

```
expand.ev(...)
```

```
ev_expand(...)
```

## Arguments

... passed to [expand.grid](#)

## Details

An ID column is added as `seq(nrow(ans))` if not supplied by the user. For `expand.ev`, defaults also added include `cmt = 1`, `time = 0`, `evid = 1`. If `total` is included, then `addl` is derived as `total - 1`. If `tinf` is included, then an infusion rate is derived for row where `tinf` is greater than zero.

## Examples

```
idata <- expand.idata(CL = c(1,2,3), VC = c(10,20,30))
```

```
doses <- expand.ev(amt = c(300,100), ii = c(12,24), cmt = 1)
```

```
infusion <- expand.ev(amt = 100, tinf = 2)
```



---

expand\_observations     *Insert observations into a data set*

---

**Description**

Insert observations into a data set

**Usage**

```
expand_observations(data, times, unique = FALSE, obs_pos = -1L)
```

**Arguments**

|         |   |
|---------|---|
| data    | a data set or event object  |
| times   | a vector of observation times   |
| unique  | 'logical'; if 'TRUE' then values for 'time' are dropped if they are found anywhere in 'data'  |
| obs_pos | determines sorting order for observations; use '-1' (default) to put observations first; otherwise, use large integer to ensure observations are placed after doses |

**Details**

Non-numeric columns will be dropped with a warning.

**Value**

A data frame

**Examples**

```
data <- expand.ev(amt = c(100,200,300))  
expand_observations(data, times = seq(0,48,2))
```

---

idata\_set     *Select and modify a idata set for simulation*

---

**Description**

The individual data set (idata\_set) is a data frame with one row for each individual in a population, specifying parameters and other individual-level data.

**Usage**

```

idata_set(x, data, ...)

## S4 method for signature 'mrgmod,data.frame'
idata_set(
  x,
  data,
  .subset = TRUE,
  .select = TRUE,
  object = NULL,
  need = NULL,
  ...
)

## S4 method for signature 'mrgmod,ANY'
idata_set(x, data, ...)

## S4 method for signature 'mrgmod,missing'
idata_set(x, object, ...)

```

**Arguments**

|         |   |
|---------|---|
| x       | model object  |
| data    | a data set that can be coerced to data.frame  |
| ...     | passed along  |
| .subset | an unquoted expression passed to <code>dplyr::filter</code> ; retain only certain rows in the data set  |
| .select | passed to <code>dplyr::select</code> ; retain only certain columns in the data set; this should be the result of a call to <code>dplyr::vars()</code> |
| object  | character name of an object existing in <code>\$ENV</code> to use for the data set  |
| need    | passed to <a href="#">inventory</a>   |

**Details**

The `idata_set` is a `data.frame` that specifies individual-level data for the problem. An ID column is required and there can be no more than one row in the data frame for each individual.

In most cases, the columns in the `idata_set` have the same names as parameters in the [param](#) list. When this is the case, the parameter set is updated as the simulation proceeds once at the start of each individual. The ‘`idata_set`’ can also be used to set initial conditions for each individual: for a compartment called CMT, make a column in `idata_set` called `CMT_0` and make the value the desired initial value for that compartment. Note that this initial condition will be over-ridden if you also set the `CMT_0` in `$MAIN`.

The most common application of `idata_set` is to specify a population or batch of simulations to do. We commonly use `idata_set` with an event object (see [ev](#)). In that case, the event gets applied to each individual in the [idata\\_set](#).

It is also possible to provide both a `data_set` and a `idata_set`. In this case, the `idata_set` is used as a parameter lookup for IDs found in the `data_set`. Remember in this case, it is the `data_set` (not the `idata_set`) that determines the number of individuals in the simulation.

An error will be generated if any parameter columns in the input `idata` set contain NA.

### See Also

[data\\_set](#), [ev](#)

### Examples

```
mod <- mrgsolve::house()

data(exidata)

exidata

mod %>%
  idata_set(exidata, ID <= 2) %>%
  ev(amt = 100) %>%
  mrgsim() %>%
  plot()

mod %>%
  idata_set(exidata) %>%
  ev(amt = 100) %>%
  mrgsim()

mod %>% ev(amt = 100) %>% mrgsim(idata=exidata)
```

---

init

*Methods for working with the model compartment list*

---

### Description

Calling `init` with the model object as the first argument will return the model initial conditions as a `numericlist` object. See [numericlist](#) for methods to deal with `cmt_list` objects.

### Usage

```
init(.x, ...)
```

## S4 method for signature 'mrgmod'

```
init(.x, .y = list(), ..., .pat = "*")
```

## S4 method for signature 'mrgsims'

```
init(.x, ...)
```

```
## S4 method for signature 'missing'
init(.x, ...)

## S4 method for signature 'list'
init(.x, ...)

## S4 method for signature 'ANY'
init(.x, ...)
```

### Arguments

|                   |   |
|-------------------|---|
| <code>.x</code>   | the model object  |
| <code>...</code>  | passed along  |
| <code>.y</code>   | list to be merged into parameter list   |
| <code>.pat</code> | a regular expression (character) to be applied as a filter when printing compartments to the screen |

### Details

Can be used to either get a compartment list object from a `mrgmod` model object or to update the compartment initial conditions in a model object. For both uses, the return value is a `cmt_list` object. For the former use, `init` is usually called to print the compartment initial conditions to the screen, but the `cmt_list` object can also be coerced to a list or numeric R object.

### Value

an object of class `cmt_list` (see [numericlist](#))

### Examples

```
## example("init")
mod <- mrgsolve::house()

init(mod)
init(mod, .pat="^C") ## may be useful for large models

class(init(mod))

init(mod)$CENT

as.list(init(mod))
as.data.frame(init(mod))
```

---

|           |   |
|-----------|---|
| inventory | <i>Check whether all required parameters needed in a model are present in an object</i> |
|-----------|---|

---

## Description

Check whether all required parameters needed in a model are present in an object

## Usage

```
inventory(x, obj, ..., .strict = FALSE)
```

## Arguments

|         |  |
|---------|--|
| x       | model object   |
| obj     | data.frame to pass to <a href="#">idata_set</a> or <a href="#">data_set</a>                        |
| ...     | capture dplyr-style parameter requirements   |
| .strict | whether to stop execution if all requirements are present (TRUE) or just warn (FALSE); see details |

## Details

If parameter requirements are not explicitly stated, the requirement defaults to all parameter names in x. Note that, by default, the inventory is not `.strict` unless the user explicitly states the parameter requirement. That is, if parameter requirements are explicitly stated, `.strict` will be set to TRUE if a value `.strict` was not passed in the call.

## Value

original mrgmod

## Examples

```
## Not run:
inventory(mod, idata, CL:V) # parameters defined, inclusively, CL through Volume
inventory(mod, idata, everything()) # all parameters
inventory(mod, idata, contains("OCC")) # all parameters containing OCC
inventory(mod, idata, -F) # all parameters except F

## End(Not run)
```

---

`is.mrgmod`*Check if an object is a model object*

---

**Description**

The function checks to see if the object is either mrgmod or packmod.

**Usage**`is.mrgmod(x)`**Arguments**

x                    any object

**Value**

TRUE if x inherits mrgsims.

---

`is.mrgsims`*Check if an object is mrgsim output*

---

**Description**

Check if an object is mrgsim output

**Usage**`is.mrgsims(x)`**Arguments**

x                    any object

**Value**

TRUE if x inherits mrgsims.

---

|        |  |
|--------|--|
| lctran | <i>Convert select upper case column names to lower case to conform to mrgsolve data expectations</i> |
|--------|--|

---

**Description**

Convert select upper case column names to lower case to conform to mrgsolve data expectations

**Usage**

```
lctran(data)
```

**Arguments**

data            an nmtran-like data frame

**Details**

Columns that will be renamed with lower case versions: AMT, II, SS, CMT, ADDL, RATE, EVID, TIME. If a lower case version of these names exist in the data set, the column will not be renamed.

**Value**

A data.frame with renamed columns

---

|        |                                     |
|--------|-------------------------------------|
| loadso | <i>Load the model shared object</i> |
|--------|-------------------------------------|

---

**Description**

Once the model is compiled, the model object can be used to re-load the model shared object (the compiled code underlying the mode) when the simulation is to be done in a different R process.

**Usage**

```
loadso(x, ...)
```

```
## S3 method for class 'mrgmod'
loadso(x, ...)
```

**Arguments**

x            the model object  
 ...         not used

## Details

The 'loadso' function most frequently needs to be used when parallelizing simulations across worker nodes. The model can be run after calling 'loadso', without requiring that it is re-compiled on worker nodes. It is likely required that the model is built (and the shared object stored) in a local directory off of the working R directory (see the second example).

## Value

The model object (invisibly).

## Examples

```
## Not run:
mod <- mread("pk1", modlib())
loadso(mod)

mod2 <- mread("pk2", modlib(), soloc = "build")
loadso(mod2)

## End(Not run)
```

---

matrix\_helpers

*Create matrices from vector input*

---

## Description

Create matrices from vector input

## Usage

```
bmat(..., correlation = FALSE, digits = -1)
```

```
cmat(..., digits = -1)
```

```
dmat(...)
```

## Arguments

|             |   |
|-------------|---|
| ...         | matrix data   |
| correlation | logical; if TRUE, off diagonal elements are assumed to be correlations and converted to covariances |
| digits      | if greater than zero, matrix is passed to signif (along with digits) prior to returning             |



**Details**

bmat makes a block matrix. cmatrix makes a correlation matrix. dmat makes a diagonal matrix.

**See Also**

[as\\_bmat](#)

[as\\_dmat](#)

**Examples**

```
dmat(1,2,3)/10
```

```
bmat(0.5,0.01,0.2)
```

```
cmatrix(0.5, 0.87,0.2)
```

---

mcode

*Write, compile, and load model code*

---

**Description**

This is a convenience function that ultimately calls [mread](#). Model code is written to a file and read back in using [mread](#).

**Usage**

```
mcode(model, code, project = getOption("mrgsolve.project", tempdir()), ...)
```

```
mcode_cache(  
  model,  
  code,  
  project = getOption("mrgsolve.project", tempdir()),  
  ...  
)
```

**Arguments**

|         |   |
|---------|---|
| model   | model name  |
| code    | character string specifying a mrgsolve model  |
| project | project name  |
| ...     | passed to <a href="#">mread</a> ; see that help topic for other arguments that can be set |

### Details

Note that the arguments are in slightly different order than `mread`. The default project is `tempdir()`.

See the `mread` help topic for discussion about caching compilation results with `mcode_cache`.

### See Also

[mread](#), [mread\\_cache](#)

### Examples

```
## Not run:
code <- '
$CMT DEPOT CENT
$PKMODEL ncmt=1, depot=TRUE
$MAIN
double CL = 1;
double V = 20;
double KA = 1;
'

mod <- mcode("example",code)

## End(Not run)
```

---

mcRNG

*Set RNG to use L'Ecuyer-CMRG*

---

### Description

Set RNG to use L'Ecuyer-CMRG

### Usage

`mcRNG()`

---

|        |                               |
|--------|-------------------------------|
| modlib | <i>Internal model library</i> |
|--------|-------------------------------|

---

## Description

Internal model library

## Usage

```
modlib(model = NULL, ..., list = FALSE)
```

## Arguments

|       |  |
|-------|--|
| model | character name of a model in the library |
| ...   | passed to <a href="#">mread_cache</a>    |
| list  | list available models                    |

## Details

See [modlib\\_details](#), [modlib\\_pk](#), [modlib\\_pkpd](#), [modlib\\_tmdd](#), [modlib\\_viral](#) for details.

Call `modlib("<modelname>")` to compile and load a mode from the library.

Call `modlib(list=TRUE)` to list available models. Once the model is loaded (see examples below), call `as.list(mod)$code` to see model code and equations.

## Examples

```
## Not run:
mod <- mread("pk1cmt", modlib())
mod <- mread("pk2cmt", modlib())
mod <- mread("pk3cmt", modlib())
mod <- mread("pk1", modlib())
mod <- mread("pk2", modlib())
mod <- mread("popex", modlib())
mod <- mread("irm1", modlib())
mod <- mread("irm2", modlib())
mod <- mread("irm3", modlib())
mod <- mread("irm4", modlib())
mod <- mread("emax", modlib())
mod <- mread("effect", modlib())
mod <- mread("tmdd", modlib())
mod <- mread("viral1", modlib())
mod <- mread("viral2", modlib())
mod <- mread("pred1", modlib())
mod <- mread("pbpk", modlib())
mod <- mread("1005", modlib()) # embedded NONMEM result

mrgsolve:::code(mod)
```

```
## End(Not run)
```

---

|                |   |
|----------------|---|
| modlib_details | <i>modlib: PK/PD Model parameters, compartments, and output variables</i> |
|----------------|---|

---

## Description

modlib: PK/PD Model parameters, compartments, and output variables

## Compartments

- EV1, EV2: extravascular dosing compartments
- CENT: central PK compartment
- PERIPH: peripheral PK compartment
- PERIPH2: peripheral PK compartment 2
- RESP: response PD compartment (irm models)

## Output variables

- CP: concentration in the central compartment (CENT/VC)
- RESP: response (emax model)

## PK parameters

- KA1, KA2: first order absorption rate constants from first and second extravascular compartment (1/time)
- CL: clearance (volume/time)
- VC: volume of distribution, central compartment (volume)
- VP: volume of distribution, peripheral compartment (volume)
- VP2: volume of distribution, peripheral compartment 2 (volume)
- Q: intercompartmental clearance (volume/time)
- Q2: intercompartmental clearance 2 (volume/time)
- VMAX: maximum rate, nonlinear process (mass/time)
- KM: Michaelis constant (mass/volume)
- K10: elimination rate constant (1/time); CL/VC
- K12: rate constant for transfer to peripheral compartment from central (1/time); Q/VC
- K21: rate constant for transfer to central compartment from peripheral (1/time); Q/VP

**PD parameters**

- E0: baseline effect (emax model)
- EMAX, IMAX: maximum effect (response)
- EC50, IC50: concentration producing 50 percent of effect (mass/volume)
- KIN: zero-order response production rate (irm models) (response/time)
- KOUT: first-order response elimination rate (irm models) (1/time)
- n: sigmoidicity factor
- KE0: rate constant for transfer to effect compartment (1/time)

---

 modlib\_pk

*modlib: Pharmacokinetic models*


---

**Description**

modlib: Pharmacokinetic models

**Arguments**

... passed to update

**Details**

See [modlib\\_details](#) for more detailed descriptions of parameters and compartments.

The pk1cmt model is parameterized in terms of CL, VC, KA1 and KA2 and uses compartments EV1, EV2, and CENT. The pk2cmt model adds a PERIPH compartment and parameters Q and VP to that of the one-compartment model. Likewise, the three-compartment model (pk3cmt) adds PERIPH2 and parameters Q2 and VP2 to that of the two-compartment models. All pk models also have parameters VMAX (defaulting to zero, no non-linear clearance) and KM.

**Value**

an object of class packmod

**Model description**

All pk models have two extravascular dosing compartments and potential for linear and nonlinear clearance.

- pk1cmt: one compartment pk model using ODEs
- pk2cmt: two compartment pk model using ODEs
- pk3cmt: three compartment pk model using ODEs
- pk1: one compartment pk model in closed-form
- pk2: two compartment pk model in closed-form
- popex: a simple population pk model

---

 modlib\_pkpd

*modlib: Pharmacokinetic / pharmacodynamic models*


---

### Description

modlib: Pharmacokinetic / pharmacodynamic models

### Details

See [modlib\\_details](#) for more detailed descriptions of parameters and compartments.

All PK/PD models include 2-compartment PK model with absorption from 2 extravascular compartments and linear + nonlinear clearance. The PK models are parameterized with CL, VC, Q, VMAX, KM, KA1 and KA2 and implement compartments EV1, EV2, CENT, PERIPH . The indirect response models have compartment RESP and the emax model has output variable RESP. PD parameters include KIN, KOUT, IC50, EC50, IMAX, EMAX, E0, and n.

Also, once the model is loaded, use [see](#) method for mrgmod to view the model code.

### Model description

- irm1 inhibition of response production
- irm2 inhibition of response loss
- irm3 stimulation of response production
- irm4 stimulation of response loss
- pd\_effect effect compartment model
- emax sigmoid emax model

---

 modlib\_tmdd

*modlib: Target mediated disposition model*


---

### Description

modlib: Target mediated disposition model

### Arguments

... passed to update

**Parameters**

- KEL: elimination rate constant
- KTP: tissue to plasma rate constant
- KPT: plasma to tissue rate constant
- VC: volume of distribution
- KA1, KA2: absorption rate constants
- KINT: internalization rate constant
- KON: association rate constant
- KOFF: dissociation rate constant
- KSYN: target synthesis rate
- KDEG: target degradation rate constant

**Compartments**

- CENT: unbound drug in central compartment
- TISS: unbound drug in tissue compartment
- REC: concentration of target
- RC: concentration of drug-target complex
- EV1, EV2: extravascular dosing compartments

**Output variables**

- CP: unbound drug in the central compartment
- TOTAL: total concentration of target (complexed and uncomplexed)

---

modlib\_viral

*modlib: HCV viral dynamics models*

---

**Description**

modlib: HCV viral dynamics models

**Models**

- viral1: viral dynamics model with single HCV species
- viral2: viral dynamics model with wild-type and mutant HCV species

### Parameters

- s: new hepatocyte synthesis rate (cells/ml/day)
- d: hepatocyte death rate constant (1/day)
- p: viral production rate constant (copies/cell/day)
- beta: new infection rate constant (ml/copy/day)
- delta: infected cell death rate constant (1/day)
- c: viral clearance rate constant (1/day)
- fit: mutant virus fitness
- N: non-target hepatocytes
- mu: forward mutation rate
- Tmax: maximum number of target hepatocytes (cells/ml)
- rho: maximum hepatocyte regeneration rate (1/day)

### Compartments

- T: uninfected target hepatocytes (cells/ml)
- I: productively infected hepatocytes (cells/ml)
- V: hepatitis C virus (copies/ml)
- IM: mutant infected hepatocytes (cells/ml)
- VM: mutant hepatitis C virus (copies/ml)
- expos: exposure metric to drive pharmacodynamic model

---

mread

*Read a model specification file*

---

### Description

mread reads and parses the mrgsolve model specification file, builds the model, and returns a model object for simulation. mread\_cache does the same, but caches the compilation result for later use.

### Usage

```
mread(  
  model,  
  project = getOption("mrgsolve.project", getwd()),  
  code = NULL,  
  file = NULL,  
  udll = TRUE,  
  ignore.stdout = TRUE,  
  raw = FALSE,  
  compile = TRUE,  
  audit = TRUE,
```



```

    quiet = getOption("mrgsolve_mread_quiet", FALSE),
    check.bounds = FALSE,
    warn = TRUE,
    soloc = getOption("mrgsolve.soloc", tempdir()),
    preclean = FALSE,
    recover = FALSE,
    ...
)

mread_cache(
  model = NULL,
  project = getOption("mrgsolve.project", getwd()),
  file = paste0(model, ".cpp"),
  code = NULL,
  soloc = getOption("mrgsolve.soloc", tempdir()),
  quiet = FALSE,
  preclean = FALSE,
  ...
)

mread_file(file, ...)

```

### Arguments

|               |   |
|---------------|---|
| model         | model name  |
| project       | location of the model specification file and any headers to be included; see also the discussion about model; this argument can be set via options() library under details as well as the <a href="#">modlib</a> help topic |
| code          | a character string with model specification code to be used instead of a model file   |
| file          | the full file name (with extension, but without path) where the model is specified  |
| udll          | use unique name for shared object   |
| ignore.stdout | passed to system call for compiling model   |
| raw           | if TRUE, return a list of raw output  |
| compile       | logical; if TRUE, the model will be built   |
| audit         | check the model specification file for errors   |
| quiet         | don't print messages when compiling   |
| check.bounds  | check boundaries of parameter list  |
| warn          | logical; if TRUE, print warning messages that may arise   |
| soloc         | the directory location where the model shared object is built and stored; see details; this argument can be set via options(); if the directory does not exist, 'mread' will attempt to create it.                          |
| preclean      | logical; if TRUE, compilation artifacts are cleaned up first  |
| recover       | if TRUE, an object will be returned in case the model shared object fails to build  |
| ...           | passed to <a href="#">update</a> ; also arguments passed to mread from <a href="#">mread_cache</a> .  |

## Details

The `model` argument is required. For typical use, the `file` argument is omitted and the value for `file` is generated from the value for `model`. To determine the source file name, `mrgsolve` will look for a file extension in `model`. A file extension is assumed when it finds a period followed by one to three alpha-numeric characters at the end of the string (e.g. `mymodel.txt` but not `my.model`). If no file extension is found, the extension `.cpp` is assumed (e.g. `file` is `<model-name>.cpp`). If a file extension is found, `file` is `<model-name>`.

Best practice is to avoid using `.` in `model` unless you are using `model` to point to the model specification file name. Otherwise, use [mread\\_file](#).

Use the `soloc` argument to specify a directory location for building the model. This is the location where the model shared object will be stored on disk. The default is a temporary directory, so compilation artifacts are lost when R restarts when the default is used. Changing `soloc` to a persistent directory location will preserve those artifacts across R restarts. Also, if simulation from a single model is being done in separate processes on separate compute nodes, it might be necessary to store these compilation artifacts in a local directory to make them accessible to the different nodes. If the `soloc` directory does not exist, `mread` will attempt to create it.

Similarly, using `mread_cache` will cache results in the temporary directory and the cache cannot be accessed after the R process is restarted.

## Model Library

`mrgsolve` comes bundled with several precoded PK, PK/PD, and other systems models that are accessible via the `mread` interface.

Models available in the library include:

- PK models: `pk1cmt`, `pk2cmt`, `pk3cmt`, `pk1`, `pk2`, `popex`, `tddd`
- PKPD models: `irm1`, `irm2`, `irm3`, `irm4`, `emax`, `effect`
- Other models: `viral1`, `viral2`

When the library model is accessed, `mrgsolve` will compile and load the model as you would for any other model. It is only necessary to reference the correct model name and point the project argument to the `mrgsolve` model library location via [modlib](#).

For more details, see [modlib\\_pk](#), [modlib\\_pkpd](#), [modlib\\_tddd](#), [modlib\\_viral](#), and [modlib\\_details](#) for more information about the state variables and parameters in each model.

## See Also

[mcode](#), [mcode\\_cache](#)

## Examples

```
## Not run:
code <- '
$PARAM CL = 1, VC = 5
$CMT CENT
$ODE dxdt_CENT = -(CL/VC)*CENT;
```

```
,  
  
mod <- mcode("ex_mread", code)  
  
mod  
  
mod %>% init(CENT=1000) %>% mrgsim %>% plot  
  
mod <- mread("irm3", modlib())  
  
mod  
  
# if the model is in the file mymodel.cpp  
mod <- mread("mymodel")  
  
# if the model is in the file mymodel.txt  
mod <- mread(file = "mymodel.txt")  
  
or  
  
mod <- mread_file("mymodel.txt")  
  
## End(Not run)
```

---

mrgsim

*Simulate from a model object*

---

## Description

This function sets up the simulation run from data stored in the model object as well as arguments passed in. Use [mrgsim\\_q](#) instead to benchmark mrgsolve or to do repeated quick simulation for tasks like parameter optimization, sensitivity analyses, or optimal design. See [mrgsim\\_variants](#) for other mrgsim-like functions that have more focused inputs. `mrgsim_df` coerces output to `data.frame` prior to returning.

## Usage

```
mrgsim(x, data = NULL, idata = NULL, events = NULL, nid = 1, ...)
```

```
mrgsim_df(..., output = "df")
```

```
do_mrgsim(  
  x,  
  data,  
  idata = no_idata_set(),
```

```

carry_out = carry.out,
carry.out = character(0),
recover = character(0),
seed = as.integer(NA),
Request = character(0),
output = NULL,
capture = NULL,
obsonly = FALSE,
obsaug = FALSE,
tgrid = NULL,
recsort = 1,
deslist = list(),
descol = character(0),
filbak = TRUE,
tad = FALSE,
nocb = TRUE,
skip_init_calc = FALSE,
ss_n = 500,
ss_fixed = FALSE,
...
)

```

### Arguments

|           |   |
|-----------|---|
| x         | the model object  |
| data      | NMTRAN-like data set (see <a href="#">data_set</a> )  |
| idata     | a matrix or data frame of model parameters, one parameter per row (see <a href="#">idata_set</a> )  |
| events    | an event object   |
| nid       | integer number of individuals to simulate; only used if idata and data are missing  |
| ...       | passed to <a href="#">update</a> and <a href="#">do_mrgsim</a>  |
| output    | if NULL (the default) a mrgsim object is returned; otherwise, pass <code>df</code> to return a <code>data.frame</code> or <code>matrix</code> to return a matrix  |
| carry_out | numeric data items to copy into the output  |
| carry.out | soon to be deprecated; use <code>carry_out</code> instead   |
| recover   | character column names in either data or idata to join back (recover) to simulated data; may be any class (e.g. numeric, character, factor, etc)  |
| seed      | deprecated  |
| Request   | compartments or captured variables to retain in the simulated output; this is different than the request slot in the model object, which refers only to model compartments  |
| capture   | character file name used for debugging (not related to <code>\$CAPTURE</code> )   |
| obsonly   | if TRUE, dosing records are not included in the output  |
| obsaug    | augment the data set with time grid observations; when TRUE and a full data set is used, the simulated output is augmented with an observation at each time in <a href="#">stime()</a> . When using <code>obsaug</code> , a flag indicating augmented observations can be requested by including <code>a.u.g</code> in <code>carry_out</code> |

|                |  |
|----------------|--|
| tgrid          | a tgrid object; or a numeric vector of simulation times or another object with an stime method   |
| recsort        | record sorting flag. Default value is 1. Possible values are 1,2,3,4: 1 and 2 put doses in a data set after padded observations at the same time; 3 and 4 put those doses before padded observations at the same time. 2 and 4 will put doses scheduled through add1 after observations at the same time; 1 and 3 put doses scheduled through add1 before observations at the same time. recsort will not change the order of your input data set if both doses and observations are given.  |
| deslist        | a list of tgrid objects  |
| descol         | the name of a column for assigning designs   |
| filbak         | carry data items backward when the first data set row has time greater than zero   |
| tad            | when TRUE a column is added to simulated output is added showing the time since the last dose. Only data records with evid == 1 will be considered doses for the purposes of tad calculation. The tad can be properly calculated with a dosing lag time in the model as long as the dosing lag time (specified in \$MAIN) is always appropriate for any subsequent doses scheduled through add1. This will always be true if the lag time doesn't change over time. But it might (possibly) not hold if the lag time changes prior to the last dose in the add1 sequence. This known limitation shouldn't affect tad calculation in most common dosing lag time implementations. |
| nocb           | if TRUE, use next observation carry backward method; otherwise, use locf.  |
| skip_init_calc | don't use \$MAIN to calculate initial conditions   |
| ss_n           | maximum number of iterations for determining steady state for the PK system; a warning will be issued if steady state is not achieved within ss_n iterations when ss_fixed is TRUE   |
| ss_fixed       | if FALSE (the default), then a warning will be issued if the system does not reach steady state within ss_n iterations given the model tolerances rtol and atol; if TRUE, the number of iterations for determining steady state are capped at ss_n and no warning will be issued if steady state has not been reached within ss_n dosing iterations. To silence warnings related to steady state, set ss_fixed to TRUE and set ss_n as the maximum number of iterations to try when advancing the system for steady state determination.   |

## Details

- Use mrgsim\_df to return a data frame rather than mrgsim object
- Both data and idata will be coerced to numeric matrix
- carry\_out can be used to insert data columns into the output data set. This is partially dependent on the nature of the data brought into the problem
- When using data and idata together, an error is generated if an ID occurs in data but not idata. Also, when looking up data in idata, ID in idata is assumed to be uniquely keyed to ID in data. No error is generated if ID is duplicated in data; parameters will be used from the first occurrence found in idata

- `carry_out`: `idata` is assumed to be individual-level and variables that are carried from `idata` are repeated throughout the individual's simulated data. Variables carried from data are carried via last-observation carry forward. NA is returned from observations that are inserted into simulated output that occur prior to the first record in data
- `recover`: this is similar to `carry_out` with respect to end result, but it uses a different process. Columns to be recovered are cached prior to running the simulation, and then joined back on to the simulated data. So, whereas `carry_out` will only accept numeric data items, `recover` can handle data frame columns of any type. There is a small decrease in performance with `recover` compared to `carry_out`, but it is likely that the performance difference is difficult to perceive (when the simulation runs very fast) or only a small fractional increase in run time when the simulation is very large. And any performance hit is likely to be well worth it in light of the convenience gain. Just think carefully about using this feature when every millisecond counts.

### Value

An object of class `mrgsims`

### See Also

`mrgsim_variants`, `mrgsim_q`

### Examples

```
## example("mrgsim")

e <- ev(amt = 1000)

mod <- mrgsolve::house()

out <- mod %>% ev(e) %>% mrgsim()

plot(out)

out <- mod %>% ev(e) %>% mrgsim(end=22)

out

data(exTheoph)

out <- mod %>% data_set(exTheoph) %>% mrgsim()

out

out <- mod %>% mrgsim(data=exTheoph)

out <- mrgsim(mod, data=exTheoph, obsonly=TRUE)

out

out <- mod %>% mrgsim(data=exTheoph, obsaug=TRUE, carry_out="a.u.g")
```

```
out

out <- mod %>% ev(e) %>% mrgsim(outvars="CP,RESP")

out

a <- ev(amt = 1000, group = 'a')
b <- ev(amt = 750, group = 'b')
data <- as_data_set(a,b)

out <- mrgsim_d(mod, data, recover="group")

out
```

---

mrgsims\_dplyr

*Methods for handling output with dplyr verbs*

---

## Description

These methods modify the data in a mrgsims object and return a data frame. Contrast with the functions in [mrgsims\\_modify](#).

## Usage

```
## S3 method for class 'mrgsims'
pull(.data, ...)

## S3 method for class 'mrgsims'
filter(.data, ...)

## S3 method for class 'mrgsims'
group_by(.data, ..., add = FALSE, .add = FALSE)

## S3 method for class 'mrgsims'
distinct(.data, ..., .keep_all = FALSE)

## S3 method for class 'mrgsims'
mutate(.data, ...)

## S3 method for class 'each'
summarise(.data, funs, ...)

## S3 method for class 'mrgsims'
summarise(.data, ...)

## S3 method for class 'mrgsims'
```

```

do(.data, ..., .dots)

## S3 method for class 'mrgsims'
select(.data, ...)

## S3 method for class 'mrgsims'
slice(.data, ...)

as_data_frame.mrgsims(.data_, ...)

## S3 method for class 'mrgsims'
as_tibble(.data_, ...)

as.tbl.mrgsims(x, ...)

```

### Arguments

|                        |  |
|------------------------|--|
| <code>.data</code>     | an mrgsims object; passed to various dplyr functions           |
| <code>...</code>       | passed to other methods  |
| <code>add</code>       | passed to <a href="#">dplyr::group_by</a> (for dplyr < 1.0.0)  |
| <code>.add</code>      | passed to <a href="#">dplyr::group_by</a> (for dplyr >= 1.0.0) |
| <code>.keep_all</code> | passed to <a href="#">dplyr::distinct</a>                      |
| <code>funcs</code>     | passed to <a href="#">dplyr::summarise_each</a>                |
| <code>.dots</code>     | passed to various dplyr functions                              |
| <code>.data_</code>    | mrgsims object   |
| <code>x</code>         | passed to <code>dplyr::as.tbl</code>                           |

### Details

For the `select_sims` function, the dots `...` must be either compartment names or variables in `$CAPTURE`. An error will be generated if no valid names are selected or the names for selection are not found in the simulated output.

### See Also

[mrgsims\\_modify](#)

### Examples

```

out <- mrgsim(house(), events = ev(amt = 100), end = 5, delta=1)

dplyr::filter(out, time==2)

dplyr::mutate(out, label = "abc")

dplyr::select(out, time, RESP, CP)

```



---

|                |  |
|----------------|--|
| mrgsims_modify | <i>Methods for modifying mrgsims objects</i> |
|----------------|--|

---

### Description

These functions modify the simulated data in an `mrgsims` object and return the modified object. Contrast with the functions in [mrgsims\\_dplyr](#).

### Usage

```
mutate_sims(.data, ...)
```

```
select_sims(.data, ...)
```

```
filter_sims(.data, ...)
```

### Arguments

|                    |  |
|--------------------|--|
| <code>.data</code> | a <code>mrgsims</code> object                              |
| <code>...</code>   | other arguments passed to the <code>dplyr</code> functions |

### See Also

[mrgsims\\_dplyr](#)

### Examples

```
out <- mrgsim(house(), events = ev(amt = 100))  
filter_sims(out, time > 2)  
mutate_sims(out, label = "abc")  
select_sims(out, RESP, CP)
```

---

|          |   |
|----------|---|
| mrgsim_q | <i>Simulate from a model object with quicker turnaround</i> |
|----------|---|

---

### Description

Use the function when you would usually use [mrgsim\\_d](#), but you need a quicker turnaround time. The timing differences might be difficult to detect for a single simulation run but could become appreciable with repeated simulation. See details for important differences in how [mrgsim\\_q](#) is invoked compared to [mrgsim](#) and [mrgsim\\_d](#). This function should always be used for benchmarking simulation time with `mrgsolve`.

**Usage**

```
mrgsim_q(
  x,
  data,
  recsort = 1,
  stime = numeric(0),
  output = "mrgsims",
  skip_init_calc = FALSE,
  simcall = 0
)
```

**Arguments**

|                |   |
|----------------|---|
| x              | a model object  |
| data           | a simulation data set   |
| recsort        | record sorting flag   |
| stime          | a numeric vector of observation times; these observation times will only be added to the output if there are no observation records in data |
| output         | output data type; if mrgsims, then the default output object is returned; if "df" then a data frame is returned                             |
| skip_init_calc | don't use \$MAIN to calculate initial conditions  |
| simcall        | not used; only the default value of 0 is allowed  |

**Details**

This function does not support the piped simulation workflow. All arguments must be passed into the function except for `x`.

A data set is required for this simulation workflow. The data set can have only dosing records or doses with observations. When the data set only includes doses, a single numeric vector of observation times should be passed in.

This simulation workflow does not support Req (request) functionality. All compartments and captured variables will always be returned in the simulation output.

This simulation workflow does not support carry-out functionality.

This simulation workflow does not accept arguments to be passed to [update](#). This must be done by a separate call to [update](#).

This simulation workflow does not support use of event objects. If an event object is needed, it should be converted to a data set prior to the simulation run (see [as\\_data\\_set](#) or [as.data.frame.ev](#)).

This simulation workflow does not support idata sets or any feature enabled by idata set use. Individual level parameters should be joined onto the data set prior to simulation. Otherwise [mrgsim\\_i](#) or [mrgsim\\_ei](#) should be used.

By default, a mrgsims object is returned (as with [mrgsim](#)). Use the `output="df"` argument to request a plain data.frame of simulated data on return.

**See Also**

[mrgsim](#), [mrgsim\\_variants](#), [qsim](#)

**Examples**

```
mod <- mrgsolve::house()
data <- expand.ev(amt = c(100,300,1000))
out <- mrgsim_q(mod,data)

out
```

---

mrgsim\_variants      *mrgsim variant functions*

---

**Description**

These functions are called by [mrgsim](#) and have explicit input requirements written into the function name. The motivation behind these variants is to give the user a clear workflow with specific, required inputs as indicated by the function name. Use [mrgsim\\_q](#) instead to benchmark mrgsolve or to do repeated quick simulation for tasks like parameter optimization, sensitivity analyses, or optimal design.

**Usage**

```
mrgsim_e(x, events, idata = NULL, data = NULL, ...)
mrgsim_d(x, data, idata = NULL, events = NULL, ...)
mrgsim_ei(x, events, idata, data = NULL, ...)
mrgsim_di(x, data, idata, events = NULL, ...)
mrgsim_i(x, idata, data = NULL, events = NULL, ...)
mrgsim_0(x, idata = NULL, data = NULL, events = NULL, ...)
```

**Arguments**

|        |  |
|--------|--|
| x      | the model object   |
| events | an event object  |
| idata  | a matrix or data frame of model parameters, one parameter per row (see <a href="#">idata_set</a> ) |
| data   | NMTRAN-like data set (see <a href="#">data_set</a> )   |
| ...    | passed to <a href="#">update</a> and <a href="#">do_mrgsim</a>                                     |

## Details

**Important:** all of these functions require that data, idata, and/or events be pass directly to the functions. They will not recognize these inputs from a pipeline.

- `mrgsim_e` simulate using an event object
- `mrgsim_ei` simulate using an event object and `idata_set`
- `mrgsim_d` simulate using a `data_set`
- `mrgsim_di` simulate using a `data_set` and `idata_set`
- `mrgsim_i` simulate using a `idata_set`
- `mrgsim_0` simulate using just the model
- `mrgsim_q` simulate from a data set with quicker turnaround (see [mrgsim\\_q](#))

## See Also

[mrgsim](#), [mrgsim\\_q](#), [qsim](#)

---

mrgsolve

*mrgsolve*

---

## Description

mrgsolve is an R package maintained under the auspices of Metrum Research Group that facilitates simulation from models based on systems of ordinary differential equations (ODE) that are typically employed for understanding pharmacokinetics, pharmacodynamics, and systems biology and pharmacology. mrgsolve consists of computer code written in the R and C++ languages, providing an interface to a C++ translation of the lsoda differential equation solver. See [aboutsolver](#) for more information.

## Resources

- Main mrgsolve resource page: <https://mrgsolve.github.io>
- User guide: [https://mrgsolve.github.io/user\\_guide](https://mrgsolve.github.io/user_guide)
- Vignettes: <https://mrgsolve.github.io/vignettes>

## Package-wide options

- `mrgolve.project`: sets the default project director ([mread](#))
- `mrgsolve.soloc`: sets the default package build directory ([mread](#))
- `mrgsolve.mread.quiet`: don't print messages during [mread](#)
- `mrgsolve.update.strict`: if TRUE, print warning when trying to update an item in the model object that doesn't exist

**Examples**

```
## example("mrgsolve")

mod <- mrgsolve::house(delta=0.1) %>% param(CL=0.5)

events <- ev(amt=1000, cmt=1, addl=5, ii=24)

events

mod

see(mod)

## Not run:
stime(mod)

## End(Not run)
param(mod)

init(mod)

out <- mod %>% ev(events) %>% mrgsim(end=168)

head(out)
tail(out)
dim(out)

plot(out, GUT+CP~.)

sims <- as.data.frame(out)

t72 <- dplyr::filter(sims, time==72)

str(t72)

idata <- data.frame(ID=c(1,2,3), CL=c(0.5,1,2),VC=12)
out <- mod %>% ev(events) %>% mrgsim(end=168, idata=idata, req="")
plot(out)

out <- mod %>% ev(events) %>% mrgsim(carry_out="amt,evid,cmt,CL")
head(out)

ev1 <- ev(amt=500, cmt=2,rate=10)
ev2 <- ev(amt=100, cmt=1, time=54, ii=8, addl=10)
events <- c(ev1+ev2)
events

out <- mod %>% ev(events) %>% mrgsim(end=180, req="")
plot(out)
```

```
## "Condensed" data set
data(extran1)
extran1

out <- mod %>% data_set(extran1) %>% mrgsim(end=200)

plot(out,CP~time|factor(ID))

## idata
data(exidata)

out <-
  mod %>%
  ev(amt=1000, cmt=1) %>%
  idata_set(exidata) %>%
  mrgsim(end=72)

plot(out, CP~., as="log10")

# Internal model library
## Not run:
mod <- mread("irm1", modlib())

mod

x <- mod %>% ev(amt=300, ii=12, addl=3) %>% mrgsim

## End(Not run)
```

---

mutate.ev

*dplyr verbs for event objects*

---

## Description

dplyr verbs for event objects

## Usage

```
## S3 method for class 'ev'
mutate(.data, ...)
```

```
## S3 method for class 'ev'
select(.data, ...)
```

```
## S3 method for class 'ev'
filter(.data, ...)
```

**Arguments**

.data            the event object  
 ...             passed to the dplyr function

---

names,mrgmod-method    *Get all names from a model object*

---

**Description**

Get all names from a model object

**Usage**

```
## S4 method for signature 'mrgmod'
names(x)
```

**Arguments**

x                the model object

**Examples**

```
mod <- mrgsolve::house()
names(mod)
```

---

nmext                    *Import model estimates from a NONMEM ext file*

---

**Description**

Import model estimates from a NONMEM ext file

**Usage**

```
nmext(
  run = NA_real_,
  project = getwd(),
  file = paste0(run, ".ext"),
  path = NULL,
  theta = TRUE,
  omega = TRUE,
  sigma = TRUE,
  olabels = NULL,
  slabels = NULL,
```

```

  oprefix = "",
  sprefix = "",
  tname = "THETA",
  oname = "...",
  sname = "...",
  read_fun = "data.table"
)

```

### Arguments

|          |  |
|----------|--|
| run      | run number                                       |
| project  | project directory                                |
| file     | deprecated; use path instead                     |
| path     | full path to NONMEM ext file                     |
| theta    | logical; if TRUE, the \$THETA vector is returned |
| omega    | logical; if TRUE, the \$OMEGA matrix is returned |
| sigma    | logical; if TRUE, the \$SIGMA matrix is returned |
| olabels  | labels for \$OMEGA                               |
| slabels  | labels for \$SIGMA                               |
| oprefix  | prefix for \$OMEGA labels                        |
| sprefix  | prefix for \$SIGMA labels                        |
| tname    | name for \$THETA                                 |
| oname    | name for \$OMEGA                                 |
| sname    | name for \$SIGMA                                 |
| read_fun | function to use when reading the ext file        |

### See Also

[nmxml](#), [read\\_nmext](#)

---

nmxml

*Import model estimates from a NONMEM xml file*

---

### Description

Import model estimates from a NONMEM xml file



**Usage**

```
nmxml(
  run = numeric(0),
  project = character(0),
  file = character(0),
  path = character(0),
  theta = TRUE,
  omega = TRUE,
  sigma = TRUE,
  olabels = NULL,
  slabels = NULL,
  oprefix = "",
  sprefix = "",
  tname = "THETA",
  oname = "...",
  sname = "...",
  index = "last",
  xpath = "../nm:estimation"
)
```

**Arguments**

|         |   |
|---------|---|
| run     | run number  |
| project | project directory   |
| file    | deprecated; use path instead  |
| path    | the complete path to the run.xml file   |
| theta   | logical; if TRUE, the \$THETA vector is returned  |
| omega   | logical; if TRUE, the \$OMEGA matrix is returned  |
| sigma   | logical; if TRUE, the \$SIGMA matrix is returned  |
| olabels | labels for \$OMEGA  |
| slabels | labels for \$SIGMA  |
| oprefix | prefix for \$OMEGA labels   |
| sprefix | prefix for \$SIGMA labels   |
| tname   | name for \$THETA  |
| oname   | name for \$OMEGA  |
| sname   | name for \$SIGMA  |
| index   | the estimation number to return; "last" will return the last estimation results; otherwise, pass an integer indicating which estimation results to return |
| xpath   | xml path containing run results; if the default doesn't work, consider using <code>../estimation</code> as an alternative; see details                    |

**Details**

If `run` and `project` are supplied, the `.xml` file is assumed to be located in `run.xml`, in directory `run` off the project directory. If `file` is supplied, `run` and `project` arguments are ignored.

This function requires that the `xml2` package be installed and loadable. If `requireNamespace("xml2")` fails, an error will be generated.

`nmxml` usually expects to find run results in the `xpath` called `./nm:estimation`. Occasionally, the run results are not stored in this namespace but no namespaces are found in the `xml` file. In this case, the user can specify the `xpath` containing run results. Consider trying `./estimation` as an alternative if the default fails.

**Value**

A list with `theta`, `omega` and `sigma` elements, depending on what was requested

**See Also**

`nmext`

**Examples**

```
if(requireNamespace("xml2")) {
  proj <- system.file("nonmem", package = "mrgsolve")
  mrgsolve::nmxml(run = 1005, project = proj)
}
```

---

numerics\_only

*Prepare data.frame for input to mrgsim*

---

**Description**

Prepare data.frame for input to mrgsim

**Usage**

```
numerics_only(x, quiet = FALSE, convert_lgl = FALSE)
```

**Arguments**

|                          |  |
|--------------------------|--|
| <code>x</code>           | a input data set   |
| <code>quiet</code>       | logical indicating whether or not warnings should be printed     |
| <code>convert_lgl</code> | if TRUE, convert logical columns with <a href="#">as.integer</a> |

---

|        |   |
|--------|---|
| obsaug | <i>Augment observations in the simulated output</i> |
|--------|---|

---

**Description**

Augment observations in the simulated output

**Usage**

```
obsaug(x, value = TRUE, ...)
```

**Arguments**

|       |  |
|-------|--|
| x     | model object   |
| value | the value for obsaug   |
| ...   | passed along There is also a obsaug argument to <a href="#">mrgsim</a> that can be set to accomplish the same thing as a call to obsaug in the pipeline. |

---

|         |  |
|---------|--|
| obsonly | <i>Collect only observations in the simulated output</i> |
|---------|--|

---

**Description**

Collect only observations in the simulated output

**Usage**

```
obsonly(x, value = TRUE, ...)
```

**Arguments**

|       |                       |
|-------|-----------------------|
| x     | model object          |
| value | the value for obsonly |
| ...   | passed along          |

**Details**

There is also a obsonly argument to [mrgsim](#) that can be set to accomplish the same thing as a call to obsonly in the pipeline.

---

 omega

 Manipulate OMEGA matrices
 

---

## Description

The primary function is `omat` that can be used to both get the `$OMEGA` matrices out of a model object and to update `$OMEGA` matrices in a model object.

## Usage

```
omat(.x, ...)

## S4 method for signature 'missing'
omat(.x, ...)

## S4 method for signature 'matrix'
omat(.x, ..., labels = list())

## S4 method for signature ``NULL``
omat(.x, ...)

## S4 method for signature 'list'
omat(.x, ...)

## S4 method for signature 'omegalist'
omat(.x, ...)

## S4 method for signature 'mrgmod'
omat(.x, ..., make = FALSE, open = FALSE)

## S4 method for signature 'mrgsims'
omat(.x, make = FALSE, ...)
```

## Arguments

|                     |  |
|---------------------|--|
| <code>.x</code>     | a matrix, list of matrices or <code>matlist</code> object  |
| <code>...</code>    | passed to other functions, including <a href="#">modMATRIX</a>   |
| <code>labels</code> | character vector of names for <code>\$OMEGA</code> elements; must be equal to number of rows/columns in the matrix |
| <code>make</code>   | logical; if <code>TRUE</code> , matrix list is rendered into a single matrix                                       |
| <code>open</code>   | passed to <a href="#">merge.list</a>   |
| <code>x</code>      | <code>matlist</code> object  |

**Examples**

```

## example("omega")
mat1 <- matrix(1)
mat2 <- diag(c(1,2,3))
mat3 <- matrix(c(0.1, 0.002, 0.002, 0.5), 2,2)
mat4 <- dmat(0.1, 0.2, 0.3, 0.4)

omat(mat1)
omat(mat1, mat2, mat3)
omat(A=mat1, B=mat2, C=mat3)

mod <- mrgsolve::house() %>% omat(mat4)

omat(mod)
omat(mod, make=TRUE)

## Not run:

$OMEGA
1 2 3

$OMEGA @block
1 0.1 2

$OMEGA \@cor
\@ prefix ETA_
\@ labels CL VC KA
0.1
0.67 0.2
0 0 0.3

## End(Not run)

```

---

outvars

*Show names of current output variables*


---

**Description**

Show names of current output variables

**Usage**

```
outvars(x, unlist = FALSE)
```

**Arguments**

|        |  |
|--------|--|
| x      | mrgmod object  |
| unlist | if TRUE then a character vector (rather than list) is returned |

**Value**

When `unlist` is `FALSE` (default) : a named list, with `cmt` showing names of output compartments and `capture` giving names of output variables in capture. When `unlist` is `TRUE`, then a single, unnamed character vector of `outvar` names is returned.

**Examples**

```
outvars(mrgsolve::house())
```

---

 param

---

*Create and work with parameter objects*


---

**Description**

See [numericlist](#) for methods to deal with `parameter_list` objects.

**Usage**

```
param(.x, ...)
```

```
## S4 method for signature 'mrgmod'
```

```
param(.x, .y = NULL, ..., .pat = "*", .strict = FALSE)
```

```
## S4 method for signature 'mrgsims'
```

```
param(.x, ...)
```

```
## S4 method for signature 'missing'
```

```
param(..., .strict = TRUE)
```

```
## S4 method for signature 'list'
```

```
param(.x, ...)
```

```
## S4 method for signature 'ANY'
```

```
param(.x, ...)
```

```
allparam(.x)
```

**Arguments**

|                      |   |
|----------------------|---|
| <code>.x</code>      | the model object  |
| <code>...</code>     | passed along or name/value pairs to update the parameters in a model object                           |
| <code>.y</code>      | list to be merged into parameter list   |
| <code>.pat</code>    | a regular expression (character) to be applied as a filter for which parameters to show when printing |
| <code>.strict</code> | if <code>TRUE</code> , all names to be updated must be found in the parameter list                    |

## Details

Can be used to either get a parameter list object from a `mrgmod` model object or to update the parameters in a model object. For both uses, the return value is a `parameter_list` object. For the former use, `param` is usually called to print the parameters to the screen, but the `parameter_list` object can also be coerced to a list or numeric R object.

Use `allparam` to get a `parameter_list` object including both model parameters and data items listed in `$FIXED`.

## Value

An object of class `parameter_list` (see [numericlist](#)).

## Examples

```
## example("param")

mod <- mrgsolve::house()

param(mod)

param(mod, .pat="^(C|F)") ## may be useful when large number of parameters

class(param(mod))

param(mod)$KA

param(mod)[["KA"]]

as.list(param(mod))

as.data.frame(param(mod))

param(mod, CL = 1.2)

new_values <- list(CL = 1.3, VC = 20.5)

param(mod, new_values)
```

---

PKMODEL

*Parse PKMODEL BLOCK data*

---

## Description

Parse PKMODEL BLOCK data

**Usage**

```
PKMODEL(
  ncmt = 1,
  depot = FALSE,
  cmt = NULL,
  trans = pick_trans(ncmt, depot),
  env = list(),
  pos = 1,
  ...
)
```

**Arguments**

|       |  |
|-------|--|
| ncmt  | number of compartments; must be 1 (one-compartment, not including a depot dosing compartment) or 2 (two-compartment model, not including a depot dosing compartment) |
| depot | logical indicating whether to add depot compartment  |
| cmt   | compartment names as comma-delimited character   |
| trans | the parameterization for the PK model; must be 1, 2, 4, or 11  |
| env   | parse environment  |
| pos   | block position number  |
| ...   | not used   |

**Details**

When using \$PKMODEL, certain symbols must be defined in the model specification depending on the value of ncmt, depot and trans.

- ncmt 1, depot FALSE, trans 2: CL, V
- ncmt 1, depot TRUE, trans 2: CL, V, KA
- ncmt 2, depot FALSE, trans 4: CL, V1, Q, V2
- ncmt 2, depot TRUE, trans 4: CL, V2, Q, V3, KA

If trans=11 is specified, use the symbols listed above for the ncmt / depot combination, but append *i* at the end (e.g. CL<sub>*i*</sub> or Q<sub>*i*</sub> or KA<sub>*i*</sub>).

If trans=1, the user must utilize the following symbols:

- pred\_CL for clearance
- pred\_V or pred\_V2 for central compartment volume of distribution
- pred\_Q for intercompartmental clearance
- pred\_V3 for peripheral compartment volume of distribution
- pred\_KA for absorption rate constant

**See Also**

[BLOCK\\_PARSE](#)



---

 plot, batch\_mrgsims, missing-method

*Plot method for mrgsims objects*


---

## Description

Plot method for mrgsims objects

## Usage

```
## S4 method for signature 'batch_mrgsims,missing'
plot(x, yval = variables(x), auto.key = list(), mincol = 3, ...)

## S4 method for signature 'batch_mrgsims,formula'
plot(
  x,
  y,
  show.grid = TRUE,
  lwd = 2,
  type = "l",
  yval = variables(x),
  auto.key = list(columns = 1),
  scales = list(y = list(relation = "free")),
  ...
)
```

## Arguments

|           |                                  |
|-----------|----------------------------------|
| x         | mrgsims object                   |
| yval      | y variables to plot              |
| auto.key  | passed to xyplot                 |
| mincol    | minimum number of columns in key |
| ...       | arguments passed to xyplot       |
| y         | a formula passed to xyplot       |
| show.grid | print grid in the plot           |
| lwd       | passed to xyplot                 |
| type      | passed to xyplot                 |
| scales    | passed to xyplot                 |

---

`plot_mrgsims`*Generate a quick plot of simulated data*

---

**Description**

Generate a quick plot of simulated data

**Usage**

```
## S4 method for signature 'mrgsims,missing'
plot(x, limit = 16, ...)

## S4 method for signature 'mrgsims,formula'
plot(
  x,
  y,
  limit = 16,
  show.grid = TRUE,
  outer = TRUE,
  type = "l",
  lwd = 2,
  ylab = "value",
  groups = ID,
  scales = list(y = list(relation = "free")),
  logy = FALSE,
  logbr = 1,
  ...
)

## S4 method for signature 'mrgsims,character'
plot(x, y, ...)
```

**Arguments**

|                        |  |
|------------------------|--|
| <code>x</code>         | mrgsims object                                       |
| <code>limit</code>     | limit the the number of panels to create             |
| <code>...</code>       | other arguments passed to xyplot                     |
| <code>y</code>         | formula used for plotting                            |
| <code>show.grid</code> | logical indicating whether or not to draw panel.grid |
| <code>outer</code>     | passed to xyplot                                     |
| <code>type</code>      | passed to xyplot                                     |
| <code>lwd</code>       | passed to xyplot                                     |
| <code>ylab</code>      | passed to xyplot                                     |
| <code>groups</code>    | passed to xyplot                                     |

|        |   |
|--------|---|
| scales | passed to xyplot  |
| logy   | plot the y variables on log scale   |
| logbr  | log scale breaks indicator; use 1 for breaks every log unit; use 3 for breaks every half log unit; use 0 for default breaks |

### Examples

```

mod <- mrgsolve::house(end=48, delta=0.2) %>% init(GUT=1000)

out <- mrgsim(mod)

plot(out)

plot(out, subset=time <=24)

plot(out, GUT+CP~.)

plot(out, CP+RESP~time, col="black", scales="same", lty=2)

## Not run:
plot(out, "CP RESP, GUT")

## End(Not run)

```

---

|           |                                       |
|-----------|---------------------------------------|
| plot_sims | <i>Plot data as an mrgsims object</i> |
|-----------|---------------------------------------|

---

### Description

Plot data as an mrgsims object

### Usage

```
plot_sims(.data, ..., .f = NULL, .dots = list())
```

### Arguments

|       |   |
|-------|---|
| .data | a data frame                              |
| ...   | unquoted column names to plot on y-axis   |
| .f    | a formula to plot                         |
| .dots | extra arguments passed to lattice::xyplot |

### Details

This function is only intended for use with data frames that were created by modifying an mrgsims object.

**Examples**

```

mod <- mrgsolve::house() %>% ev(amt = 100)

out <- mrgsim(mod)
out_df <- dplyr::mutate(out, time <= 72)

plot(out)
plot_sims(out, CP, RESP)

## Not run:
plot_sims(out, .f = ~ CP + RESP)
plot_sims(out, .f = CP + RESP ~ time)

## End(Not run)

```

---

qsim

*Basic, simple simulation from model object*


---

**Description**

This is just a lighter version of [mrgsim](#), with fewer options. See [Details](#).

**Usage**

```

qsim(
  x,
  data,
  idata = no_idata_set(),
  obsonly = FALSE,
  tgrid = NULL,
  recsort = 1,
  tad = FALSE,
  Req = NULL,
  outvars = Req,
  skip_init_calc = FALSE,
  output = "mrgsims"
)

```

**Arguments**

|         |  |
|---------|--|
| x       | the model object   |
| data    | can be either event object or data set   |
| idata   | a matrix or data frame of model parameters, one parameter per row (see <a href="#">idata_set</a> ) |
| obsonly | if TRUE, dosing records are not included in the output   |

|                |  |
|----------------|--|
| tgrid          | a tgrid object; or a numeric vector of simulation times or another object with an stime method   |
| recsort        | record sorting flag. Default value is 1. Possible values are 1,2,3,4: 1 and 2 put doses in a data set after padded observations at the same time; 3 and 4 put those doses before padded observations at the same time. 2 and 4 will put doses scheduled through add1 after observations at the same time; 1 and 3 put doses scheduled through add1 before observations at the same time. recsort will not change the order of your input data set if both doses and observations are given.  |
| tad            | when TRUE a column is added to simulated output is added showing the time since the last dose. Only data records with evid == 1 will be considered doses for the purposes of tad calculation. The tad can be properly calculated with a dosing lag time in the model as long as the dosing lag time (specified in \$MAIN) is always appropriate for any subsequent doses scheduled through add1. This will always be true if the lag time doesn't change over time. But it might (possibly) not hold if the lag time changes prior to the last dose in the add1 sequence. This known limitation shouldn't affect tad calculation in most common dosing lag time implementations. |
| Req            | synonym for outvars  |
| outvars        | output items to request; if missing, then only captured items will be returned in the output   |
| skip_init_calc | don't use \$MAIN to calculate initial conditions   |
| output         | output data type; the default is mrgsims, which returns the default output object; other options include df (for data.frame) or matrix   |

## Details

There is no pipeline interface for this function; all configuration options (see Arguments) must be passed as formal arguments to the function. You can't carry\_out, Request specific columns, or pass items in for update. Some other limitations, but only convenience-related. See Arguments for available options. Specifically, there is no . . . argument for this function. Use the [mrgsolve::update](#) method to update the model object.

## See Also

[mrgsim\\_q](#), [mrgsim](#), [mrgsim\\_variants](#)

## Examples

```
mod <- mrgsolve::house()
dose <- ev(amt = 100)
out <- qsim(mod,dose)
```

---

`read_nmext`*Extract estimates from NONMEM ext file*

---

**Description**

Extract estimates from NONMEM ext file

**Usage**

```
read_nmext(  
  run = NA_real_,  
  project = getwd(),  
  file = paste0(run, ".ext"),  
  path = NULL,  
  read_fun = c("data.table", "read.table")  
)
```

**Arguments**

|                       |                                      |
|-----------------------|--------------------------------------|
| <code>run</code>      | a run number or run identifier       |
| <code>project</code>  | the NONMEM project directory         |
| <code>file</code>     | the ext file name                    |
| <code>path</code>     | full path and file name for ext file |
| <code>read_fun</code> | function to read the ext file        |

**Value**

A list with `param`, `omega`, and `sigma` in a format ready to be used to update a model object.

**Examples**

```
project <- system.file("nonmem", package = "mrgsolve")  
est <- read_nmext(1005, project = project)  
est$param  
est$omega  
est$sigma
```

---

|              |  |
|--------------|--|
| realize_addl | <i>Make addl doses explicit in an event object or data set</i> |
|--------------|--|

---

### Description

Make addl doses explicit in an event object or data set

### Usage

```
realize_addl(x, ...)

## S3 method for class 'data.frame'
realize_addl(
  x,
  warn = FALSE,
  mark_new = FALSE,
  fill = c("inherit", "na", "locf"),
  ...
)

## S3 method for class 'ev'
realize_addl(x, ...)
```

### Arguments

|          |  |
|----------|--|
| x        | a data_set data frame or an ev object (see details)  |
| ...      | not used   |
| warn     | if TRUE a warning is issued if no ADDL or addl column is found   |
| mark_new | if TRUE, a flag is added to indicate new columns   |
| fill     | specifies how to handle non-dose related data columns in new data set records; this option is critical when handling data sets with time-varying, non-dose-related data items; see details |

### Details

If no addl column is found the data frame is returned and a warning is issued if warn is true. If ii, time, or evid are missing, an error is generated.

Use caution when passing in data that has non-dose-related data columns that vary within a subject and pay special attention to the fill argument. By definition, realize\_addl will add new rows to your data frame and it is not obvious how the non-dose-related data should be handled in these new rows. When inherit is chosen, the new records have non-dose-related data that is identical to the originating dose record. This should be fine when these data items are not varying with time, but will present a problem when the data are varying with time. When locf is chosen, the missing data are filled in with NA and an last observation carry forward operation is applied to **every** column in the data set. This may not be what you want if you already had missing values in the input data set and want to preserve that missingness. When na is chosen, the missing data are filled in with

NA and no locf operation is applied. But note that these missing values may be problematic for a mrgsolve simulation run. If you have any time-varying columns or missing data in your data set, be sure to check that the output from this function is what you were expecting.

---

|        |                                     |
|--------|-------------------------------------|
| render | <i>Render a model to a document</i> |
|--------|-------------------------------------|

---

### Description

Render a model to a document

### Usage

```
render(x, ...)

## S4 method for signature 'character'
render(x, project = NULL, ...)

## S4 method for signature 'mrgmod'
render(x, ...)

dorender(model, project, template = NULL, compile = TRUE, ...)
```

### Arguments

|          |   |
|----------|---|
| x        | model object or the model name                            |
| ...      | passed to <code>rmarkdown::render</code>                  |
| project  | the directory containing the <code>.cpp</code> model file |
| model    | model name  |
| template | template document   |
| compile  | logical; if true, the model will be compiled to run       |

### Examples

```
## Not run:
mod <- mrgsolve::house()
mrgsolve::render(mod)
mrgsolve::render("irm2", modlib())

## End(Not run)
```



---

|     |                                 |
|-----|---------------------------------|
| Req | <i>Request simulated output</i> |
|-----|---------------------------------|

---

### Description

Use this function to select, by name, either compartments or derived variables that have been captured (see [CAPTURE](#)) into the simulated output.

### Usage

```
Req(x, ...)  
  
req(x, ...)  
  
## S3 method for class 'mrgmod'  
req(x, ...)
```

### Arguments

|     |  |
|-----|--|
| x   | model object                                   |
| ... | unquoted names of compartments or tabled items |

### Details

There is also a Req argument to [mrgsim](#) that can be set to accomplish the same thing as a call to Req in the pipeline.

Note the difference between req and Req: the former only selects compartments to appear in output while the latter selects both compartments and captured items. Also, when there are items are explicitly listed in Req, all other compartments or captured items not listed there are ignored. But when compartments are selected with req all of the captured items are returned. Remember that req is strictly for compartments.

### Examples

```
mod <- mrgsolve::house()  
  
mod %>% Req(CP,RESP) %>% ev(amt=1000) %>% mrgsim
```

---

 reserved

*Reserved words*


---

**Description**

Reserved words

**Usage**

```
reserved()
```

**Details**

Note: this function is not exported; you must go into the `mrgsolve` namespace by using the `mrgsolve:::` prefix.

**Examples**

```
mrgsolve:::reserved()
```

---

 revar

*Get model random effect variances and covariances*


---

**Description**

Get model random effect variances and covariances

**Usage**

```
revar(x, ...)
```

```
## S4 method for signature 'mrgmod'
revar(x, ...)
```

**Arguments**

|                  |              |
|------------------|--------------|
| <code>x</code>   | model object |
| <code>...</code> | passed along |

---

|     |  |
|-----|--|
| see | <i>Print model code to the console</i> |
|-----|--|

---

**Description**

Print model code to the console

**Usage**

```
see(x, ...)
```

```
## S4 method for signature 'mrgmod'
```

```
see(x, raw = FALSE, ...)
```

**Arguments**

|     |                     |
|-----|---------------------|
| x   | model object        |
| ... | passed along        |
| raw | return the raw code |

**Value**

invisible NULL

---

|       |                                  |
|-------|----------------------------------|
| sigma | <i>Manipulate SIGMA matrices</i> |
|-------|----------------------------------|

---

**Description**

The primary function is `smat` that can be used to both get the `$SIGMA` matrices out of a model object and to update `$SIGMA` matrices in a model object.

**Usage**

```
smat(.x, ...)
```

```
## S4 method for signature 'missing'
```

```
smat(.x, ...)
```

```
## S4 method for signature 'matrix'
```

```
smat(.x, ..., labels = list())
```

```
## S4 method for signature 'list'
```

```
smat(.x, ...)
```

```
## S4 method for signature 'sigalist'
smat(.x, ...)

## S4 method for signature 'mrgmod'
smat(.x, ..., make = FALSE, open = FALSE)

## S4 method for signature '`NULL`'
smat(.x, ...)

## S4 method for signature 'mrgsims'
smat(.x, make = FALSE, ...)
```

### Arguments

|        |   |
|--------|---|
| .x     | a matrix, list of matrices or matlist object  |
| ...    | passed to other functions, including <a href="#">modMATRIX</a>  |
| labels | character vector of names for \$SIGMA elements; must be equal to number of rows/columns in the matrix |
| make   | logical; if TRUE, matrix list is rendered into a single matrix  |
| open   | passed to <a href="#">merge.list</a>  |
| x      | matlist object  |

### Examples

```
## example("sigma")
mat1 <- matrix(1)
mat2 <- diag(c(1,2))
mat3 <- matrix(c(0.1, 0.002, 0.002, 0.5), 2,2)
mat4 <- dmat(0.1, 0.2, 0.3, 0.4)

smat(mat1)
smat(mat1, mat2, mat3)
smat(A=mat1, B=mat2, C=mat3)

mod <- mrgsolve::house() %>% smat(mat1)

smat(mod)
smat(mod, make=TRUE)
```

---

simargs

*Access or clear arguments for calls to mrgsim*

---

### Description

Access or clear arguments for calls to mrgsim

**Usage**

```
simargs(x, which = NULL, clear = FALSE, ...)
```

**Arguments**

|       |   |
|-------|---|
| x     | model object  |
| which | character with length 1 naming a single arg to get                    |
| clear | logical indicating whether or not to clear args from the model object |
| ...   | passed along  |

**Value**

If `clear` is `TRUE`, the argument list is cleared and the model object is returned. Otherwise, the argument list is returned.

**Examples**

```
mod <- mrgsolve::house()
mod %>% Req(CP,RESP) %>% carry_out(evid,WT,FLAG) %>% simargs
```

---

soloc

*Return the location of the model shared object*

---

**Description**

Return the location of the model shared object

**Usage**

```
soloc(x, short = FALSE)
```

**Arguments**

|       |  |
|-------|--|
| x     | model object   |
| short | logical; if <code>TRUE</code> , <code>soloc</code> will be rendered with a short path name |

**Examples**

```
mod <- mrgsolve::house()
soloc(mod)
```

---

 solversettings

*Optional inputs for lsoda*


---

### Description

These are settings for the differential equation solver (lsoda) that can be accessed via the R interface. The code listing below is taken directly from the lsoda source code.

### Details

The following items can be set

- hmax (HMAX below); decrease hmax when you want to limit how big of a step the solver can take when integrating from one time to the next time. However be aware that smaller hmax will result in longer run times.
- hmin (HMIN below); don't fiddle with this unless you know what you're doing.
- ixpr (IXPR below)
- maxsteps (MXSTEP below); increase this number when the solver has a long interval between two integration times (e.g. when observation records are far apart).
- mxhnil (MXHNIL below); don't usually modify this one
- atol - the absolute solver tolerance; decrease this number (e.g. to 1E-10 or 1E-20 or 1E-50) when the value in a compartment can get extremely small; without this extra (lower) tolerance, the value can get so low that the number can randomly become negative. However be aware that more precision here will result in longer run times.
- rtol - the relative solver tolerances; decrease this number when you want a more precise solution. However be aware that more precision here will result in longer run times.

### See Also

[aboutsolver](#), [update](#)

---

 summary.mrgmod

*Print summary of a mrgmod object*


---

### Description

Print summary of a mrgmod object

### Usage

```
## S3 method for class 'mrgmod'
summary(object, ...)
```

**Arguments**

|        |                 |
|--------|-----------------|
| object | a mrgmod object |
| ...    | not used        |

---

|        |  |
|--------|--|
| tscale | <i>Re-scale time in the simulated output</i> |
|--------|--|

---

**Description**

Re-scale time in the simulated output

**Usage**

```
tscale(x, value = 1, ...)
```

**Arguments**

|       |                                    |
|-------|------------------------------------|
| x     | model object                       |
| value | value by which time will be scaled |
| ...   | passed along                       |

**Details**

There is also a `tscale` argument to `mrgsim` that can be set to accomplish the same thing as a call to `tscale` in the pipeline.

**Examples**

```
# The model is in hours:  
mod <- mrgsolve::house()  
  
# The output is in days:  
mod %>% tscale(1/24) %>% mrgsim
```

---

|        |                                |
|--------|--------------------------------|
| update | <i>Update the model object</i> |
|--------|--------------------------------|

---

### Description

After the model object is created, update various attributes.

### Usage

```
## S4 method for signature 'mrgmod'
update(object, ..., merge = TRUE, open = FALSE, data = NULL, strict = TRUE)

## S4 method for signature 'omegalist'
update(object, y, ...)

## S4 method for signature 'sigmalist'
update(object, y, ...)

## S4 method for signature 'parameter_list'
update(object, .y, ...)
```

### Arguments

|        |   |
|--------|---|
| object | a model object  |
| ...    | named items to update   |
| merge  | logical indicating to merge (rather than replace) new and existing attributes   |
| open   | logical; used only when merge is TRUE and parameter list or initial conditions list is being updated; if FALSE, no new items will be added; if TRUE, the parameter list may expand. |
| data   | a list of items to update; this list is combined with any items passed in via ...   |
| strict | if TRUE, then an error will be generated if there is attempt to update a non-existent item  |
| y      | another object involved in update   |
| .y     | data to update  |

### Details

Slots that can be updated:

- verbose
- debug
- preclean
- mindt
- digits



- atol - absolute solver tolerance; see [solversettings](#)
- rtol - relative solver tolerance; see [solversettings](#)
- ss\_rtol - relative tolerance when finding steady state
- ss\_atol - absolute tolerance when finding steady state
- ixpr - see IXPR in [solversettings](#)
- mxhnil - see MXHNIL in [solversettings](#)
- hmin - see HMIN in [solversettings](#)
- hmax - see HMAX in [solversettings](#)
- maxsteps - see MXSTEP in [solversettings](#)
- start, end, delta, add
- tscale
- request
- param
- init
- omega
- sigma
- outvars

### Value

The updated model object is returned.

### See Also

[update](#), [mrgmod-class](#), [within](#)

### Examples

```
## Not run:  
mod <- mrgsolve::house()  
  
mod <- update(mod, end=120, delta=4, param=list(CL=19.1))  
  
## End(Not run)
```

---

|                |  |
|----------------|--|
| valid_data_set | <i>Validate and prepare a data sets for simulation</i> |
|----------------|--|

---

### Description

This function is called by mrgsim. Users may also call this function to pre-validate data when the same data set is used for repeated simulation.

### Usage

```
valid_data_set(x, m = NULL, verbose = FALSE, quiet = FALSE)
```

```
valid_data_set.matrix(x, verbose = FALSE)
```

### Arguments

|         |                                      |
|---------|--------------------------------------|
| x       | data.frame or matrix                 |
| m       | a model object                       |
| verbose | logical                              |
| quiet   | if TRUE, messages will be suppressed |

### Value

A matrix with non-numeric columns dropped; if x is a data.frame with character cmt column comprised of valid compartment names and m is a model object, the cmt column will be converted to the corresponding compartment number.

### See Also

[valid\\_idata\\_set](#), [idata\\_set](#), [data\\_set](#)

### Examples

```
mod <- mrgsolve::house()
data(exTheoph)
d <- valid_data_set(exTheoph,mod)
```

---

|                 |  |
|-----------------|--|
| valid_idata_set | <i>Validate and prepare idata data sets for simulation</i> |
|-----------------|--|

---

**Description**

Validate and prepare idata data sets for simulation

**Usage**

```
valid_idata_set(x, m, verbose = FALSE, quiet = FALSE)
```

**Arguments**

|         |                                      |
|---------|--------------------------------------|
| x       | data.frame or matrix                 |
| m       | a model object                       |
| verbose | logical                              |
| quiet   | if TRUE, messages will be suppressed |

**Value**

A numeric matrix with class `valid_idata_set`.

**See Also**

[valid\\_data\\_set](#), [idata\\_set](#), [data\\_set](#)

---

|        |  |
|--------|--|
| within | <i>Update parameters, initials, and settings within a model object</i> |
|--------|--|

---

**Description**

The main use case for using [within](#) rather than [update](#) or [param](#) or [init](#) is when you want to update to a new value that is calculated from the existing value. See the example in details

**Usage**

```
## S3 method for class 'mrgmod'
within(data, expr, ...)
```

**Arguments**

|      |  |
|------|--|
| data | an object with class <code>mrgmod</code>   |
| expr | expressions evaluated in an environment containing various model object components, including parameters, initial conditions, and others (see details) |
| ...  | not used   |

**Details**

Other model object slots that can be updated: start, end, delta, add, rtol, atol, hmax, maxsteps. These are include for convenience, but we expect that most of the time these will get updated through the update method.

**See Also**

[update](#)

**Examples**

```
mod <- mrgsolve::house()
mod2 <- within(mod, {CL <- CL * 1.5})

mod$CL
mod2$CL
```

---

\$.ev-method

*Select columns from an ev object*

---

**Description**

Select columns from an ev object

**Usage**

```
## S4 method for signature 'ev'
x$name

## S4 method for signature 'ev'
x[[i, exact = TRUE]]
```

**Arguments**

|       |                      |
|-------|----------------------|
| x     | ev object            |
| name  | column to select     |
| i     | an element to select |
| exact | not used             |

---

\$.mrgmod-method      *Select parameter values from a model object*

---

### **Description**

The \$ and [[ operators get the value of a single parameter in the model. The [ gets several values, returning a named list.

### **Usage**

```
## S4 method for signature 'mrgmod'
x$name

## S4 method for signature 'mrgmod'
x[[i, exact = TRUE]]

## S4 method for signature 'mrgmod'
x[i]
```

### **Arguments**

|       |                      |
|-------|----------------------|
| x     | mrgmod object        |
| name  | parameter to take    |
| i     | an element to select |
| exact | not used             |

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