

# Package ‘pharmr’

June 15, 2023

**Encoding** UTF-8

**Version** 0.96.0

**Date** 2023-06-15

**Title** Interface to the 'PharmPy' 'Pharmacometrics' Library

**Maintainer** Rikard Nordgren <rikard.nordgren@farmaci.uu.se>

**Depends** R (>= 3.6.0), altair (>= 4.0.0)

**SystemRequirements** Python (>= 3.9.0)

**Imports** reticulate (>= 1.19), utils

**Suggests** testthat, magrittr, here, knitr

**NeedsCompilation** no

**Description** Interface to the 'PharmPy' 'pharmacometrics' library. The 'Reticulate' package is used to interface Python from R.

**Config/reticulate** list( packages = list( list(package = ``altair"),  
list(package = ``pharmPy-core") ) )

**URL** <https://github.com/pharmPy/pharmr>

**BugReports** <https://github.com/pharmPy/pharmr/issues>

**License** BSD\_2\_clause + file LICENSE

**RoxygenNote** 7.2.3

**Author** Rikard Nordgren [aut, cre, cph],  
Stella Belin [aut, cph],  
Mats O. Karlsson [sad],  
Andrew C. Hooker [sad],  
Sebastian Ueckert [sad] (<<https://orcid.org/0000-0002-3712-0255>>),  
Simon Carter [rev],  
Simon Buatois [rev],  
João A. Abrantes [rev],  
F. Hoffmann-La Roche Ltd. [fnd]

**Repository** CRAN

**Date/Publication** 2023-06-15 13:50:11 UTC

**R topics documented:**

add_allometry . . . . .	6
add_covariance_step . . . . .	7
add_covariate_effect . . . . .	8
add_estimation_step . . . . .	11
add_iiv . . . . .	12
add_individual_parameter . . . . .	14
add_iov . . . . .	14
add_lag_time . . . . .	15
add_metabolite . . . . .	16
add_peripheral_compartment . . . . .	17
add_pk_iiv . . . . .	18
add_population_parameter . . . . .	19
add_time_after_dose . . . . .	20
append_estimation_step_options . . . . .	20
bump_model_number . . . . .	21
calculate_aic . . . . .	22
calculate_bic . . . . .	22
calculate_corr_from_cov . . . . .	23
calculate_corr_from_prec . . . . .	24
calculate_cov_from_corrse . . . . .	25
calculate_cov_from_prec . . . . .	26
calculate_epsilon_gradient_expression . . . . .	27
calculate_eta_gradient_expression . . . . .	28
calculate_eta_shrinkage . . . . .	29
calculate_individual_parameter_statistics . . . . .	30
calculate_individual_shrinkage . . . . .	31
calculate_parameters_from_ucp . . . . .	32
calculate_pk_parameters_statistics . . . . .	33
calculate_prec_from_corrse . . . . .	34
calculate_prec_from_cov . . . . .	35
calculate_se_from_cov . . . . .	36
calculate_se_from_prec . . . . .	37
calculate_ucp_scale . . . . .	38
check_dataset . . . . .	38
check_high_correlations . . . . .	39
check_parameters_near_bounds . . . . .	40
check_pharmpy . . . . .	41
cleanup_model . . . . .	41
convert_model . . . . .	42
create_basic_pk_model . . . . .	43
create_config_template . . . . .	44
create_joint_distribution . . . . .	44
create_report . . . . .	45
create_results . . . . .	45
create_rng . . . . .	46
create_symbol . . . . .	47

deidentify_data . . . . .	47
display_odes . . . . .	48
drop_columns . . . . .	49
drop_dropped_columns . . . . .	50
evaluate_epsilon_gradient . . . . .	50
evaluate_eta_gradient . . . . .	51
evaluate_expression . . . . .	52
evaluate_individual_prediction . . . . .	53
evaluate_population_prediction . . . . .	54
evaluate_weighted_residuals . . . . .	55
expand_additional_doses . . . . .	56
find_clearance_parameters . . . . .	56
find_volume_parameters . . . . .	57
fit . . . . .	58
fix_or_unfix_parameters . . . . .	58
fix_parameters . . . . .	59
fix_parameters_to . . . . .	60
get_baselines . . . . .	61
get_bioavailability . . . . .	62
get_cmt . . . . .	62
get_concentration_parameters_from_data . . . . .	63
get_config_path . . . . .	63
get_covariate_baselines . . . . .	64
get_doseid . . . . .	65
get_doses . . . . .	65
get_dv_symbol . . . . .	66
get_evid . . . . .	67
get_ids . . . . .	67
get_individual_parameters . . . . .	68
get_individual_prediction_expression . . . . .	69
get_initial_conditions . . . . .	69
get_lag_times . . . . .	70
get_mdv . . . . .	71
get_model_code . . . . .	71
get_model_covariates . . . . .	72
get_number_of_individuals . . . . .	72
get_number_of_observations . . . . .	73
get_number_of_observations_per_individual . . . . .	74
get_observations . . . . .	75
get_observation_expression . . . . .	76
get_omegas . . . . .	76
get_pk_parameters . . . . .	77
get_population_prediction_expression . . . . .	78
get_rv_parameters . . . . .	79
get_sigmas . . . . .	79
get_thetas . . . . .	80
get_unit_of . . . . .	81
get_zero_order_inputs . . . . .	82

<code>greekify_model</code>	82
<code>has_additive_error_model</code>	83
<code>has_combined_error_model</code>	84
<code>has_covariate_effect</code>	85
<code>has_first_order_elimination</code>	85
<code>has_linear_odes</code>	86
<code>has_linear_odes_with_real_eigenvalues</code>	87
<code>has_michaelis_menten_elimination</code>	87
<code>has_mixed_mm_fo_elimination</code>	88
<code>has_odes</code>	89
<code>has_proportional_error_model</code>	89
<code>has_random_effect</code>	90
<code>has_weighted_error_model</code>	91
<code>has_zero_order_absorption</code>	92
<code>has_zero_order_elimination</code>	92
<code>install_pharmpy</code>	93
<code>install_pharmpy_devel</code>	94
<code>is_linearized</code>	94
<code>is_real</code>	95
<code>list_time_varying_covariates</code>	96
<code>load_example_model</code>	96
<code>load_example_modelfit_results</code>	97
<code>make_declarative</code>	98
<code>mu_reference_model</code>	98
<code>omit_data</code>	99
<code>plot_individual_predictions</code>	100
<code>plot_iofv_vs_iofv</code>	100
<code>plot_transformed_eta_distributions</code>	101
<code>predict_influential_individuals</code>	101
<code>predict_influential_outliers</code>	102
<code>predict_outliers</code>	103
<code>print_fit_summary</code>	104
<code>print_model_code</code>	104
<code>print_model_symbols</code>	105
<code>print_pharmpy_version</code>	105
<code>rank_models</code>	106
<code>read_dataset_from_datainfo</code>	107
<code>read_model</code>	107
<code>read_modelfit_results</code>	108
<code>read_model_from_string</code>	108
<code>read_results</code>	109
<code>remove_covariance_step</code>	110
<code>remove_covariate_effect</code>	111
<code>remove_error_model</code>	111
<code>remove_estimation_step</code>	112
<code>remove_iiv</code>	113
<code>remove_iov</code>	114
<code>remove_lag_time</code>	115

remove_loq_data . . . . .	115
remove_peripheral_compartment . . . . .	116
remove_unused_parameters_and_rvs . . . . .	117
rename_symbols . . . . .	117
resample_data . . . . .	118
reset_index . . . . .	119
reset_indices_results . . . . .	119
resume_tool . . . . .	120
retrieve_final_model . . . . .	120
retrieve_models . . . . .	121
run_allometry . . . . .	122
run_amd . . . . .	123
run_bootstrap . . . . .	124
run_covsearch . . . . .	125
run_estmethod . . . . .	126
run_iivsearch . . . . .	127
run_iovsearch . . . . .	128
run_modelfit . . . . .	129
run_modelsearch . . . . .	130
run_ruvsearch . . . . .	131
run_tool . . . . .	132
sample_individual_estimates . . . . .	132
sample_parameters_from_covariance_matrix . . . . .	133
sample_parameters_uniformly . . . . .	135
set_additive_error_model . . . . .	136
set_bolus_absorption . . . . .	137
set_combined_error_model . . . . .	138
set_covariates . . . . .	139
set_dtbs_error_model . . . . .	139
set_dvid . . . . .	140
set_estimation_step . . . . .	140
set_evaluation_step . . . . .	141
set_first_order_absorption . . . . .	142
set_first_order_elimination . . . . .	143
set_iiv_on_ruv . . . . .	144
set_initial_condition . . . . .	145
set_initial_estimates . . . . .	145
set_lower_bounds . . . . .	146
set_michaelis_menten_elimination . . . . .	147
set_mixed_mm_fo_elimination . . . . .	148
set_name . . . . .	149
set_ode_solver . . . . .	149
set_peripheral_compartments . . . . .	150
set_power_on_ruv . . . . .	151
set_proportional_error_model . . . . .	152
set_seq_zo_fo_absorption . . . . .	153
set_time_varying_error_model . . . . .	154
set_tmdd . . . . .	155

set_transit_compartments . . . . .	155
set_upper_bounds . . . . .	156
set_weighted_error_model . . . . .	157
set_zero_order_absorption . . . . .	158
set_zero_order_elimination . . . . .	158
set_zero_order_input . . . . .	159
simplify_expression . . . . .	160
solve_ode_system . . . . .	161
split_joint_distribution . . . . .	161
summarize_errors . . . . .	162
summarize_individuals . . . . .	163
summarize_individuals_count_table . . . . .	164
summarize_modelfit_results . . . . .	165
transform_blq . . . . .	165
transform_etas_boxcox . . . . .	167
transform_etas_john_draper . . . . .	168
transform_etas_tdist . . . . .	169
translate_nmtran_time . . . . .	170
unconstrain_parameters . . . . .	170
undrop_columns . . . . .	171
unfix_parameters . . . . .	172
unfix_parameters_to . . . . .	173
update_initial_individual_estimates . . . . .	174
update_inits . . . . .	174
use_thetas_for_error_stdev . . . . .	175
write_csv . . . . .	176
write_model . . . . .	177
write_results . . . . .	177

**Index****179**


---

add_allometry	<i>add_allometry</i>
---------------	----------------------

---

**Description**

Add allometric scaling of parameters

Add an allometric function to each listed parameter. The function will be  $P=P*(X/Z)**T$  where P is the parameter, X the allometric\_variable, Z the reference\_value and T is a theta. Default is to automatically use clearance and volume parameters.

**Usage**

```
add_allometry(
  model,
  allometric_variable = "WT",
  reference_value = 70,
```

```

    parameters = NULL,
    initials = NULL,
    lower_bounds = NULL,
    upper_bounds = NULL,
    fixed = TRUE
  )

```

### Arguments

model	(Model) Pharmpy model
allometric_variable	(str) Value to use for allometry (X above)
reference_value	(str or numeric) Reference value (Z above)
parameters	(array(str) (optional)) Parameters to use or NULL (default) for all available CL, Q and V parameters
initials	(array(numeric) (optional)) Initial estimates for the exponents. Default is to use 0.75 for CL and Qs and 1 for Vs
lower_bounds	(array(numeric) (optional)) Lower bounds for the exponents. Default is 0 for all parameters
upper_bounds	(array(numeric) (optional)) Upper bounds for the exponents. Default is 2 for all parameters
fixed	(logical) Whether the exponents should be fixed

### Value

(Model) Pharmpy model object

### Examples

```

## Not run:
model <- load_example_model("pheno")
model <- add_allometry(model, allometric_variable='WGT')
model$statements$before_odes

## End(Not run)

```

---

add\_covariance\_step    *add\_covariance\_step*

---

### Description

Adds covariance step to the final estimation step

**Usage**

```
add_covariance_step(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(Model) PharmPy model object

**See Also**

```
add_estimation_step  
set_estimation_step  
remove_estimation_step  
append_estimation_step_options  
remove_covariance_step  
set_evaluation_step
```

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- set_estimation_step(model, 'FOCE', cov=FALSE)  
model <- add_covariance_step(model)  
ests <- model$estimation_steps  
ests[1]  
  
## End(Not run)
```

---

add\_covariate\_effect *add\_covariate\_effect*

---

**Description**

Adds covariate effect to :class:pharmPy.model.

The following effects have templates:

- Linear function for continuous covariates (*lin*)
- Function:

math::

coveff = 1 + theta \* (cov - median)

- Init: 0.001



- Upper:
- If median of covariate equals minimum:  $:math:100,000$
- Otherwise:  $:math:\frac{1}{\{\{median\} - \{min\}\}}$
- Lower:
- If median of covariate equals maximum:  $:math:-100,000$
- Otherwise:  $:math:\frac{1}{\{\{median\} - \{max\}\}}$
- Linear function for categorical covariates (*cat*)
- Function:
- If covariate is most common category:

math::

coveff = 1

- For each additional category:

math::

coveff = 1 + theta

- Init:  $:math:0.001$
- Upper:  $:math:100,000$
- Lower:  $:math:-100,000$
- Piecewise linear function/"hockey-stick", continuous covariates only (*piece\_lin*)
- Function:
- If cov <= median:

math::

coveff = 1 + theta1 \* (cov - median)

- If cov > median:

math::

coveff = 1 + theta2 \* (cov - median)

- Init:  $:math:0.001$
- Upper:
- For first state:  $:math:\frac{1}{\{\{median\} - \{min\}\}}$
- Otherwise:  $:math:100,000$
- Lower:
- For first state:  $:math:-100,000$
- Otherwise:  $:math:\frac{1}{\{\{median\} - \{max\}\}}$
- Exponential function, continuous covariates only (*exp*)
- Function:

math::

coveff = exp(theta \* (cov - median))

- Init:
- If lower > 0.001 or upper < 0.001:  $\frac{\{\text{upper}\} - \{\text{lower}\}}{2}$
- If estimated init is 0:  $\frac{\{\text{upper}\}}{2}$
- Otherwise:  $0.001$
- Upper:
- If min - median = 0 or max - median = 0:  $100$
- Otherwise:

math::

min(fraclog(0.01)min - median, fraclog(100)max - median)

- Lower:
- If min - median = 0 or max - median = 0:  $0.01$
- Otherwise:

math::

max(fraclog(0.01)max - median, fraclog(100)min - median)

- Power function, continuous covariates only (*pow*)
- Function:

math::

coveff = (fraccovmedian)^theta

- Init:  $0.001$
- Upper:  $100,000$
- Lower:  $-100$

## Usage

```
add_covariate_effect(
  model,
  parameter,
  covariate,
  effect,
  operation = "*",
  allow_nested = FALSE
)
```

**Arguments**

model	(Model) Pharmpy model to add covariate effect to.
parameter	(str) Name of parameter to add covariate effect to.
covariate	(str) Name of covariate.
effect	(str) Type of covariate effect. May be abbreviated covariate effect (see above) or custom.
operation	(str) Whether the covariate effect should be added or multiplied (default).
allow_nested	(logical) Whether to allow adding a covariate effect when one already exists for the input parameter-covariate pair.

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- add_covariate_effect(model, "CL", "APGR", "exp")
model$statements$before_odes$full_expression("CL")

## End(Not run)
```

---

add\_estimation\_step    *add\_estimation\_step*

---

**Description**

Add estimation step

Adds estimation step for a model in a given index. Methods currently supported are: FO, FOCE, ITS, LAPLACE, IMPMAP, IMP, SAEM

**Usage**

```
add_estimation_step(model, method, idx = NULL, ...)
```

**Arguments**

model	(Model) Pharmpy model
method	(str) estimation method to change to
idx	(numeric (optional)) index of estimation step (starting from 0), default is NULL (adds step at the end)
...	Arguments to pass to EstimationStep (such as interaction, evaluation)

**Value**

(Model) Pharmpy model object

**See Also**

set\_estimation\_step  
remove\_estimation\_step  
append\_estimation\_step\_options  
add\_covariance\_step  
remove\_covariance\_step  
set\_evaluation\_step

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
opts <- list('NITER'=1000, 'ISAMPLE'=100)  
model <- add_estimation_step(model, "IMP", tool_options=opts)  
ests <- model$estimation_steps  
length(ests)  
ests[2]  
  
## End(Not run)
```

---

add\_iiv

*add\_iiv*

---

**Description**

Adds IIVs to :class:pharmpy.model.

Effects that currently have templates are:

- Additive (*add*)
- Proportional (*prop*)
- Exponential (*exp*)
- Logit (*log*)

For all except exponential the operation input is not needed. Otherwise user specified input is supported. Initial estimates for new etas are 0.09.

**Usage**

```
add_iiv(  
  model,  
  list_of_parameters,  
  expression,  
  operation = "*",  
  initial_estimate = 0.09,  
  eta_names = NULL  
)
```

**Arguments**

`model` (Model) Pharmpy model to add new IIVs to.

`list_of_parameters` (array(str) or str) Name/names of parameter to add new IIVs to.

`expression` (array(str) or str) Effect/effects on eta. Either abbreviated (see above) or custom.

`operation` (str) Whether the new IIV should be added or multiplied (default).

`initial_estimate` (numeric) Value of initial estimate of parameter. Default is 0.09

`eta_names` (array(str) (optional)) Custom name/names of new eta

**Value**

(Model) Pharmpy model object

**See Also**

```
add_pk_iiv  
add_iov  
remove_iiv  
remove_iov
```

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- remove_iiv(model, "CL")  
model <- add_iiv(model, "CL", "add")  
model$statements$find_assignment("CL")  
  
## End(Not run)
```

---

```
add_individual_parameter
      add_individual_parameter
```

---

**Description**

Add an individual or pk parameter to a model

**Usage**

```
add_individual_parameter(model, name)
```

**Arguments**

model	(Model) PharmPy model
name	(str) Name of individual/pk parameter

**Value**

(Model) PharmPy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- add_individual_parameter(model, "KA")
model$statements$find_assignment("KA")

## End(Not run)
```

---

```
add_iov      add_iov
```

---

**Description**

Adds IOVs to :class:pharmPy.model.

Initial estimate of new IOVs are 10% of the IIV eta it is based on.

**Usage**

```
add_iov(
  model,
  occ,
  list_of_parameters = NULL,
  eta_names = NULL,
  distribution = "disjoint"
)
```

**Arguments**

model	(Model) Pharmpy model to add new IOVs to.
occ	(str) Name of occasion column.
list_of_parameters	(array(str) or str (optional)) List of names of parameters and random variables. Accepts random variable names, parameter names, or a mix of both.
eta_names	(array(str) or str (optional)) Custom names of new etas. Must be equal to the number of input etas times the number of categories for occasion.
distribution	(str) The distribution that should be used for the new etas. Options are 'disjoint' for disjoint normal distributions, 'joint' for joint normal distribution, 'explicit' for an explicit mix of joint and disjoint distributions, and 'same-as-iiv' for copying the distribution of IIV etas.

**Value**

(Model) Pharmpy model object

**See Also**

add\_iiv  
 add\_pk\_iiv  
 remove\_iiv  
 remove\_iov

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- add_iov(model, "TIME", "CL")
model$statements$find_assignment("CL")

## End(Not run)
```

---

add\_lag\_time                      *add\_lag\_time*

---

**Description**

Add lag time to the dose compartment of model.

Initial estimate for lag time is set the previous lag time if available, otherwise it is set to the time of first observation/2.

**Usage**

```
add_lag_time(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(Model) PharmPy model object

**See Also**

set\_transit\_compartments

remove\_lag\_time

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- add_lag_time(model)

## End(Not run)
```

---

add\_metabolite      *add\_metabolite*

---

**Description**

Adds a metabolite compartment to a model

The flow from the central compartment to the metabolite compartment will be unidirectional.

**Usage**

```
add_metabolite(model, drug_dvid = 1)
```

**Arguments**

model (Model) PharmPy model

drug\_dvid (numeric) DVID for drug (assuming all other DVIDs being for metabolites)

**Value**

(Model) PharmPy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- add_metabolite(model)

## End(Not run)
```



---

```
add_peripheral_compartment
      add_peripheral_compartment
```

---

## Description

Add a peripheral distribution compartment to model

The rate of flow from the central to the peripheral compartment will be parameterized as  $QP_n / VC$  where  $VC$  is the volume of the central compartment. The rate of flow from the peripheral to the central compartment will be parameterized as  $QP_n / VP_n$  where  $VP_n$  is the volume of the added peripheral compartment.

Initial estimates:

```
=====  
1 :math:{CL} = {CL'}, :math:{VC} = {VC'}, :math:{QP1} = {CL'} and :math:{VP1} = {VC'} * 0.05  
2 :math:{QP1} = {QP1'} * 0.1, :math:{VP1} = {VP1'}, :math:{QP2} = {QP1'} * 0.9 and  
:math:{VP2} = {VP1'}===== n =====
```

## Usage

```
add_peripheral_compartment(model)
```

## Arguments

```
model          (Model) PharmPy model
```

## Value

```
(Model) PharmPy model object
```

## See Also

```
set_peripheral_compartment  
remove_peripheral_compartment
```

## Examples

```
## Not run:  
model <- load_example_model("pheno")  
model <- add_peripheral_compartment(model)  
model$statements$ode_system  
  
## End(Not run)
```

---

`add_pk_iiv``add_pk_iiv`

---

### Description

Adds IIVs to all PK parameters in `:class:pharmpy.model`.

Will add exponential IIVs to all parameters that are included in the ODE.

### Usage

```
add_pk_iiv(model, initial_estimate = 0.09)
```

### Arguments

`model` (Model) Pharmpy model to add new IIVs to.

`initial_estimate`  
(numeric) Value of initial estimate of parameter. Default is 0.09

### Value

(Model) Pharmpy model object

### See Also

`add_iiv`

`add_iov`

`remove_iiv`

`remove_iov`

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
model <- set_first_order_absorption(model)  
model$statements$find_assignment("MAT")  
model <- add_pk_iiv(model)  
model$statements$find_assignment("MAT")  
  
## End(Not run)
```

---

```
add_population_parameter  
    add_population_parameter
```

---

## Description

Add a new population parameter to the model

## Usage

```
add_population_parameter(  
  model,  
  name,  
  init,  
  lower = NULL,  
  upper = NULL,  
  fix = FALSE  
)
```

## Arguments

<code>model</code>	(Model) PharmPy model
<code>name</code>	(str) Name of the new parameter
<code>init</code>	(numeric) Initial estimate of the new parameter
<code>lower</code>	(numeric (optional)) Lower bound of the new parameter
<code>upper</code>	(numeric (optional)) Upper bound of the new parameter
<code>fix</code>	(logical) Should the new parameter be fixed?

## Value

(Model) PharmPy model object

## Examples

```
## Not run:  
model <- load_example_model("pheno")  
model <- add_population_parameter(model, 'POP_KA', 2)  
model$parameters  
  
## End(Not run)
```

---

```
add_time_after_dose  add_time_after_dose
```

---

**Description**

Calculate and add a TAD column to the dataset"

**Usage**

```
add_time_after_dose(model)
```

**Arguments**

model (Model) Pharmpy model

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- add_time_after_dose(model)  
  
## End(Not run)
```

---

```
append_estimation_step_options  
  append_estimation_step_options
```

---

**Description**

Append estimation step options  
Appends options to an existing estimation step.

**Usage**

```
append_estimation_step_options(model, tool_options, idx)
```

**Arguments**

model (Model) Pharmpy model  
tool\_options (list(str=any)) any additional tool specific options  
idx (numeric) index of estimation step (starting from 0)

**Value**

(Model) Pharmpy model object

**See Also**

add\_estimation\_step  
set\_estimation\_step  
remove\_estimation\_step  
add\_covariance\_step  
remove\_covariance\_step  
set\_evaluation\_step

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
opts <- list('NITER'=1000, 'ISAMPLE'=100)  
model <- append_estimation_step_options(model, tool_options=opts, idx=0)  
est <- model$estimation_steps[1]  
length(est$tool_options)  
  
## End(Not run)
```

---

bump\_model\_number      *bump\_model\_number*

---

**Description**

If the model name ends in a number increase it

If path is set increase the number until no file exists with the same name in path. If model name does not end in a number do nothing.

**Usage**

```
bump_model_number(model, path = NULL)
```

**Arguments**

model                    (Model) Pharmpy model object  
path                    (str) Default is to not look for files.

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- model$replace(name="run2")
model <- bump_model_number(model)
model$name

## End(Not run)
```

---

calculate_aic	<i>calculate_aic</i>
---------------	----------------------

---

**Description**

Calculate AIC

$$\text{AIC} = -2\text{LL} + 2 * n_{\text{estimated\_parameters}}$$
**Usage**

```
calculate_aic(model, likelihood)
```

**Arguments**

model	(Model) PharmPy model object
likelihood	(numeric) -2LL

**Value**

(numeric) AIC of model fit

---

calculate_bic	<i>calculate_bic</i>
---------------	----------------------

---

**Description**

Calculate BIC

Different variations of the BIC can be calculated:

- | mixed (default) |  $\text{BIC} = -2\text{LL} + n_{\text{random\_parameters}} * \log(n_{\text{individuals}}) + | n_{\text{fixed\_parameters}} * \log(n_{\text{observations}})$
- | fixed |  $\text{BIC} = -2\text{LL} + n_{\text{estimated\_parameters}} * \log(n_{\text{observations}})$
- | random |  $\text{BIC} = -2\text{LL} + n_{\text{estimated\_parameters}} * \log(n_{\text{individuals}})$
- | iiv |  $\text{BIC} = -2\text{LL} + n_{\text{estimated\_iiv\_omega\_parameters}} * \log(n_{\text{individuals}})$

**Usage**

```
calculate_bic(model, likelihood, type = NULL)
```

**Arguments**

```
model          (Model) Pharmpy model object
likelihood     (numeric) -2LL to use
type           (str (optional)) Type of BIC to calculate. Default is the mixed effects.
```

**Value**

(numeric) BIC of model fit

**Examples**

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_modelfit_results("pheno")
ofv <- results$ofv
calculate_bic(model, ofv)
calculate_bic(model, ofv, type='fixed')
calculate_bic(model, ofv, type='random')
calculate_bic(model, ofv, type='iiv')

## End(Not run)
```

---

```
calculate_corr_from_cov
      calculate_corr_from_cov
```

---

**Description**

Calculate correlation matrix from a covariance matrix

**Usage**

```
calculate_corr_from_cov(cov)
```

**Arguments**

```
cov          (data.frame) Covariance matrix
```

**Value**

(data.frame) Correlation matrix

**See Also**

calculate\_se\_from\_cov : Standard errors from covariance matrix  
calculate\_se\_from\_prec : Standard errors from precision matrix  
calculate\_cov\_from\_prec : Covariance matrix from precision matrix  
calculate\_cov\_from\_corrse : Covariance matrix from correlation matrix and standard errors  
calculate\_prec\_from\_cov : Precision matrix from covariance matrix  
calculate\_prec\_from\_corrse : Precision matrix from correlation matrix and standard errors  
calculate\_corr\_from\_prec : Correlation matrix from precision matrix

**Examples**

```
## Not run:  
results <- load_example_modelfit_results("pheno")  
cov <- results$covariance_matrix  
cov  
calculate_corr_from_cov(cov)  
  
## End(Not run)
```

---

```
calculate_corr_from_prec  
      calculate_corr_from_prec
```

---

**Description**

Calculate correlation matrix from a precision matrix

**Usage**

```
calculate_corr_from_prec(precision_matrix)
```

**Arguments**

precision\_matrix  
(data.frame) Precision matrix

**Value**

(data.frame) Correlation matrix



**See Also**

calculate\_se\_from\_cov : Standard errors from covariance matrix  
calculate\_se\_from\_prec : Standard errors from precision matrix  
calculate\_corr\_from\_cov : Correlation matrix from covariance matrix  
calculate\_cov\_from\_prec : Covariance matrix from precision matrix  
calculate\_cov\_from\_corrse : Covariance matrix from correlation matrix and standard errors  
calculate\_prec\_from\_cov : Precision matrix from covariance matrix  
calculate\_prec\_from\_corrse : Precision matrix from correlation matrix and standard errors

**Examples**

```
## Not run:  
results <- load_example_modelfit_results("pheno")  
prec <- results$precision_matrix  
prec  
calculate_corr_from_prec(prec)  
  
## End(Not run)
```

---

```
calculate_cov_from_corrse  
      calculate_cov_from_corrse
```

---

**Description**

Calculate covariance matrix from a correlation matrix and standard errors

**Usage**

```
calculate_cov_from_corrse(corr, se)
```

**Arguments**

corr            (data.frame) Correlation matrix  
se              (array) Standard errors

**Value**

(data.frame) Covariance matrix

**See Also**

calculate\_se\_from\_cov : Standard errors from covariance matrix  
calculate\_se\_from\_prec : Standard errors from precision matrix  
calculate\_corr\_from\_cov : Correlation matrix from covariance matrix  
calculate\_cov\_from\_prec : Covariance matrix from precision matrix  
calculate\_prec\_from\_cov : Precision matrix from covariance matrix  
calculate\_prec\_from\_corrse : Precision matrix from correlation matrix and standard errors  
calculate\_corr\_from\_prec : Correlation matrix from precision matrix

**Examples**

```
## Not run:  
results <- load_example_modelfit_results("pheno")  
corr <- results$correlation_matrix  
se <- results$standard_errors  
corr  
calculate_cov_from_corrse(corr, se)  
  
## End(Not run)
```

---

```
calculate_cov_from_prec  
      calculate_cov_from_prec
```

---

**Description**

Calculate covariance matrix from a precision matrix

**Usage**

```
calculate_cov_from_prec(precision_matrix)
```

**Arguments**

```
precision_matrix  
  (data.frame) Precision matrix
```

**Value**

(data.frame) Covariance matrix

**See Also**

calculate\_se\_from\_cov : Standard errors from covariance matrix  
calculate\_se\_from\_prec : Standard errors from precision matrix  
calculate\_corr\_from\_cov : Correlation matrix from covariance matrix  
calculate\_cov\_from\_corrse : Covariance matrix from correlation matrix and standard errors  
calculate\_prec\_from\_cov : Precision matrix from covariance matrix  
calculate\_prec\_from\_corrse : Precision matrix from correlation matrix and standard errors  
calculate\_corr\_from\_prec : Correlation matrix from precision matrix

**Examples**

```
## Not run:  
results <- load_example_modelfit_results("pheno")  
prec <- results$precision_matrix  
prec  
calculate_cov_from_prec(prec)  
  
## End(Not run)
```

---

```
calculate_epsilon_gradient_expression  
    calculate_epsilon_gradient_expression
```

---

**Description**

Calculate the symbolic expression for the epsilon gradient  
This function currently only support models without ODE systems

**Usage**

```
calculate_epsilon_gradient_expression(model)
```

**Arguments**

model (Model) PharmPy model object

**Value**

(Expression) Symbolic expression

**See Also**

calculate\_eta\_gradient\_expression : Eta gradient

**Examples**

```
## Not run:  
model <- load_example_model("pheno_linear")  
calculate_epsilon_gradient_expression(model)  
  
## End(Not run)
```

---

```
calculate_eta_gradient_expression  
    calculate_eta_gradient_expression
```

---

**Description**

Calculate the symbolic expression for the eta gradient

This function currently only support models without ODE systems

**Usage**

```
calculate_eta_gradient_expression(model)
```

**Arguments**

model (Model) PharmPy model object

**Value**

(Expression) Symbolic expression

**See Also**

calculate\_epsilon\_gradient\_expression : Epsilon gradient

**Examples**

```
## Not run:  
model <- load_example_model("pheno_linear")  
calculate_eta_gradient_expression(model)  
  
## End(Not run)
```

---

```
calculate_eta_shrinkage  
    calculate_eta_shrinkage
```

---

### Description

Calculate eta shrinkage for each eta

### Usage

```
calculate_eta_shrinkage(  
  model,  
  parameter_estimates,  
  individual_estimates,  
  sd = FALSE  
)
```

### Arguments

model	(Model) Pharnpy model
parameter_estimates	(array) Parameter estimates
individual_estimates	(data.frame) Table of individual (eta) estimates
sd	(logical) Calculate shrinkage on the standard deviation scale (default is to calculate on the variance scale)

### Value

(Series) Shrinkage for each eta

### See Also

calculate\_individual\_shrinkage

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
results <- load_example_model_fit_results("pheno")  
pe <- results$parameter_estimates  
ie <- results$individual_estimates  
calculate_eta_shrinkage(model, pe, ie)  
calculate_eta_shrinkage(model, pe, ie, sd=TRUE)  
  
## End(Not run)
```

---

```
calculate_individual_parameter_statistics
      calculate_individual_parameter_statistics
```

---

## Description

Calculate statistics for individual parameters

Calculate the mean (expected value of the distribution), variance (variance of the distribution) and standard error for individual parameters described by arbitrary expressions. Any dataset column or variable used in the model can be used in the expression. The exception being that variables that depends on the solution of the ODE system cannot be used. If covariates are used in the expression the statistics of the parameter is calculated at the median value of each covariate as well as at the 5:th and 95:th percentiles. If no parameter uncertainty is available for the model the standard error will not be calculated.

## Usage

```
calculate_individual_parameter_statistics(
  model,
  expr_or_exprs,
  parameter_estimates,
  covariance_matrix = NULL,
  rng = NULL
)
```

## Arguments

model	(Model) A previously estimated model
expr_or_exprs	(array(str) or str) Parameter estimates
parameter_estimates	(array) Parameter uncertainty covariance matrix
covariance_matrix	(data.frame (optional)) sympy expression or iterable of str or sympy expressions Expressions or equations for parameters of interest. If equations are used the names of the left hand sides will be used as the names of the parameters.
rng	(numeric (optional)) Random number generator or integer seed

## Value

(data.frame) A DataFrame of statistics indexed on parameter and covariate value.

## Examples

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_modelfit_results("pheno")
```

```
rng <- create_rng(23)
pe <- results$parameter_estimates
cov <- results$covariance_matrix
calculate_individual_parameter_statistics(model, "K=CL/V", pe, cov, rng=rng)

## End(Not run)
```

---

```
calculate_individual_shrinkage
      calculate_individual_shrinkage
```

---

### Description

Calculate the individual eta-shrinkage

Definition:  $\eta_{shr} = (\text{var}(\eta) / \omega)$

### Usage

```
calculate_individual_shrinkage(
  model,
  parameter_estimates,
  individual_estimates_covariance
)
```

### Arguments

`model` (Model) PharmPy model  
`parameter_estimates` (array) Parameter estimates of model  
`individual_estimates_covariance` (data.frame) Uncertainty covariance matrices of individual estimates

### Value

(DataFrame) Shrinkage for each eta and individual

### See Also

`calculate_eta_shrinkage`

## Examples

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_model_fit_results("pheno")
pe <- results$parameter_estimates
covs <- results$individual_estimates_covariance
calculate_individual_shrinkage(model, pe, covs)

## End(Not run)
```

---

```
calculate_parameters_from_ucp
      calculate_parameters_from_ucp
```

---

## Description

Scale parameter values from ucp to normal scale

## Usage

```
calculate_parameters_from_ucp(model, scale, ucps)
```

## Arguments

model	(Model) PharmPy model
scale	(UCPScale) A parameter scale
ucps	(array or list(str=numeric)) Series of parameter values

## Value

(data.frame) Parameters on the normal scale

## See Also

calculate\_ucp\_scale : Calculate the scale for conversion from ucps

## Examples

```
## Not run:
model <- load_example_model("pheno")
scale <- calculate_ucp_scale(model)
values <- {'PTVCL': 0.1, 'PTVV': 0.1, 'THETA_3': 0.1, 'IVCL': 0.1, 'IVV': 0.1, 'SIGMA_1_1': 0.1}
calculate_parameters_from_ucp(model, scale, values)

## End(Not run)
```



---

```
calculate_pk_parameters_statistics  
    calculate_pk_parameters_statistics
```

---

### Description

Calculate statistics for common pharmacokinetic parameters

Calculate the mean (expected value of the distribution), variance (variance of the distribution) and standard error for some individual pre-defined pharmacokinetic parameters.

### Usage

```
calculate_pk_parameters_statistics(  
  model,  
  parameter_estimates,  
  covariance_matrix = NULL,  
  rng = NULL  
)
```

### Arguments

model	(Model) A previously estimated model
parameter_estimates	(array) Parameter estimates
covariance_matrix	(data.frame (optional)) Parameter uncertainty covariance matrix
rng	(numeric (optional)) Random number generator or seed

### Value

(data.frame) A DataFrame of statistics indexed on parameter and covariate value.

### See Also

calculate\_individual\_parameter\_statistics : Calculation of statistics for arbitrary parameters

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
results <- load_example_model_fit_results("pheno")  
rng <- create_rng(23)  
pe <- results$parameter_estimates  
cov <- results$covariance_matrix  
calculate_pk_parameters_statistics(model, pe, cov, rng=rng)  
  
## End(Not run)
```

calculate\_prec\_from\_corrse  
*calculate\_prec\_from\_corrse*

---

**Description**

Calculate precision matrix from a correlation matrix and standard errors

**Usage**

```
calculate_prec_from_corrse(corr, se)
```

**Arguments**

corr            (data.frame) Correlation matrix  
se              (array) Standard errors

**Value**

(data.frame) Precision matrix

**See Also**

calculate\_se\_from\_cov : Standard errors from covariance matrix  
calculate\_se\_from\_prec : Standard errors from precision matrix  
calculate\_corr\_from\_cov : Correlation matrix from covariance matrix  
calculate\_cov\_from\_prec : Covariance matrix from precision matrix  
calculate\_cov\_from\_corrse : Covariance matrix from correlation matrix and standard errors  
calculate\_prec\_from\_cov : Precision matrix from covariance matrix  
calculate\_corr\_from\_prec : Correlation matrix from precision matrix

**Examples**

```
## Not run:  
results <- load_example_modelfit_results("pheno")  
corr <- results$correlation_matrix  
se <- results$standard_errors  
corr  
calculate_prec_from_corrse(corr, se)  
  
## End(Not run)
```

---

```
calculate_prec_from_cov  
    calculate_prec_from_cov
```

---

**Description**

Calculate precision matrix from a covariance matrix

**Usage**

```
calculate_prec_from_cov(cov)
```

**Arguments**

cov (data.frame) Covariance matrix

**Value**

(data.frame) Precision matrix

**See Also**

calculate\_se\_from\_cov : Standard errors from covariance matrix

calculate\_se\_from\_prec : Standard errors from precision matrix

calculate\_corr\_from\_cov : Correlation matrix from covariance matrix

calculate\_cov\_from\_prec : Covariance matrix from precision matrix

calculate\_cov\_from\_corrse : Covariance matrix from correlation matrix and standard errors

calculate\_prec\_from\_corrse : Precision matrix from correlation matrix and standard errors

calculate\_corr\_from\_prec : Correlation matrix from precision matrix

**Examples**

```
## Not run:  
results <- load_example_modelfit_results("pheno")  
cov <- results$covariance_matrix  
cov  
calculate_prec_from_cov(cov)  
  
## End(Not run)
```

---

calculate\_se\_from\_cov *calculate\_se\_from\_cov*

---

### Description

Calculate standard errors from a covariance matrix

### Usage

```
calculate_se_from_cov(cov)
```

### Arguments

cov (data.frame) Input covariance matrix

### Value

(data.frame) Standard errors

### See Also

calculate\_se\_from\_prec : Standard errors from precision matrix  
calculate\_corr\_from\_cov : Correlation matrix from covariance matrix  
calculate\_cov\_from\_prec : Covariance matrix from precision matrix  
calculate\_cov\_from\_corrse : Covariance matrix from correlation matrix and standard errors  
calculate\_prec\_from\_cov : Precision matrix from covariance matrix  
calculate\_prec\_from\_corrse : Precision matrix from correlation matrix and standard errors  
calculate\_corr\_from\_prec : Correlation matrix from precision matrix

### Examples

```
## Not run:  
results <- load_example_modelfit_results("pheno")  
cov <- results$covariance_matrix  
cov  
calculate_se_from_cov(cov)  
  
## End(Not run)
```

---

```
calculate_se_from_prec  
    calculate_se_from_prec
```

---

**Description**

Calculate standard errors from a precision matrix

**Usage**

```
calculate_se_from_prec(precision_matrix)
```

**Arguments**

```
precision_matrix  
    (data.frame) Input precision matrix
```

**Value**

(data.frame) Standard errors

**See Also**

calculate\_se\_from\_cov : Standard errors from covariance matrix  
calculate\_corr\_from\_cov : Correlation matrix from covariance matrix  
calculate\_cov\_from\_prec : Covariance matrix from precision matrix  
calculate\_cov\_from\_corrse : Covariance matrix from correlation matrix and standard errors  
calculate\_prec\_from\_cov : Precision matrix from covariance matrix  
calculate\_prec\_from\_corrse : Precision matrix from correlation matrix and standard errors  
calculate\_corr\_from\_prec : Correlation matrix from precision matrix

**Examples**

```
## Not run:  
results <- load_example_modelfit_results("pheno")  
prec <- results$precision_matrix  
prec  
calculate_se_from_prec(prec)  
  
## End(Not run)
```

---

calculate\_ucp\_scale     *calculate\_ucp\_scale*

---

**Description**

Calculate a scale for unconstrained parameters for a model

The UCPScale object can be used to calculate unconstrained parameters back into the normal parameter space.

**Usage**

```
calculate_ucp_scale(model)
```

**Arguments**

model                    (Model) Model for which to calculate an ucp scale

**Value**

(UCPScale) A scale object

**See Also**

calculate\_parameters\_from\_ucp : Calculate parameters from ucp:s

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
scale <- calculate_ucp_scale(model)  
  
## End(Not run)
```

---

check\_dataset             *check\_dataset*

---

**Description**

Check dataset for consistency across a set of rules

**Usage**

```
check_dataset(model, dataframe = FALSE, verbose = FALSE)
```

**Arguments**

model	(Model) PharmPy model object
dataframe	(logical) TRUE to return a DataFrame instead of printing to the console
verbose	(logical) Print out all rules checked if TRUE else print only failed rules

**Value**

(data.frame) Only returns a DataFrame is dataframe=TRUE

---

check\_high\_correlations  
*check\_high\_correlations*

---

**Description**

Check for highly correlated parameter estimates

**Usage**

```
check_high_correlations(model, cor, limit = 0.9)
```

**Arguments**

model	(Model) PharmPy model object
cor	(data.frame) Estimated correlation matrix
limit	(numeric) Lower limit for a high correlation

**Value**

(data.frame) Correlation values indexed on pairs of parameters for (absolute) correlations above limit

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
results <- load_example_model_fit_results("pheno")  
cor <- results$correlation_matrix  
check_high_correlations(model, cor, limit=0.3)  
  
## End(Not run)
```

---

```
check_parameters_near_bounds  
    check_parameters_near_bounds
```

---

### Description

Check if any estimated parameter value is close to its bounds

### Usage

```
check_parameters_near_bounds(  
  model,  
  values,  
  zero_limit = 0.001,  
  significant_digits = 2  
)
```

### Arguments

model	(Model) PharmPy model object
values	(array) Series of values with index a subset of parameter names.
zero_limit	(numeric) maximum distance to 0 bounds
significant_digits	(numeric) maximum distance to non-zero bounds in number of significant digits

### Value

(data.frame) Logical Series with same index as values

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
results <- load_example_modelfit_results("pheno")  
check_parameters_near_bounds(model, results$parameter_estimates)  
  
## End(Not run)
```



---

check_pharmpy	<i>Checks version of Pharmpy/pharmr</i>
---------------	---

---

**Description**

Checks whether Pharmpy and pharmr has the same version

**Usage**

```
check_pharmpy(pharmpy_version)
```

**Arguments**

pharmpy\_version  
(str) version number as string

---

cleanup_model	<i>cleanup_model</i>
---------------	----------------------

---

**Description**

Perform various cleanups of a model

This is what is currently done

- Make model statements declarative, i.e. only one assignment per symbol
- Inline all assignments of one symbol, e.g. X = Y

**Usage**

```
cleanup_model(model)
```

**Arguments**

model (Model) Pharmpy model object

**Value**

(Model) Reference to the same model

**Note**

When creating NONMEM code from the cleaned model Pharmpy might need to add certain assignments to make it in line with what NONMEM requires.

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model$statements
model <- cleanup_model(model)
model$statements

## End(Not run)
```

---

convert_model	<i>convert_model</i>
---------------	----------------------

---

**Description**

Convert model to other format

Note that the operation is not done inplace.

**Usage**

```
convert_model(model, to_format)
```

**Arguments**

model	(Model) Model to convert
to_format	(str) Name of format to convert into. Currently supported 'generic', 'nlmixr', 'nonmem', and 'rxode'

**Value**

(Model) New model object with new underlying model format

**Examples**

```
## Not run:
model <- load_example_model("pheno")
converted_model <- convert_model(model, "nlmixr")

## End(Not run)
```

---

`create_basic_pk_model create_basic_pk_model`

---

## Description

Creates a basic pk model of given type

## Usage

```
create_basic_pk_model(  
  modeltype,  
  dataset_path = NULL,  
  cl_init = 0.01,  
  vc_init = 1,  
  mat_init = 0.1  
)
```

## Arguments

<code>modeltype</code>	(str) Type of PK model to create. Supported are 'oral' and 'iv'
<code>dataset_path</code>	(str (optional)) Optional path to a dataset
<code>cl_init</code>	(numeric) Initial estimate of the clearance parameter
<code>vc_init</code>	(numeric) Initial estimate of the central volume parameter
<code>mat_init</code>	(numeric) Initial estimate of the mean absorption time parameter (if applicable)

## Value

(Model) PharmPy model object

## Examples

```
## Not run:  
model <- create_basic_pk_model('oral')  
  
## End(Not run)
```

---

```
create_config_template
    create_config_template
```

---

**Description**

Create a basic config file template

If a configuration file already exists it will not be overwritten

**Usage**

```
create_config_template()
```

**Examples**

```
## Not run:
create_config_template()

## End(Not run)
```

---

```
create_joint_distribution
    create_joint_distribution
```

---

**Description**

Combines some or all etas into a joint distribution.

The etas must be IIVs and cannot be fixed. Initial estimates for covariance between the etas is dependent on whether the model has results from a previous run. In that case, the correlation will be calculated from individual estimates, otherwise correlation will be set to 10%.

**Usage**

```
create_joint_distribution(model, rvs = NULL, individual_estimates = NULL)
```

**Arguments**

model	(Model) Pharnpy model
rvs	(array(str) (optional)) Sequence of etas or names of etas to combine. If NULL, all etas that are IIVs and non-fixed will be used (full block). NULL is default.
individual_estimates	(data.frame (optional)) Optional individual estimates to use for calculation of initial estimates

**Value**

(Model) Pharmpy model object

**See Also**

split\_joint\_distribution : split etas into separate distributions

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model$random_variables$etas  
model <- create_joint_distribution(model, c('ETA_1', 'ETA_2'))  
model$random_variables$etas  
  
## End(Not run)
```

---

create_report	<i>create_report</i>
---------------	----------------------

---

**Description**

Create standard report for results  
The report will be an html created at specified path.

**Usage**

```
create_report(results, path)
```

**Arguments**

results	(Results) Results for which to create report
path	(str) Path to report file

---

create_results	<i>create_results</i>
----------------	-----------------------

---

**Description**

Create/recalculate results object given path to run directory

**Usage**

```
create_results(path, ...)
```

**Arguments**

path (str) Path to run directory  
 ... Arguments to pass to tool specific create results function

**Value**

(Results) Results object for tool

**See Also**

read\_results

**Examples**

```
## Not run:
res <- create_results("frem_dir1")

## End(Not run)
```

---

create_rng	<i>create_rng</i>
------------	-------------------

---

**Description**

Create a new random number generator

Pharmpy functions that use random sampling take a random number generator or seed as input. This function can be used to create a default new random number generator.

**Usage**

```
create_rng(seed = NULL)
```

**Arguments**

seed (numeric (optional)) Seed for the random number generator or NULL (default) for a randomized seed. If seed is generator it will be passed through.

**Value**

(Generator) Initialized numpy random number generator object

**Examples**

```
## Not run:
rng <- create_rng(23)
rng$standard_normal()

## End(Not run)
```

---

create_symbol	<i>create_symbol</i>
---------------	----------------------

---

**Description**

Create a new unique variable symbol given a model

**Usage**

```
create_symbol(model, stem, force_numbering = FALSE)
```

**Arguments**

model	(Model) PharmPy model object
stem	(str) First part of the new variable name
force_numbering	(logical) Forces addition of number to name even if variable does not exist, e.g. COVEFF → COVEFF1

**Value**

(Symbol) Created symbol with unique name

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
create_symbol(model, "TEMP")  
create_symbol(model, "TEMP", force_numbering=TRUE)  
create_symbol(model, "CL")  
  
## End(Not run)
```

---

deidentify_data	<i>deidentify_data</i>
-----------------	------------------------

---

**Description**

Deidentify a dataset

Two operations are performed on the dataset:

1. All ID numbers are randomized from the range 1 to n
2. All columns containing dates will have the year changed

The year change is done by letting the earliest year in the dataset be used as a reference and by maintaining leap years. The reference year will either be 1901, 1902, 1903 or 1904 depending on its distance to the closest preceding leap year.

**Usage**

```
deidentify_data(df, id_column = "ID", date_columns = NULL)
```

**Arguments**

df (data.frame) A dataset  
id\_column (str) Name of the id column  
date\_columns (array(str) (optional)) Names of all date columns

**Value**

(data.frame) Deidentified dataset

---

display_odes	<i>display_odes</i>
--------------	---------------------

---

**Description**

Displays the ordinary differential equation system

**Usage**

```
display_odes(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(ODEDisplayer) A displayable object

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
display_odes(model)  
  
## End(Not run)
```



---

drop_columns	<i>drop_columns</i>
--------------	---------------------

---

## Description

Drop columns from the dataset or mark as dropped

## Usage

```
drop_columns(model, column_names, mark = FALSE)
```

## Arguments

model	(Model) Pharmpy model object
column_names	(array(str) or str) List of column names or one column name to drop or mark as dropped
mark	(logical) Default is to remove column from dataset. Set this to TRUE to only mark as dropped

## Value

(Model) Pharmpy model object

## See Also

drop\_dropped\_columns : Drop all columns marked as drop

undrop\_columns : Undrop columns of model

## Examples

```
## Not run:  
model <- load_example_model("pheno")  
model <- drop_columns(model, c('WGT', 'APGR'))  
vector(model$dataset$columns)  
  
## End(Not run)
```

---

drop\_dropped\_columns    *drop\_dropped\_columns*

---

### Description

Drop columns marked as dropped from the dataset

NM-TRAN date columns will not be dropped by this function even if marked as dropped. Columns not specified in the datainfo (\$INPUT for NONMEM) will also be dropped from the dataset.

### Usage

```
drop_dropped_columns(model)
```

### Arguments

model                    (Model) PharmPy model object

### Value

(Model) PharmPy model object

### See Also

drop\_columns : Drop specific columns or mark them as drop

### Examples

```
## Not run:
model <- load_example_model("pheno")
model <- drop_dropped_columns(model)
vector(model$dataset$columns)

## End(Not run)
```

---

evaluate\_epsilon\_gradient  
                          *evaluate\_epsilon\_gradient*

---

### Description

Evaluate the numeric epsilon gradient

The gradient is evaluated at the current model parameter values or optionally at the given parameter values. The gradient is done for each data record in the model dataset or optionally using the dataset argument. The gradient is done at the current eta values or optionally at the given eta values.

This function currently only support models without ODE systems

**Usage**

```
evaluate_epsilon_gradient(  
  model,  
  etas = NULL,  
  parameters = NULL,  
  dataset = NULL  
)
```

**Arguments**

model	(Model) PharmPy model
etas	(data.frame (optional)) Optional list of eta values
parameters	(list(str=numeric) (optional)) Optional list of parameters and values
dataset	(data.frame (optional)) Optional dataset

**Value**

(data.frame) Gradient

**See Also**

evaluate\_eta\_gradient : Evaluate the eta gradient

**Examples**

```
## Not run:  
model <- load_example_model("pheno_linear")  
results <- load_example_model_fit_results("pheno_linear")  
etas <- results$individual_estimates  
evaluate_epsilon_gradient(model, etas=etas)  
  
## End(Not run)
```

---

evaluate\_eta\_gradient *evaluate\_eta\_gradient*

---

**Description**

Evaluate the numeric eta gradient

The gradient is evaluated at the current model parameter values or optionally at the given parameter values. The gradient is done for each data record in the model dataset or optionally using the dataset argument. The gradient is done at the current eta values or optionally at the given eta values.

This function currently only support models without ODE systems

**Usage**

```
evaluate_eta_gradient(model, etas = NULL, parameters = NULL, dataset = NULL)
```

**Arguments**

model	(Model) Pharmpy model
etas	(data.frame (optional)) Optional list of eta values
parameters	(list(str=numeric) (optional)) Optional list of parameters and values
dataset	(data.frame (optional)) Optional dataset

**Value**

(data.frame) Gradient

**See Also**

evaluate\_epsilon\_gradient : Evaluate the epsilon gradient

**Examples**

```
## Not run:
model <- load_example_model("pheno_linear")
results <- load_example_modelfit_results("pheno_linear")
etas <- results$individual_estimates
evaluate_eta_gradient(model, etas=etas)

## End(Not run)
```

---

evaluate\_expression    *evaluate\_expression*

---

**Description**

Evaluate expression using model

Calculate the value of expression for each data record. The expression can contain dataset columns, variables in model and population parameters. If the model has parameter estimates these will be used. Initial estimates will be used for non-estimated parameters.

**Usage**

```
evaluate_expression(model, expression, parameter_estimates = NULL)
```

**Arguments**

model	(Model) Pharmpy model
expression	(str) Expression to evaluate
parameter_estimates	(list(str=numeric) (optional)) Parameter estimates to use instead of initial estimates

**Value**

(data.frame) A series of one evaluated value for each data record

**Examples**

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_modelfit_results("pheno")
pe <- results$parameter_estimates
evaluate_expression(model, "TVCL*1000", parameter_estimates=pe)

## End(Not run)
```

---

```
evaluate_individual_prediction
      evaluate_individual_prediction
```

---

**Description**

Evaluate the numeric individual prediction

The prediction is evaluated at the current model parameter values or optionally at the given parameter values. The evaluation is done for each data record in the model dataset or optionally using the dataset argument. The evaluation is done at the current eta values or optionally at the given eta values.

This function currently only support models without ODE systems

**Usage**

```
evaluate_individual_prediction(
  model,
  etas = NULL,
  parameters = NULL,
  dataset = NULL
)
```

**Arguments**

model	(Model) PharmPy model
etas	(data.frame (optional)) Optional list of eta values
parameters	(list(str=numeric) (optional)) Optional list of parameters and values
dataset	(data.frame (optional)) Optional dataset

**Value**

(data.frame) Individual predictions

**See Also**

evaluate\_population\_prediction : Evaluate the population prediction

**Examples**

```
## Not run:
model <- load_example_model("pheno_linear")
results <- load_example_modelfit_results("pheno_linear")
etas <- results$individual_estimates
evaluate_individual_prediction(model, etas=etas)

## End(Not run)
```

---

```
evaluate_population_prediction
      evaluate_population_prediction
```

---

**Description**

Evaluate the numeric population prediction

The prediction is evaluated at the current model parameter values or optionally at the given parameter values. The evaluation is done for each data record in the model dataset or optionally using the dataset argument.

This function currently only support models without ODE systems

**Usage**

```
evaluate_population_prediction(model, parameters = NULL, dataset = NULL)
```

**Arguments**

model	(Model) PharmPy model
parameters	(list(str=numeric) (optional)) Optional list of parameters and values
dataset	(data.frame (optional)) Optional dataset

**Value**

(data.frame) Population predictions

**See Also**

evaluate\_individual\_prediction : Evaluate the individual prediction

## Examples

```
## Not run:
model <- load_example_model("pheno_linear")
results <- load_example_model_fit_results("pheno_linear")
pe <- results$parameter_estimates
evaluate_population_prediction(model, parameters=list(pe))

## End(Not run)
```

---

```
evaluate_weighted_residuals
      evaluate_weighted_residuals
```

---

## Description

Evaluate the weighted residuals

The residuals is evaluated at the current model parameter values or optionally at the given parameter values. The residuals is done for each data record in the model dataset or optionally using the dataset argument.

This function currently only support models without ODE systems

## Usage

```
evaluate_weighted_residuals(model, parameters = NULL, dataset = NULL)
```

## Arguments

model	(Model) PharmPy model
parameters	(list(str=numeric) (optional)) Optional list of parameters and values
dataset	(data.frame (optional)) Optional dataset

## Value

(data.frame) WRES

## Examples

```
## Not run:
model <- load_example_model("pheno_linear")
results <- load_example_model_fit_results("pheno_linear")
parameters <- results$parameter_estimates
evaluate_weighted_residuals(model, parameters=list(parameters))

## End(Not run)
```

expand\_additional\_doses  
*expand\_additional\_doses*

---

**Description**

Expand additional doses into separate dose records

**Usage**

```
expand_additional_doses(model, flag = FALSE)
```

**Arguments**

model	(Model) PharmPy model object
flag	(logical) TRUE to add a boolean EXPANDED column to mark added records. In this case all columns in the original dataset will be kept. Care needs to be taken to handle the new dataset.

**Value**

(Model) PharmPy model object

---

find\_clearance\_parameters  
*find\_clearance\_parameters*

---

**Description**

Find clearance parameters in model

**Usage**

```
find_clearance_parameters(model)
```

**Arguments**

model	(Model) PharmPy model
-------	-----------------------

**Value**

(vector) A vector of clearance parameters



### Examples

```
## Not run:  
model <- load_example_model("pheno")  
find_clearance_parameters(model)  
  
## End(Not run)
```

---

```
find_volume_parameters  
                          find_volume_parameters
```

---

### Description

Find volume parameters in model

### Usage

```
find_volume_parameters(model)
```

### Arguments

model                    (Model) PharmPy model

### Value

(vector) A vector of volume parameters

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
find_volume_parameters(model)  
  
## End(Not run)
```

fit

*fit*

---

**Description**

Fit models.

**Usage**

```
fit(model_or_models, tool = NULL)
```

**Arguments**

model\_or\_models

(Model or array(Model)) List of models or one single model

tool

(str (optional)) Estimation tool to use. NULL to use default

**Value**

(ModelfitResults | vector of ModelfitResults) ModelfitResults for the model or models

**See Also**

run\_tool

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
results <- fit(model)  
  
## End(Not run)
```

---

fix\_or\_unfix\_parameters*fix\_or\_unfix\_parameters*

---

**Description**

Fix or unfix parameters

Set fixedness of parameters to specified values

**Usage**

```
fix_or_unfix_parameters(model, parameters)
```

**Arguments**

model (Model) PharmPy model  
 parameters (list(str=logical)) Set fix/unfix for these parameters

**Value**

(Model) PharmPy model object

**See Also**

fix\_parameters : Fix parameters  
 unfix\_parameters : Unfixing parameters  
 fix\_parameters\_to : Fixing parameters and setting a new initial estimate in the same function  
 unfix\_parameters\_to : Unfixing parameters and setting a new initial estimate in the same function

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model$parameters['PTVCL']
model <- fix_or_unfix_parameters(model, list('PTVCL'=TRUE))
model$parameters['PTVCL']

## End(Not run)
```

---

<code>fix_parameters</code>	<i>fix_parameters</i>
-----------------------------	-----------------------

---

**Description**

Fix parameters  
 Fix all listed parameters

**Usage**

```
fix_parameters(model, parameter_names)
```

**Arguments**

model (Model) PharmPy model  
 parameter\_names (array(str) or str) one parameter name or a vector of parameter names

**Value**

(Model) Pharmpy model object

**See Also**

fix\_or\_unfix\_parameters : Fix or unfix parameters (given boolean)

fix\_parameters\_to : Fixing and setting parameter initial estimates in the same function

unfix\_paramaters : Unfixing parameters

unfix\_paramaters\_to : Unfixing parameters and setting a new initial estimate in the same function

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model$parameters['PTVCL']
model <- fix_parameters(model, 'PTVCL')
model$parameters['PTVCL']

## End(Not run)
```

---

fix\_parameters\_to      *fix\_parameters\_to*

---

**Description**

Fix parameters to

Fix all listed parameters to specified value/values

**Usage**

```
fix_parameters_to(model, inits)
```

**Arguments**

model                    (Model) Pharmpy model

inits                    (list(str=numeric)) Inits for all parameters to fix and set init

**Value**

(Model) Pharmpy model object

**See Also**

fix\_parameters : Fix parameters  
fix\_or\_unfix\_parameters : Fix or unfix parameters (given boolean)  
unfix\_paramaters : Unfixing parameters  
unfix\_paramaters\_to : Unfixing parameters and setting a new initial estimate in the same function

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model$parameters['PTVCL']  
model <- fix_parameters_to(model, {'PTVCL': 0.5})  
model$parameters['PTVCL']  
  
## End(Not run)
```

---

get_baselines	<i>get_baselines</i>
---------------	----------------------

---

**Description**

Baselines for each subject.  
Baseline is taken to be the first row even if that has a missing value.

**Usage**

```
get_baselines(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(data.frame) Dataset with the baselines

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_baselines(model)  
  
## End(Not run)
```

---

get\_bioavailability    *get\_bioavailability*

---

**Description**

Get bioavailability of doses for all compartments

**Usage**

```
get_bioavailability(model)
```

**Arguments**

model                    (Model) PharmPy model

**Value**

(list) Dictionary from compartment name to bioavailability expression

---

get\_cmt                    *get\_cmt*

---

**Description**

Get the cmt (compartment) column from the model dataset

If a cmt column is present this will be extracted otherwise a cmt column will be created.

**Usage**

```
get_cmt(model)
```

**Arguments**

model                    (Model) PharmPy model

**Value**

(data.frame) CMT

---

```
get_concentration_parameters_from_data  
    get_concentration_parameters_from_data
```

---

**Description**

Create a dataframe with concentration parameters  
Note that all values are directly calculated from the dataset

**Usage**

```
get_concentration_parameters_from_data(model)
```

**Arguments**

model (Model) PharmPy model object

**Value**

(data.frame) Concentration parameters

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_concentration_parameters_from_data(model)  
  
## End(Not run)
```

---

```
get_config_path    get_config_path
```

---

**Description**

Returns path to the user config path

**Usage**

```
get_config_path()
```

**Value**

(str or NULL) Path to user config or NULL if file does not exist

**Examples**

```
## Not run:  
get_config_path()  
  
## End(Not run)
```

---

```
get_covariate_baselines  
    get_covariate_baselines
```

---

**Description**

Return a dataframe with baselines of all covariates for each id.  
Baseline is taken to be the first row even if that has a missing value.

**Usage**

```
get_covariate_baselines(model)
```

**Arguments**

model                    (Model) PharmPy model

**Value**

(data.frame) covariate baselines

**See Also**

get\_baselines : baselines for all data columns

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- set_covariates(model, c("WGT", "APGR"))  
get_covariate_baselines(model)  
  
## End(Not run)
```



---

get_doseid	<i>get_doseid</i>
------------	-------------------

---

**Description**

Get a DOSEID series from the dataset with an id of each dose period starting from 1

If a dose and observation exist at the same time point the observation will be counted towards the previous dose.

**Usage**

```
get_doseid(model)
```

**Arguments**

model (Model) Pharmpy model

**Value**

(data.frame) DOSEIDs

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_doseid(model)  
  
## End(Not run)
```

---

get_doses	<i>get_doses</i>
-----------	------------------

---

**Description**

Get a series of all doses

Indexed with ID and TIME

**Usage**

```
get_doses(model)
```

**Arguments**

model (Model) Pharmpy model

**Value**

(data.frame) doses

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_doses(model)  
  
## End(Not run)
```

---

get_dv_symbol	<i>get_dv_symbol</i>
---------------	----------------------

---

**Description**

Get the symbol for a certain dvid or dv and check that it is valid

**Usage**

```
get_dv_symbol(model, dv = NULL)
```

**Arguments**

model	(Model) PharmPy model
dv	(str or numeric (optional)) Either a dv symbol, str or dvid. If NULL (default) return the only or first dv.

**Value**

(sympy.Symbol) DV symbol

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_dv_symbol(model, "Y")  
get_dv_symbol(model, 1)  
  
## End(Not run)
```

---

get_evid	<i>get_evid</i>
----------	-----------------

---

**Description**

Get the evid from model dataset

If an event column is present this will be extracted otherwise an evid column will be created.

**Usage**

```
get_evid(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(data.frame) EVID

---

get_ids	<i>get_ids</i>
---------	----------------

---

**Description**

Retrieve a vector of all subject ids of the dataset

**Usage**

```
get_ids(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(vector) All subject ids

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_ids(model)  
  
## End(Not run)
```

```
get_individual_parameters  
    get_individual_parameters
```

---

### Description

Retrieves all parameters with IIV or IOV in `:class:pharmpy.model`.

### Usage

```
get_individual_parameters(model, level = "all")
```

### Arguments

`model` (Model) Pharmpy model to retrieve the individuals parameters from  
`level` (str) The variability level to look for: 'iiv', 'ioi', or 'all' (default)

### Value

(vector`str`) A vector of the parameter names as strings

### See Also

```
get_pk_parameters  
get_rv_parameters  
has_random_effect
```

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
get_individual_parameters(model)  
get_individual_parameters(model, 'iiv')  
get_individual_parameters(model, 'ioi')  
  
## End(Not run)
```

---

`get_individual_prediction_expression`  
*get\_individual\_prediction\_expression*

---

**Description**

Get the full symbolic expression for the modelled individual prediction  
This function currently only support models without ODE systems

**Usage**

```
get_individual_prediction_expression(model)
```

**Arguments**

model (Model) PharmPy model object

**Value**

(Expression) Symbolic expression

**See Also**

`get_population_prediction_expression` : Get full symbolic epression for the population prediction

**Examples**

```
## Not run:  
model <- load_example_model("pheno_linear")  
get_individual_prediction_expression(model)  
  
## End(Not run)
```

---

`get_initial_conditions`  
*get\_initial\_conditions*

---

**Description**

Get initial conditions for the ode system  
Default initial conditions at t=0 for amounts is 0

**Usage**

```
get_initial_conditions(model, dosing = FALSE)
```

**Arguments**

model (Model) PharmPy model  
dosing (logical) Set to TRUE to add dosing as initial conditions

**Value**

(list) Initial conditions

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_initial_conditions(model)  
get_initial_conditions(model, dosing=TRUE)  
  
## End(Not run)
```

---

*get\_lag\_times*                      *get\_lag\_times*

---

**Description**

Get lag times for all compartments

**Usage**

```
get_lag_times(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(list) Dictionary from compartment name to lag time expression

---

get_mdv	<i>get_mdv</i>
---------	----------------

---

**Description**

Get MDVs from dataset

**Usage**

```
get_mdv(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(data.frame) MDVs

---

get_model_code	<i>get_model_code</i>
----------------	-----------------------

---

**Description**

Get the model code of the underlying model language

**Usage**

```
get_model_code(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(str) Model code

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_model_code(model)  
  
## End(Not run)
```

---

```
get_model_covariates  get_model_covariates
```

---

**Description**

List of covariates used in model

A covariate in the model is here defined to be a data item affecting the model prediction excluding dosing items that are not used in model code.

**Usage**

```
get_model_covariates(model, strings = FALSE)
```

**Arguments**

model	(Model) PharmPy model
strings	(logical) Return strings instead of symbols? FALSE (default) will give symbols

**Value**

(vector) Covariate symbols or names

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_model_covariates(model)  
get_model_covariates(model, strings=TRUE)  
  
## End(Not run)
```

---

```
get_number_of_individuals  
  get_number_of_individuals
```

---

**Description**

Retrieve the number of individuals in the model dataset

**Usage**

```
get_number_of_individuals(model)
```

**Arguments**

model	(Model) PharmPy model
-------	-----------------------



**Value**

(integer) Number of individuals in the model dataset

**Note**

For NONMEM models this is the number of individuals of the active dataset, i.e. after filtering of IGNORE and ACCEPT and removal of individuals with no observations.

**See Also**

*get\_number\_of\_observations* : Get the number of observations in a dataset

*get\_number\_of\_observations\_per\_individual* : Get the number of observations per individual in a dataset

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_number_of_individuals(model)  
  
## End(Not run)
```

---

*get\_number\_of\_observations*  
*get\_number\_of\_observations*

---

**Description**

Retrieve the total number of observations in the model dataset

**Usage**

```
get_number_of_observations(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(integer) Number of observations in the model dataset

**Note**

For NONMEM models this is the number of observations of the active dataset, i.e. after filtering of IGNORE and ACCEPT and removal of individuals with no observations.

**See Also**

`get_number_of_individuals` : Get the number of individuals in a dataset

`get_number_of_observations_per_individual` : Get the number of observations per individual in a dataset

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_number_of_observations(model)  
  
## End(Not run)
```

---

```
get_number_of_observations_per_individual  
    get_number_of_observations_per_individual
```

---

**Description**

Number of observations for each individual

**Usage**

```
get_number_of_observations_per_individual(model)
```

**Arguments**

`model` (Model) PharmPy model

**Value**

(data.frame) Number of observations in the model dataset

**Note**

For NONMEM models this is the individuals and number of observations of the active dataset, i.e. after filtering of IGNORE and ACCEPT and removal of individuals with no observations.

**See Also**

`get_number_of_individuals` : Get the number of individuals in a dataset

`get_number_of_observations_per_individual` : Get the number of observations per individual in a dataset

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_number_of_observations_per_individual(model)  
  
## End(Not run)
```

---

get_observations	<i>get_observations</i>
------------------	-------------------------

---

**Description**

Get observations from dataset

**Usage**

```
get_observations(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(data.frame) Observations indexed over ID and TIME

**See Also**

get\_number\_of\_observations : get the number of observations  
get\_number\_of\_observations\_per\_individual : get the number of observations per individual

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_observations(model)  
  
## End(Not run)
```

```
get_observation_expression  
    get_observation_expression
```

---

**Description**

Get the full symbolic expression for the observation according to the model  
This function currently only support models without ODE systems

**Usage**

```
get_observation_expression(model)
```

**Arguments**

model (Model) PharmPy model object

**Value**

(Expression) Symbolic expression

**Examples**

```
## Not run:  
model <- load_example_model("pheno_linear")  
expr <- get_observation_expression(model)  
sympy$pprint(expr)  
  
## End(Not run)
```

---

```
get_omegas    get_omegas
```

---

**Description**

Get all omegas (variability parameters) of a model

**Usage**

```
get_omegas(model)
```

**Arguments**

model (Model) PharmPy model object

**Value**

(Parameters) A copy of all omega parameters

**See Also**

`get_thetas` : Get theta parameters  
`get_sigmas` : Get sigma parameters

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_omegas(model)  
  
## End(Not run)
```

---

`get_pk_parameters`      *get\_pk\_parameters*

---

**Description**

Retrieves PK parameters in `:class:pharmpy.model`.

**Usage**

```
get_pk_parameters(model, kind = "all")
```

**Arguments**

`model`                    (Model) Pharmpy model to retrieve the PK parameters from  
`kind`                    (str) The type of parameter to retrieve: 'absorption', 'distribution', 'elimination',  
or 'all' (default).

**Value**

(vector`str`) A vector of the PK parameter names of the given model

**See Also**

`get_individual_parameters`  
`get_rv_parameters`

**Examples**

```
## Not run:
model <- load_example_model("pheno")
get_pk_parameters(model)
get_pk_parameters(model, 'absorption')
get_pk_parameters(model, 'distribution')
get_pk_parameters(model, 'elimination')

## End(Not run)
```

---

```
get_population_prediction_expression
      get_population_prediction_expression
```

---

**Description**

Get the full symbolic expression for the modelled population prediction

This function currently only support models without ODE systems

**Usage**

```
get_population_prediction_expression(model)
```

**Arguments**

model (Model) PharmPy model object

**Value**

(Expression) Symbolic expression

**See Also**

`get_individual_prediction_expression` : Get full symbolic expression for the individual prediction

**Examples**

```
## Not run:
model <- load_example_model("pheno_linear")
get_population_prediction_expression(model)

## End(Not run)
```

---

get_rv_parameters	<i>get_rv_parameters</i>
-------------------	--------------------------

---

**Description**

Retrieves parameters in :class:pharmpy.model given a random variable.

**Usage**

```
get_rv_parameters(model, rv)
```

**Arguments**

model	(Model) Pharmpy model to retrieve parameters from
rv	(str) Name of random variable to retrieve

**Value**

(vectorstr) A vector of parameter names for the given random variable

**See Also**

has\_random\_effect  
get\_pk\_parameters  
get\_individual\_parameters

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_rv_parameters(model, 'ETA_1')  
  
## End(Not run)
```

---

get_sigmas	<i>get_sigmas</i>
------------	-------------------

---

**Description**

Get all sigmas (residual error variability parameters) of a model

**Usage**

```
get_sigmas(model)
```

**Arguments**

model (Model) PharmPy model object

**Value**

(Parameters) A copy of all sigma parameters

**See Also**

*get\_thetas* : Get theta parameters

*get\_omegas* : Get omega parameters

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_sigmas(model)  
  
## End(Not run)
```

---

*get\_thetas*

*get\_thetas*

---

**Description**

Get all thetas (structural parameters) of a model

**Usage**

```
get_thetas(model)
```

**Arguments**

model (Model) PharmPy model object

**Value**

(Parameters) A copy of all theta parameters

**See Also**

*get\_omegas* : Get omega parameters

*get\_sigmas* : Get sigma parameters



**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_thetas(model)  
  
## End(Not run)
```

---

get_unit_of	<i>get_unit_of</i>
-------------	--------------------

---

**Description**

Derive the physical unit of a variable in the model

Unit information for the dataset needs to be available. The variable can be defined in the code, a dataset column, a parameter or a random variable.

**Usage**

```
get_unit_of(model, variable)
```

**Arguments**

model	(Model) PharmPy model object
variable	(str) Find physical unit of this variable

**Value**

(unit expression) A sympy physics.units expression

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_unit_of(model, "Y")  
get_unit_of(model, "V")  
get_unit_of(model, "WGT")  
  
## End(Not run)
```

`get_zero_order_inputs` *get\_zero\_order\_inputs*

---

**Description**

Get zero order inputs for all compartments

**Usage**

```
get_zero_order_inputs(model)
```

**Arguments**

`model` (Model) PharmPy model

**Value**

(sympy.Matrix) Vector of inputs

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
get_zero_order_inputs(model)  
  
## End(Not run)
```

---

`greekify_model` *greekify\_model*

---

**Description**

Convert to using greek letters for all population parameters

**Usage**

```
greekify_model(model, named_subscripts = FALSE)
```

**Arguments**

`model` (Model) PharmPy model  
`named_subscripts` (logical) Use previous parameter names as subscripts. Default is to use integer subscripts

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model$statements
model <- greekify_model(cleanup_model(model))
model$statements

## End(Not run)
```

---

has\_additive\_error\_model  
*has\_additive\_error\_model*

---

**Description**

Check if a model has an additive error model

Multiple dependent variables are supported. By default the only (in case of one) or the first (in case of many) dependent variable is going to be checked.

**Usage**

```
has_additive_error_model(model, dv = NULL)
```

**Arguments**

model	(Model) The model to check
dv	(str or numeric (optional)) Name or DVID of dependent variable. NULL for the default (first or only)

**Value**

(logical) TRUE if the model has an additive error model and FALSE otherwise

**See Also**

has\_proportional\_error\_model : Check if a model has a proportional error model

has\_combined\_error\_model : Check if a model has a combined error model

has\_weighted\_error\_model : Check if a model has a weighted error model

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
has_additive_error_model(model)  
  
## End(Not run)
```

---

```
has_combined_error_model  
      has_combined_error_model
```

---

**Description**

Check if a model has a combined additive and proportional error model

Multiple dependent variables are supported. By default the only (in case of one) or the first (in case of many) dependent variable is going to be checked.

**Usage**

```
has_combined_error_model(model, dv = NULL)
```

**Arguments**

model	(Model) The model to check
dv	(str or numeric (optional)) Name or DVID of dependent variable. NULL for the default (first or only)

**Value**

(logical) TRUE if the model has a combined error model and FALSE otherwise

**See Also**

`has_additive_error_model` : Check if a model has an additive error model

`has_proportional_error_model` : Check if a model has a proportional error model

`has_weighted_error_model` : Check if a model has a weighted error model

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
has_combined_error_model(model)  
  
## End(Not run)
```

---

has\_covariate\_effect    *has\_covariate\_effect*

---

### Description

Tests if an instance of :class:pharmpy.model has a given covariate effect.

### Usage

```
has_covariate_effect(model, parameter, covariate)
```

### Arguments

model	(Model) Pharmpy model to check for covariate effect.
parameter	(str) Name of parameter.
covariate	(str) Name of covariate.

### Value

(logical) Whether input model has a covariate effect of the input covariate on the input parameter.

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
has_covariate_effect(model, "CL", "APGR")  
  
## End(Not run)
```

---

has\_first\_order\_elimination  
                          *has\_first\_order\_elimination*

---

### Description

Check if the model describes first order elimination

This function relies on heuristics and will not be able to detect all possible ways of coding the first order elimination.

### Usage

```
has_first_order_elimination(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(logical) TRUE if model has describes first order elimination

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
has_first_order_elimination(model)  
  
## End(Not run)
```

---

*has\_linear\_odes*      *has\_linear\_odes*

---

**Description**

Check if model has a linear ODE system

**Usage**

```
has_linear_odes(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(logical) TRUE if model has an ODE system that is linear

**See Also**

`has_odes`  
`has_linear_odes_with_real_eigenvalues`

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
has_linear_odes(model)  
  
## End(Not run)
```

---

has\_linear\_odes\_with\_real\_eigenvalues  
*has\_linear\_odes\_with\_real\_eigenvalues*

---

**Description**

Check if model has a linear ode system with real eigenvalues

**Usage**

```
has_linear_odes_with_real_eigenvalues(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(logical) TRUE if model has an ODE system that is linear

**See Also**

has\_odes

has\_linear\_odes

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
has_linear_odes_with_real_eigenvalues(model)  
  
## End(Not run)
```

---

has\_michaelis\_menten\_elimination  
*has\_michaelis\_menten\_elimination*

---

**Description**

Check if the model describes Michaelis-Menten elimination

This function relies on heuristics and will not be able to detect all possible ways of coding the Michaelis-Menten elimination.

**Usage**

```
has_michaelis_menten_elimination(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(logical) TRUE if model has describes Michaelis-Menten elimination

**Examples**

```
## Not run:
model <- load_example_model("pheno")
has_michaelis_menten_elimination(model)
model <- set_michaelis_menten_elimination(model)
has_michaelis_menten_elimination(model)

## End(Not run)
```

---

```
has_mixed_mm_fo_elimination
      has_mixed_mm_fo_elimination
```

---

**Description**

Check if the model describes mixed Michaelis-Menten and first order elimination

This function relies on heuristics and will not be able to detect all possible ways of coding the mixed Michaelis-Menten and first order elimination.

**Usage**

```
has_mixed_mm_fo_elimination(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(logical) TRUE if model has describes Michaelis-Menten elimination

**Examples**

```
## Not run:
model <- load_example_model("pheno")
has_mixed_mm_fo_elimination(model)
model <- set_mixed_mm_fo_elimination(model)
has_mixed_mm_fo_elimination(model)

## End(Not run)
```



---

has_odes	<i>has_odes</i>
----------	-----------------

---

**Description**

Check if model has an ODE system

**Usage**

```
has_odes(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(logical) TRUE if model has an ODE system

**See Also**

has\_linear\_odes  
has\_linear\_odes\_with\_real\_eigenvalues

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
has_odes(model)  
  
## End(Not run)
```

---

has_proportional_error_model	<i>has_proportional_error_model</i>
------------------------------	-------------------------------------

---

**Description**

Check if a model has a proportional error model

Multiple dependent variables are supported. By default the only (in case of one) or the first (in case of many) dependent variable is going to be checked.

**Usage**

```
has_proportional_error_model(model, dv = NULL)
```

**Arguments**

`model` (Model) The model to check  
`dv` (str or numeric (optional)) Name or DVID of dependent variable. NULL for the default (first or only)

**Value**

(logical) TRUE if the model has a proportional error model and FALSE otherwise

**See Also**

`has_additive_error_model` : Check if a model has an additive error model  
`has_combined_error_model` : Check if a model has a combined error model  
`has_weighted_error_model` : Check if a model has a weighted error model

**Examples**

```
## Not run:
model <- load_example_model("pheno")
has_proportional_error_model(model)

## End(Not run)
```

---

`has_random_effect`      *has\_random\_effect*

---

**Description**

Decides whether the given parameter of a `:class:pharmpy.model` has a random effect.

**Usage**

```
has_random_effect(model, parameter, level = "all")
```

**Arguments**

`model` (Model) Input Pharmpy model  
`parameter` (str) Input parameter  
`level` (str) The variability level to look for: 'iiv', 'iov', or 'all' (default)

**Value**

(logical) Whether the given parameter has a random effect

**See Also**

get\_individual\_parameters  
get\_rv\_parameters

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
has_random_effect(model, 'S1')  
has_random_effect(model, 'CL', 'iiv')  
has_random_effect(model, 'CL', 'iov')  
  
## End(Not run)
```

---

has\_weighted\_error\_model  
*has\_weighted\_error\_model*

---

**Description**

Check if a model has a weighted error model

**Usage**

```
has_weighted_error_model(model)
```

**Arguments**

model (Model) The model to check

**Value**

(logical) TRUE if the model has a weighted error model and FALSE otherwise

**See Also**

has\_additive\_error\_model : Check if a model has an additive error model  
has\_combined\_error\_model : Check if a model has a combined error model  
has\_proportional\_error\_model : Check if a model has a proportional error model

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
has_weighted_error_model(model)  
  
## End(Not run)
```

---

```
has_zero_order_absorption  
  has_zero_order_absorption
```

---

**Description**

Check if ode system describes a zero order absorption  
currently defined as having Infusion dose with rate not in dataset

**Usage**

```
has_zero_order_absorption(model)
```

**Arguments**

model (Model) Pharnpy model

**Value**

(Model) Reference to same model

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
has_zero_order_absorption(model)  
  
## End(Not run)
```

---

```
has_zero_order_elimination  
  has_zero_order_elimination
```

---

**Description**

Check if the model describes zero-order elimination  
This function relies on heuristics and will not be able to detect all possible ways of coding the zero-order elimination.

**Usage**

```
has_zero_order_elimination(model)
```

**Arguments**

model (Model) Pharmpy model

**Value**

(logical) TRUE if model has describes zero order elimination

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
has_zero_order_elimination(model)  
model <- set_zero_order_elimination(model)  
has_zero_order_elimination(model)  
  
## End(Not run)
```

---

install\_pharmpy      *Install Pharmpy*

---

**Description**

Install the pharmpy-core python package into virtual environment. Uses the same Pharmpy version as pharmr.

**Usage**

```
install_pharmpy(envname = "r-reticulate", method = "auto")
```

**Arguments**

envname (str) name of environment. Default is r-reticulate

method (str) type of environment type (virtualenv, conda). Default is auto (virtualenv is not available on Windows)

---

`install_pharmpy_devel` *Install Pharmpy (with specified version)*

---

### **Description**

Install the pharmpy-core python package into virtual environment.

### **Usage**

```
install_pharmpy_devel(  
    envname = "r-reticulate",  
    method = "auto",  
    version = "same"  
)
```

### **Arguments**

<code>envname</code>	(str) name of environment. Default is r-reticulate
<code>method</code>	(str) type of environment type (virtualenv, conda). Default is auto (virtualenv is not available on Windows)
<code>version</code>	(str) which pharmpy version to use (use 'same' for most cases)

---

`is_linearized` *is\_linearized*

---

### **Description**

Determine if a model is linearized

### **Usage**

```
is_linearized(model)
```

### **Arguments**

<code>model</code>	(Model) Pharmpy model
--------------------	-----------------------

### **Value**

(logical) TRUE if model has been linearized and FALSE otherwise

**Examples**

```
## Not run:  
model1 <- load_example_model("pheno")  
is_linearized(model1)  
model2 <- load_example_model("pheno_linear")  
is_linearized(model2)  
  
## End(Not run)
```

---

is_real	<i>is_real</i>
---------	----------------

---

**Description**

Determine if an expression is real valued given constraints of a model

**Usage**

```
is_real(model, expr)
```

**Arguments**

model	(Model) PharmPy model
expr	(str) Expression to test

**Value**

(logical or NULL) TRUE if expression is real, FALSE if not and NULL if unknown

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
is_real(model, "CL")  
  
## End(Not run)
```

---

```
list_time_varying_covariates  
    list_time_varying_covariates
```

---

**Description**

Return a vector of names of all time varying covariates

**Usage**

```
list_time_varying_covariates(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(vector) Names of all time varying covariates

**See Also**

get\_covariate\_baselines : get baselines for all covariates

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
list_time_varying_covariates(model)  
  
## End(Not run)
```

---

```
load_example_model    load_example_model
```

---

**Description**

Load an example model

Load an example model from models built into PharmPy

**Usage**

```
load_example_model(name)
```

**Arguments**

name (str) Name of the model. Currently available models are "pheno" and "pheno\_linear"



**Value**

(Model) Loaded model object

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model$statements  
  
## End(Not run)
```

---

```
load_example_modelfit_results  
                          load_example_modelfit_results
```

---

**Description**

Load the modelfit results of an example model

Load the modelfit results of an example model built into Pharmpy

**Usage**

```
load_example_modelfit_results(name)
```

**Arguments**

name                   (str) Name of the model. Currently available models are "pheno" and "pheno\_linear"

**Value**

(ModelfitResults) Loaded modelfit results object

**Examples**

```
## Not run:  
results <- load_example_modelfit_results("pheno")  
results$parameter_estimates  
  
## End(Not run)
```

---

make\_declarative      *make\_declarative*

---

### Description

Make the model statements declarative  
Each symbol will only be declared once.

### Usage

```
make_declarative(model)
```

### Arguments

model                    (Model) Pharmpy model

### Value

(Model) Pharmpy model object

### Examples

```
## Not run:
model <- load_example_model("pheno")
model$statements$before_odes
model <- make_declarative(model)
model$statements$before_odes

## End(Not run)
```

---

mu\_reference\_model      *mu\_reference\_model*

---

### Description

Convert model to use mu-referencing

Mu-referencing an eta is to separately define its actual mu (mean) parameter. For example:  $CL = \theta_1 e^{\eta_1}$  with  $\eta_1$  following a zero-mean normal distribution would give  $\mu_1 = \log\{\theta_1\}$  and  $CL = e^{\mu_1 + \eta_1}$

### Usage

```
mu_reference_model(model)
```

**Arguments**

model (Model) Pharmpy model object

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- mu_reference_model(model)
model$statements$before_odes

## End(Not run)
```

---

omit\_data

*omit\_data*


---

**Description**

Iterate over omissions of a certain group in a dataset. One group is omitted at a time.

**Usage**

```
omit_data(dataset_or_model, group, name_pattern = "omitted_{}")
```

**Arguments**

dataset\_or\_model (data.frame or Model) Dataset or model for which to omit records

group (str) Name of the column to use for grouping

name\_pattern (str) Name to use for generated datasets. A number starting from 1 will be put in the placeholder.

**Value**

(iterator) Iterator yielding tuples of models/dataframes and the omitted group

---

```
plot_individual_predictions
      plot_individual_predictions
```

---

**Description**

Plot DV and predictions grouped on individuals

**Usage**

```
plot_individual_predictions(model, predictions, individuals = NULL)
```

**Arguments**

model	(Model) Previously run PharmPy model.
predictions	(data.frame) One column for each type of prediction
individuals	(array(numeric) (optional)) A vector of individuals to include. NULL for all individuals

**Value**

(alt.Chart) Plot

---

```
plot_iofv_vs_iofv      plot_iofv_vs_iofv
```

---

**Description**

Plot individual OFV of two models against each other

**Usage**

```
plot_iofv_vs_iofv(iofv1, iofv2, name1, name2)
```

**Arguments**

iofv1	(array) Estimated iOFV of the first model
iofv2	(array) Estimated iOFV of the second model
name1	(str) Name of first model
name2	(str) Name of second model

**Value**

(alt.Chart) Scatterplot

---

```
plot_transformed_eta_distributions
    plot_transformed_eta_distributions
```

---

**Description**

Plot transformed eta distributions for all transformed etas

**Usage**

```
plot_transformed_eta_distributions(
    model,
    parameter_estimates,
    individual_estimates
)
```

**Arguments**

model (Model) Previously run PharmPy model.  
parameter\_estimates (array or list(str=numeric)) Parameter estimates of model fit  
individual\_estimates (data.frame) Individual estimates for etas

**Value**

(alt.Chart) Plot

---

```
predict_influential_individuals
    predict_influential_individuals
```

---

**Description**

Predict influential individuals for a model using a machine learning model.

**Usage**

```
predict_influential_individuals(model, results, cutoff = 3.84)
```

**Arguments**

model (Model) PharmPy model  
results (ModelfitResults) Results for model  
cutoff (numeric) Cutoff threshold for a dofv signalling an influential individual

**Value**

(data.frame) Dataframe over the individuals with a `dofv` column containing the raw predicted delta-OFV and an `influential` column with a boolean to tell whether the individual is influential or not.

**See Also**

`predict_influential_outliers`  
`predict_outliers`

---

`predict_influential_outliers`  
*predict\_influential\_outliers*

---

**Description**

Predict influential outliers for a model using a machine learning model.

**Usage**

```
predict_influential_outliers(  
  model,  
  results,  
  outlier_cutoff = 3,  
  influential_cutoff = 3.84  
)
```

**Arguments**

`model` (Model) Pharmpy model  
`results` (ModelfitResults) Results for model  
`outlier_cutoff` (numeric) Cutoff threshold for a residual signalling an outlier  
`influential_cutoff` (numeric) Cutoff threshold for a `dofv` signalling an influential individual

**Value**

(data.frame) Dataframe over the individuals with a `outliers` and `dofv` columns containing the raw predictions and `influential`, `outlier` and `influential_outlier` boolean columns.

**See Also**

`predict_influential_individuals`  
`predict_outliers`

---

predict_outliers	<i>predict_outliers</i>
------------------	-------------------------

---

## Description

Predict outliers for a model using a machine learning model.

See the [:ref:simeval <Individual OFV summary>](#) documentation for a definition of the residual

## Usage

```
predict_outliers(model, results, cutoff = 3)
```

## Arguments

model	(Model) Pharnpy model
results	(ModelfitResults) ModelfitResults for the model
cutoff	(numeric) Cutoff threshold for a residual signalling an outlier

## Value

(data.frame) Dataframe over the individuals with a residual column containing the raw predicted residuals and a outlier column with a boolean to tell whether the individual is an outlier or not.

## See Also

[predict\\_influential\\_individuals](#)  
[predict\\_influential\\_outliers](#)

## Examples

```
## Not run:  
model <- load_example_model("pheno")  
results <- model$modelfit_results  
predict_outliers(model, results)  
  
## End(Not run)
```

---

`print_fit_summary`      *print\_fit\_summary*

---

**Description**

Print a summary of the model fit

**Usage**

```
print_fit_summary(model)
```

**Arguments**

`model`              (Model) PharmPy model object

---

`print_model_code`      *print\_model\_code*

---

**Description**

Print the model code of the underlying model language

**Usage**

```
print_model_code(model)
```

**Arguments**

`model`              (Model) PharmPy model

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
print_model_code(model)  
  
## End(Not run)
```



---

print\_model\_symbols    *print\_model\_symbols*

---

**Description**

Print all symbols defined in a model

Symbols will be in one of the categories thetas, etas, omegas, epsilons, sigmas, variables and data columns

**Usage**

```
print_model_symbols(model)
```

**Arguments**

model                    (Model) PharmPy model object

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
print_model_symbols(model)  
  
## End(Not run)
```

---

print\_pharmPy\_version    *Print pharmPy version*

---

**Description**

Print the pharmPy version pharMr uses.

**Usage**

```
print_pharmPy_version()
```

---

rank_models	<i>rank_models</i>
-------------	--------------------

---

**Description**

Ranks a vector of models

Ranks a vector of models with a given ranking function

**Usage**

```
rank_models(
  base_model,
  models,
  errors_allowed = NULL,
  rank_type = "ofv",
  cutoff = NULL,
  bic_type = "mixed"
)
```

**Arguments**

base_model	(Model) Base model to compare to
models	(array(Model)) List of models
errors_allowed	(array(str) (optional)) List of errors that are allowed for ranking. Currently available is: rounding_errors and maxevals_exceeded. Default is NULL
rank_type	(str) Name of ranking type. Available options are 'ofv', 'aic', 'bic', 'lrt' (OFV with LRT)
cutoff	(numeric (optional)) Value to use as cutoff. If using LRT, cutoff denotes p-value. Default is NULL
bic_type	(str) Type of BIC to calculate. Default is the mixed effects.

**Value**

(data.frame) DataFrame of the ranked models

**Examples**

```
## Not run:
model_1 <- load_example_model("pheno")
model_2 <- load_example_model("pheno_linear")
rank_models(model_1, c(model_2),
  errors_allowed=c('rounding_errors'),
  rank_type='lrt')

## End(Not run)
```

---

read\_dataset\_from\_datainfo  
*read\_dataset\_from\_datainfo*

---

**Description**

Read a dataset given a datainfo object or path to a datainfo file

**Usage**

```
read_dataset_from_datainfo(datainfo, datatype = NULL)
```

**Arguments**

datainfo (DataInfo or str) A datainfo object or a path to a datainfo object  
datatype (str (optional)) A string to specify dataset type

**Value**

(data.frame) The dataset

---

read\_model *read\_model*

---

**Description**

Read model from file

**Usage**

```
read_model(path)
```

**Arguments**

path (str) Path to model

**Value**

(Model) Read model object

**See Also**

read\_model\_from\_database : Read model from database  
read\_model\_from\_string : Read model from string

**Examples**

```
## Not run:  
model <- read_model("/home/run1$mod")  
  
## End(Not run)
```

---

read\_modelfit\_results *read\_modelfit\_results*

---

**Description**

Read results from external tool for a model

**Usage**

```
read_modelfit_results(path)
```

**Arguments**

path (str) Path to model file

**Value**

(ModelfitResults) Results object

---

read\_model\_from\_string  
*read\_model\_from\_string*

---

**Description**

Read model from the model code in a string

**Usage**

```
read_model_from_string(code)
```

**Arguments**

code (str) Model code to read

**Value**

(Model) Pharmpy model object

**See Also**

read\_model : Read model from file  
read\_model\_from\_database : Read model from database

**Examples**

```
## Not run:  
s <- "$PROBLEM  
$INPUT ID DV TIME  
$DATA file$csv  
$PRED  
Y=THETA(1)+ETA(1)+ERR(1)  
$THETA 1  
$OMEGA 0.1  
$SIGMA 1  
$ESTIMATION METHOD=1"  
read_model_from_string(s)  
  
## End(Not run)
```

---

read\_results

*read\_results*

---

**Description**

Read results object from file

**Usage**

```
read_results(path)
```

**Arguments**

path (str) Path to results file

**Value**

(Results) Results object for tool

**See Also**

create\_results

**Examples**

```
## Not run:  
res <- read_results("results$json")  
  
## End(Not run)
```

---

remove\_covariance\_step  
*remove\_covariance\_step*

---

**Description**

Removes covariance step to the final estimation step

**Usage**

```
remove_covariance_step(model)
```

**Arguments**

model            (Model) PharmPy model

**Value**

(Model) PharmPy model object

**See Also**

```
add_estimation_step  
set_estimation_step  
remove_estimation_step  
append_estimation_step_options  
add_covariance_step  
set_evaluation_step
```

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- remove_covariance_step(model)  
ests <- model$estimation_steps  
ests[1]  
  
## End(Not run)
```

---

```
remove_covariate_effect  
    remove_covariate_effect
```

---

**Description**

Remove a covariate effect from an instance of :class:pharmpy.model.

**Usage**

```
remove_covariate_effect(model, parameter, covariate)
```

**Arguments**

model	(Model) Pharmpy model from which to remove the covariate effect.
parameter	(str) Name of parameter.
covariate	(str) Name of covariate.

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
has_covariate_effect(model, "CL", "WGT")  
model <- remove_covariate_effect(model, "CL", "WGT")  
has_covariate_effect(model, "CL", "WGT")  
  
## End(Not run)
```

---

```
remove_error_model    remove_error_model
```

---

**Description**

Remove error model.

**Usage**

```
remove_error_model(model)
```

**Arguments**

model	(Model) Remove error model for this model
-------	---

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model$statements$find_assignment("Y")
model <- remove_error_model(model)
model$statements$find_assignment("Y")

## End(Not run)
```

---

remove\_estimation\_step  
*remove\_estimation\_step*

---

**Description**

Remove estimation step

**Usage**

```
remove_estimation_step(model, idx)
```

**Arguments**

model (Model) Pharmpy model  
idx (numeric) index of estimation step to remove (starting from 0)

**Value**

(Model) Pharmpy model object

**See Also**

add\_estimation\_step  
set\_estimation\_step  
append\_estimation\_step\_options  
add\_covariance\_step  
remove\_covariance\_step  
set\_evaluation\_step



**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- remove_estimation_step(model, 0)
ests <- model$estimation_steps
length(ests)

## End(Not run)
```

---

remove\_iiv

*remove\_iiv*


---

**Description**

Removes all IIV etas given a vector with eta names and/or parameter names.

**Usage**

```
remove_iiv(model, to_remove = NULL)
```

**Arguments**

model	(Model) Pharmpy model to create block effect on.
to_remove	(array(str) or str (optional)) Name/names of etas and/or name/names of individual parameters to remove. If NULL, all etas that are IIVs will be removed. NULL is default.

**Value**

(Model) Pharmpy model object

**See Also**

```
remove_iov
add_iiv
add_iov
add_pk_iiv
```

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- remove_iiv(model)
model$statements$find_assignment("CL")
model <- load_example_model("pheno")
model <- remove_iiv(model, "V")
model$statements$find_assignment("V")

## End(Not run)
```

---

remove_iov	<i>remove_iov</i>
------------	-------------------

---

### Description

Removes all IOV etas given a vector with eta names.

### Usage

```
remove_iov(model, to_remove = NULL)
```

### Arguments

model	(Model) Pharmpy model to remove IOV from.
to_remove	(array(str) or str (optional)) Name/names of IOV etas to remove, e.g. 'ETA_IOV_1_1'. If NULL, all etas that are IOVs will be removed. NULL is default.

### Value

(Model) Pharmpy model object

### See Also

add\_iiv

add\_iov

remove\_iiv

add\_pk\_iiv

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
model <- remove_iov(model)  
  
## End(Not run)
```

---

remove_lag_time	<i>remove_lag_time</i>
-----------------	------------------------

---

**Description**

Remove lag time from the dose compartment of model.

**Usage**

```
remove_lag_time(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(Model) PharmPy model object

**See Also**

set\_transit\_compartments

add\_lag\_time

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- remove_lag_time(model)  
  
## End(Not run)
```

---

remove_loq_data	<i>remove_loq_data</i>
-----------------	------------------------

---

**Description**

Remove loq data records from the dataset  
Does nothing if none of the limits is specified.

**Usage**

```
remove_loq_data(model, lloq = NULL, uloq = NULL)
```

**Arguments**

model (Model) Pharmpy model object  
 lloq (numeric (optional)) Lower limit of quantification. Default not specified.  
 uloq (numeric (optional)) Upper limit of quantification. Default not specified.

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- remove_loq_data(model, lloq=10, uloq=40)
length(model$dataset)

## End(Not run)
```

---

```
remove_peripheral_compartment
      remove_peripheral_compartment
```

---

**Description**

Remove a peripheral distribution compartment from model

Initial estimates:

```
=====  

2 :math:{CL} = {CL '}, :math:{QP1} = {CL ' } and :math:{VP1} = {VC ' } * 0.053 :math:{QP1} = ({QP1 ' } + {QP2 ' }) / 2,  

:math:{VP1} = {VP1 ' } + {VP2 ' }=====  

===== n =====
```

**Usage**

```
remove_peripheral_compartment(model)
```

**Arguments**

model (Model) Pharmpy model

**Value**

(Model) Pharmpy model object

**See Also**

```
set_peripheral_compartment
add_peripheral_compartment
```

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- set_peripheral_compartments(model, 2)
model <- remove_peripheral_compartment(model)
model$statements$ode_system

## End(Not run)
```

---

```
remove_unused_parameters_and_rvs
      remove_unused_parameters_and_rvs
```

---

**Description**

Remove any parameters and rvs that are not used in the model statements

**Usage**

```
remove_unused_parameters_and_rvs(model)
```

**Arguments**

model (Model) PharmPy model object

**Value**

(Model) PharmPy model object

---

```
rename_symbols      rename_symbols
```

---

**Description**

Rename symbols in the model  
Make sure that no name clash occur.

**Usage**

```
rename_symbols(model, new_names)
```

**Arguments**

model (Model) PharmPy model object  
new\_names (list(str=str)) From old name or symbol to new name or symbol

**Value**

(Model) Pharmpy model object

---

resample_data	<i>resample_data</i>
---------------	----------------------

---

**Description**

Iterate over resamples of a dataset.

The dataset will be grouped on the group column then groups will be selected randomly with or without replacement to form a new dataset. The groups will be renumbered from 1 and upwards to keep them separated in the new dataset.

**Usage**

```
resample_data(
    dataset_or_model,
    group,
    resamples = 1,
    stratify = NULL,
    sample_size = NULL,
    replace = FALSE,
    name_pattern = "resample_{}",
    name = NULL
)
```

**Arguments**

dataset_or_model	(data.frame or Model) Dataset or Model to use
group	(str) Name of column to group by
resamples	(numeric) Number of resamples (iterations) to make
stratify	(str (optional)) Name of column to use for stratification. The values in the stratification column must be equal within a group so that the group can be uniquely determined. A ValueError exception will be raised otherwise.
sample_size	(numeric (optional)) The number of groups that should be sampled. The default is the number of groups. If using stratification the default is to sample using the proportion of the stratas in the dataset. A list of specific sample sizes for each strata can also be supplied.
replace	(logical) A boolean controlling whether sampling should be done with or without replacement
name_pattern	(str) Name to use for generated datasets. A number starting from 1 will be put in the placeholder.
name	(str (optional)) Option to name pattern in case of only one resample

**Value**

(iterator) An iterator yielding tuples of a resampled DataFrame and a vector of resampled groups in order

---

reset_index	<i>Reset index</i>
-------------	--------------------

---

**Description**

Reset index of dataframe.

Reset index from a multi indexed data.frame so that index is added as columns

**Usage**

```
reset_index(df)
```

**Arguments**

df                    A data.frame converted from python using reticulate

---

reset_indices_results	<i>Reset result indices</i>
-----------------------	-----------------------------

---

**Description**

Resets indices in dataframes within Results-objects when needed

**Usage**

```
reset_indices_results(res)
```

**Arguments**

res                    A Pharmpy results object

---

resume_tool	<i>resume_tool</i>
-------------	--------------------

---

**Description**

Resume tool workflow from tool database path

**Usage**

```
resume_tool(path)
```

**Arguments**

path (str) The path to the tool database

**Value**

(Results) Results object for tool

**Examples**

```
## Not run:  
res <- resume_tool("resmod_dir1")  
  
## End(Not run)
```

---

retrieve_final_model	<i>retrieve_final_model</i>
----------------------	-----------------------------

---

**Description**

Retrieve final model from a result object

**Usage**

```
retrieve_final_model(res)
```

**Arguments**

res (Results) A results object

**Value**

(Model) Reference to final model



**See Also**

retrieve\_models

**Examples**

```
## Not run:  
res <- read_results("results$json")  
model <- retrieve_final_model(res)  
  
## End(Not run)
```

---

retrieve\_models      *retrieve\_models*

---

**Description**

Retrieve models after a tool run  
Any models created and run by the tool can be retrieved.

**Usage**

```
retrieve_models(source, names = NULL)
```

**Arguments**

source            (str or Results) Source where to find models. Can be a path (as str or Path), a results object, or a ToolDatabase/ModelDatabase  
names            (array(str) (optional)) List of names of the models to retrieve or NULL for all

**Value**

(vector) List of retrieved model objects

**See Also**

retrieve\_final\_model

**Examples**

```
## Not run:  
tooldir_path <- 'path/to/tool/directory'  
models <- retrieve_models(tooldir_path, names=c('run1'))  
  
## End(Not run)
```

---

run_allometry	<i>run_allometry</i>
---------------	----------------------

---

### Description

Run allometry tool. For more details, see :ref:allometry.

### Usage

```
run_allometry(
    model = NULL,
    results = NULL,
    allometric_variable = "WT",
    reference_value = 70,
    parameters = NULL,
    initials = NULL,
    lower_bounds = NULL,
    upper_bounds = NULL,
    fixed = TRUE,
    ...
)
```

### Arguments

model	(Model (optional)) Pharnpy model
results	(ModelfitResults (optional)) Results for model
allometric_variable	(str) Name of the variable to use for allometric scaling (default is WT)
reference_value	(str or numeric) Reference value for the allometric variable (default is 70)
parameters	(array(str) (optional)) Parameters to apply scaling to (default is all CL, Q and V parameters)
initials	(array(numeric) (optional)) Initial estimates for the exponents. (default is to use 0.75 for CL and Qs and 1 for Vs)
lower_bounds	(array(numeric) (optional)) Lower bounds for the exponents. (default is 0 for all parameters)
upper_bounds	(array(numeric) (optional)) Upper bounds for the exponents. (default is 2 for all parameters)
fixed	(logical) Should the exponents be fixed or not. (default TRUE)
...	Arguments to pass to tool

### Value

(AllometryResults) Allometry tool result object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_modelfit_results("pheno")
run_allometry(model=model, results=results, allometric_variable='WGT')

## End(Not run)
```

run\_amd

*run\_amd***Description**

Run Automatic Model Development (AMD) tool

Runs structural modelsearch, IIV building, and ruvsearch

**Usage**

```
run_amd(
  input,
  results = NULL,
  modeltype = "pk_oral",
  cl_init = 0.01,
  vc_init = 1,
  mat_init = 0.1,
  search_space = NULL,
  lloq_method = NULL,
  lloq_limit = NULL,
  order = NULL,
  allometric_variable = NULL,
  occasion = NULL,
  path = NULL,
  resume = FALSE
)
```

**Arguments**

input	(Model or str) Read model object/Path to a dataset
results	(ModelfitResults (optional)) Results of input if input is a model
modeltype	(str) Type of model to build. Either 'pk_oral' or 'pk_iv'
cl_init	(numeric) Initial estimate for the population clearance
vc_init	(numeric) Initial estimate for the central compartment population volume
mat_init	(numeric) Initial estimate for the mean absorption time (not for iv models)
search_space	(str (optional)) MFL for search space for structural model

lloq_method	(str (optional)) Method for how to remove LOQ data. See transform_blq for vector of available methods
lloq_limit	(str (optional)) Lower limit of quantification. If NULL LLOQ column from dataset will be used
order	(array(str) (optional)) Runorder of components
allometric_variable	(str (optional)) Variable to use for allometry
occasion	(str (optional)) Name of occasion column
path	(str (optional)) Path to run AMD in
resume	(logical (optional)) Whether to allow resuming previous run

**Value**

(Model) Reference to the same model object

**See Also**

run\_iiv  
run\_tool

**Examples**

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_modelfit_results("pheno")
run_amd(model, results=results)

## End(Not run)
```

---

run\_bootstrap

*run\_bootstrap*


---

**Description**

Run bootstrap tool

**Usage**

```
run_bootstrap(model, results = NULL, resamples = 1, ...)
```

**Arguments**

model	(Model) Pharmspy model
results	(ModelfitResults (optional)) Results for model
resamples	(numeric) Number of bootstrap resample
...	Arguments to pass to tool

**Value**

(BootstrapResults) Bootstrap tool result object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_modelfit_results("pheno")
run_bootstrap(model, res, resamples=500)

## End(Not run)
```

---

run_covsearch	<i>run_covsearch</i>
---------------	----------------------

---

**Description**

Run COVsearch tool. For more details, see :ref:covsearch.

**Usage**

```
run_covsearch(
  effects,
  p_forward = 0.05,
  p_backward = 0.01,
  max_steps = -1,
  algorithm = "scm-forward-then-backward",
  results = NULL,
  model = NULL,
  ...
)
```

**Arguments**

effects	(str or array(array(str or array(str)))) MFL of covariate effects to try
p_forward	(numeric) The p-value to use in the likelihood ratio test for forward steps
p_backward	(numeric) The p-value to use in the likelihood ratio test for backward steps
max_steps	(numeric) The maximum number of search steps to make
algorithm	(str) The search algorithm to use. Currently 'scm-forward' and 'scm-forward-then-backward' are supported.
results	(ModelfitResults (optional)) Results of model
model	(Model (optional)) Pharnpy mode
...	Arguments to pass to tool

**Value**

(COVSearchResults) COVsearch tool result object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_modelfit_results("pheno")
effects <- 'COVARIATE(c(CL, V), c(AGE, WT), EXP)'
res <- run_covsearch(effects, model=model, results=results)

## End(Not run)
```

---

run\_estmethod

*run\_estmethod*

---

**Description**

Run estmethod tool.

**Usage**

```
run_estmethod(
  algorithm,
  methods = NULL,
  solvers = NULL,
  results = NULL,
  model = NULL,
  ...
)
```

**Arguments**

algorithm	(str) The algorithm to use (can be 'exhaustive', 'exhaustive_with_update' or 'exhaustive_only_eval')
methods	(array(str) or str (optional)) List of estimation methods to test. Can be specified as 'all', a vector of methods, or NULL (to not test any estimation method)
solvers	(array(str) or str (optional)) List of solver to test. Can be specified as 'all', a vector of solvers, or NULL (to not test any solver)
results	(ModelfitResults (optional)) Results for model
model	(Model (optional)) Pharnpy mode
...	Arguments to pass to tool

**Value**

(EstMethodResults) Estmethod tool result object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_modelfit_results("pheno")
methods <- c('imp', 'saem')
run_estmethod('reduced', methods=methods, solvers='all', results=results, model=model)

## End(Not run)
```

---

run\_iivsearch

*run\_iivsearch*


---

**Description**

Run IIVsearch tool. For more details, see [:ref:iivsearch](#).

**Usage**

```
run_iivsearch(
  algorithm,
  iiv_strategy = "no_add",
  rank_type = "bic",
  cutoff = NULL,
  results = NULL,
  model = NULL,
  ...
)
```

**Arguments**

algorithm	(str) Which algorithm to run (brute_force, brute_force_no_of_etas, brute_force_block_structure)
iiv_strategy	(str) If/how IIV should be added to start model. Possible strategies are 'no_add', 'add_diagonal', or 'fullblock'. Default is 'no_add'
rank_type	(str) Which ranking type should be used (OFV, AIC, BIC). Default is BIC
cutoff	(numeric (optional)) Cutoff for which value of the ranking function that is considered significant. Default is NULL (all models will be ranked)
results	(ModelfitResults (optional)) Results for model
model	(Model (optional)) PharmPy mode
...	Arguments to pass to tool

**Value**

(IIVSearchResults) IIVsearch tool result object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_modelfit_results("pheno")
run_iivsearch('brute_force', results=results, model=model)

## End(Not run)
```

---

run_iovsearch	<i>run_iovsearch</i>
---------------	----------------------

---

**Description**

Run IOVsearch tool. For more details, see `:ref:iovsearch`.

**Usage**

```
run_iovsearch(
  column = "OCC",
  list_of_parameters = NULL,
  rank_type = "bic",
  cutoff = NULL,
  distribution = "same-as-iiv",
  results = NULL,
  model = NULL,
  ...
)
```

**Arguments**

<code>column</code>	(str) Name of column in dataset to use as occasion column (default is 'OCC')
<code>list_of_parameters</code>	(array(str) (optional)) List of parameters to test IOV on, if none all parameters with IIV will be tested (default)
<code>rank_type</code>	(str) Which ranking type should be used (OFV, AIC, BIC). Default is BIC
<code>cutoff</code>	(numeric (optional)) Cutoff for which value of the ranking type that is considered significant. Default is NULL (all models will be ranked)
<code>distribution</code>	(str) Which distribution added IOVs should have (default is same-as-iiv)
<code>results</code>	(ModelfitResults (optional)) Results for model
<code>model</code>	(Model (optional)) Pharmpy mode
<code>...</code>	Arguments to pass to tool

**Value**

(IOVSearchResults) IOVSearch tool result object



**Examples**

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_modelfit_results("pheno")
run_iovsearch('OCC', results=results, model=model)

## End(Not run)
```

---

run_modelfit	<i>run_modelfit</i>
--------------	---------------------

---

**Description**

Run modelfit tool.

**Usage**

```
run_modelfit(model_or_models = NULL, n = NULL, tool = NULL, ...)
```

**Arguments**

model_or_models	(Model or array(Model) (optional)) A vector of models are one single model object
n	(numeric (optional)) Number of models to fit. This is only used if the tool is going to be combined with other tools.
tool	(str (optional)) Which tool to use for fitting. Currently 'nonmem' or 'nlmixr' can be used
...	Arguments to pass to tool

**Value**

(ModelfitResults) Modelfit tool result object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
run_modelfit(model)

## End(Not run)
```

---

run_modelsearch	<i>run_modelsearch</i>
-----------------	------------------------

---

### Description

Run Modelsearch tool. For more details, see [:ref:modelsearch](#).

### Usage

```
run_modelsearch(
  search_space,
  algorithm,
  iiv_strategy = "absorption_delay",
  rank_type = "bic",
  cutoff = NULL,
  results = NULL,
  model = NULL,
  ...
)
```

### Arguments

search_space	(str) Search space to test
algorithm	(str) Algorithm to use (e.g. exhaustive)
iiv_strategy	(str) If/how IIV should be added to candidate models. Possible strategies are 'no_add', 'add_diagonal', 'fullblock', or 'absorption_delay'. Default is 'absorption_delay'
rank_type	(str) Which ranking type should be used (OFV, AIC, BIC). Default is BIC
cutoff	(numeric (optional)) Cutoff for which value of the ranking function that is considered significant. Default is NULL (all models will be ranked)
results	(ModelfitResults (optional)) Results for model
model	(Model (optional)) Pharmpy mode
...	Arguments to pass to tool

### Value

(ModelSearchResults) Modelsearch tool result object

### Examples

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_modelfit_results("pheno")
run_modelsearch('ABSORPTION(Z0);PERIPHERALS(1)', 'exhaustive', results=results, model=model)

## End(Not run)
```

---

run_ruvsearch	<i>run_ruvsearch</i>
---------------	----------------------

---

## Description

Run the ruvsearch tool. For more details, see [:ref:ruvsearch](#).

## Usage

```
run_ruvsearch(  
  model = NULL,  
  results = NULL,  
  groups = 4,  
  p_value = 0.05,  
  skip = NULL,  
  ...  
)
```

## Arguments

model	(Model (optional)) Pharmpy model
results	(ModelfitResults (optional)) Results of model
groups	(numeric) The number of bins to use for the time varying models
p_value	(numeric) The p-value to use for the likelihood ratio test
skip	(array(str) (optional)) A vector of models to not attempt
...	Arguments to pass to tool

## Value

(RUVSearchResults) Ruvsearch tool result object

## Examples

```
## Not run:  
model <- load_example_model("pheno")  
results <- load_example_modelfit_results("pheno")  
run_ruvsearch(model=model, results=results)  
  
## End(Not run)
```

---

run_tool	<i>run_tool</i>
----------	-----------------

---

**Description**

Run tool workflow

**Usage**

```
run_tool(name, ...)
```

**Arguments**

name	(str) Name of tool to run
...	Arguments to pass to tool

**Value**

(Results) Results object for tool

**Examples**

```
## Not run:
model <- load_example_model("pheno")
res <- run_tool("ruvsearch", model)

## End(Not run)
```

---

sample_individual_estimates	<i>sample_individual_estimates</i>
-----------------------------	------------------------------------

---

**Description**

Sample individual estimates given their covariance.

**Usage**

```
sample_individual_estimates(
  model,
  individual_estimates,
  individual_estimates_covariance,
  parameters = NULL,
  samples_per_id = 100,
  rng = NULL
)
```

**Arguments**

**model** (Model) Pharmpy model  
**individual\_estimates** (data.frame) Individual estimates to use  
**individual\_estimates\_covariance** (data.frame) Uncertainty covariance of the individual estimates  
**parameters** (array(str) (optional)) A vector of a subset of individual parameters to sample. Default is NULL, which means all.  
**samples\_per\_id** (numeric) Number of samples per individual  
**rng** (numeric (optional)) Random number generator or seed

**Value**

(data.frame) Pool of samples in a DataFrame

**See Also**

**sample\_parameters\_from\_covariance\_matrix** : Sample parameter vectors using the uncertainty covariance matrix  
**sample\_parameters\_uniformly** : Sample parameter vectors using uniform distribution

**Examples**

```

## Not run:
model <- load_example_model("pheno")
results <- load_example_modelfit_results("pheno")
rng <- create_rng(23)
ie <- results$individual_estimates
iec <- results$individual_estimates_covariance
sample_individual_estimates(model, ie, iec, samples_per_id=2, rng=rng)

## End(Not run)

```

---

```

sample_parameters_from_covariance_matrix
  sample_parameters_from_covariance_matrix

```

---

**Description**

Sample parameter vectors using the covariance matrix  
 If parameters is not provided all estimated parameters will be used

**Usage**

```
sample_parameters_from_covariance_matrix(
  model,
  parameter_estimates,
  covariance_matrix,
  force_posdef_samples = NULL,
  force_posdef_covmatrix = FALSE,
  n = 1,
  rng = NULL
)
```

**Arguments**

`model` (Model) Input model

`parameter_estimates` (array) Parameter estimates to use as means in sampling

`covariance_matrix` (data.frame) Parameter uncertainty covariance matrix

`force_posdef_samples` (numeric (optional)) Set to how many iterations to do before forcing all samples to be positive definite. NULL is default and means never and 0 means always

`force_posdef_covmatrix` (logical) Set to TRUE to force the input covariance matrix to be positive definite

`n` (numeric) Number of samples

`rng` (numeric (optional)) Random number generator

**Value**

(data.frame) A dataframe with one sample per row

**See Also**

`sample_parameters_uniformly` : Sample parameter vectors using uniform distribution

`sample_individual_estimates` : Sample individual estimates given their covariance

**Examples**

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_modelfit_results("pheno")
rng <- create_rng(23)
cov <- results$covariance_matrix
pe <- results$parameter_estimates
sample_parameters_from_covariance_matrix(model, pe, cov, n=3, rng=rng)

## End(Not run)
```

---

```
sample_parameters_uniformly
    sample_parameters_uniformly
```

---

## Description

Sample parameter vectors using uniform sampling

Each parameter value will be randomly sampled from a uniform distribution with the bounds being estimate  $\pm$  estimate \* fraction.

## Usage

```
sample_parameters_uniformly(  
    model,  
    parameter_estimates,  
    fraction = 0.1,  
    force_posdef_samples = NULL,  
    n = 1,  
    rng = NULL  
)
```

## Arguments

model	(Model) PharmPy model
parameter_estimates	(array) Parameter estimates for parameters to use
fraction	(numeric) Fraction of estimate value to use for distribution bounds
force_posdef_samples	(numeric (optional)) Number of samples to reject before forcing variability parameters to give positive definite covariance matrices.
n	(numeric) Number of samples
rng	(numeric (optional)) Random number generator or seed

## Value

(data.frame) samples

## See Also

sample\_parameters\_from\_covariance\_matrix : Sample parameter vectors using the uncertainty covariance matrix

sample\_individual\_estimates : Sample individual estimates given their covariance

**Examples**

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_model_fit_results("pheno")
rng <- create_rng(23)
pe <- results$parameter_estimates
sample_parameters_uniformly(model, pe, n=3, rng=rng)

## End(Not run)
```

---

```
set_additive_error_model
```

```
set_additive_error_model
```

---

**Description**

Set an additive error model. Initial estimate for new sigma is  $(\min(DV)/2)^2$ .

The error function being applied depends on the data transformation. The table displays some examples.

	Data transformation	Additive error
$f + \epsilon_1$		$y$
$\log(f) + \frac{\epsilon_1}{f}$	$\log(y)$	

**Usage**

```
set_additive_error_model(model, dv = NULL, data_trans = NULL, series_terms = 2)
```

**Arguments**

model	(Model) Set error model for this model
dv	(str or numeric (optional)) Name or DVID of dependent variable. NULL for the default (first or only)
data_trans	(str (optional)) A data transformation expression or NULL (default) to use the transformation specified by the model. Series expansion will be used for approximation.
series_terms	(numeric) Number of terms to use for the series expansion approximation for data transformation.

**Value**

(Model) PharmPy model object

**See Also**

set\_proportional\_error\_model : Proportional error model

set\_combined\_error\_model : Combined error model



**Examples**

```
## Not run:
model <- load_example_model("pheno")
model$statements$find_assignment("Y")
model <- set_additive_error_model(model)
model$statements$find_assignment("Y")
model <- load_example_model("pheno")
model$statements$find_assignment("Y")
model <- set_additive_error_model(model, data_trans="log(Y)")
model$statements$find_assignment("Y")

## End(Not run)
```

---

set\_bolus\_absorption    *set\_bolus\_absorption*

---

**Description**

Set or change to bolus absorption rate.

Currently lagtime together with bolus absorption is not supported.

**Usage**

```
set_bolus_absorption(model)
```

**Arguments**

model                    (Model) Model to set or change absorption rate

**Value**

(Model) Pharnpy model object

**See Also**

set\_zero\_order\_absorption

set\_first\_order\_absorption

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- set_bolus_absorption(model)
model$statements$ode_system

## End(Not run)
```

---

```
set_combined_error_model
      set_combined_error_model
```

---

### Description

Set a combined error model. Initial estimates for new sigmas are  $(\min(DV)/2)^2$  for proportional and 0.09 for additive.

The error function being applied depends on the data transformation.

Data transformation	Combined error
$y$	$f + f \cdot \epsilon_1 + \epsilon_2$
$\log(y)$	$\log(f) + \epsilon_1 + \frac{\epsilon_2}{f}$

### Usage

```
set_combined_error_model(model, dv = NULL, data_trans = NULL)
```

### Arguments

model	(Model) Set error model for this model
dv	(str or numeric (optional)) Name or DVID of dependent variable. NULL for the default (first or only)
data_trans	(str (optional)) A data transformation expression or NULL (default) to use the transformation specified by the model.

### Value

(Model) Pharmpy model object

### See Also

set\_additive\_error\_model : Additive error model  
 set\_proportional\_error\_model: Proportional error model

### Examples

```
## Not run:
model <- remove_error_model(load_example_model("pheno"))
model <- set_combined_error_model(model)
model$statements$find_assignment("Y")
model <- remove_error_model(load_example_model("pheno"))
model <- set_combined_error_model(model, data_trans="log(Y)")
model$statements$find_assignment("Y")

## End(Not run)
```

---

set_covariates	<i>set_covariates</i>
----------------	-----------------------

---

**Description**

Set columns in the dataset to be covariates in the datainfo

**Usage**

```
set_covariates(model, covariates)
```

**Arguments**

model	(Model) Pharmpy model
covariates	(array(str)) List of column names

**Value**

(Model) Pharmpy model object

---

set_dtbs_error_model	<i>set_dtbs_error_model</i>
----------------------	-----------------------------

---

**Description**

Dynamic transform both sides

**Usage**

```
set_dtbs_error_model(model, fix_to_log = FALSE)
```

**Arguments**

model	(Model) Pharmpy model
fix_to_log	(logical) Set to TRUE to fix lambda and zeta to 0, i.e. emulating log-transformed data

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- set_dtbs_error_model(model)

## End(Not run)
```

---

set_dvid	<i>set_dvid</i>
----------	-----------------

---

**Description**

Set a column to act as DVID. Replace DVID if one is already set.

**Usage**

```
set_dvid(model, name)
```

**Arguments**

model	(Model) Pharmpy model
name	(str) Name of DVID column

**Value**

(Model) Pharmpy model object

---

set_estimation_step	<i>set_estimation_step</i>
---------------------	----------------------------

---

**Description**

Set estimation step

Sets estimation step for a model. Methods currently supported are: FO, FOCE, ITS, LAPLACE, IMPMAP, IMP, SAEM, BAYES

**Usage**

```
set_estimation_step(model, method, idx = 0, ...)
```

**Arguments**

model	(Model) Pharmpy model
method	(str) estimation method to change to
idx	(numeric) index of estimation step, default is 0 (first estimation step)
...	Arguments to pass to EstimationStep (such as interaction, evaluation)

**Value**

(Model) Pharmpy model object

**See Also**

add\_estimation\_step  
remove\_estimation\_step  
append\_estimation\_step\_options  
add\_covariance\_step  
remove\_covariance\_step  
set\_evaluation\_step

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
opts <- list('NITER'=1000, 'ISAMPLE'=100)  
model <- set_estimation_step(model, "IMP", evaluation=TRUE, tool_options=opts)  
model$estimation_steps[1]  
  
## End(Not run)
```

---

set\_evaluation\_step    *set\_evaluation\_step*

---

**Description**

Set estimation step

Sets estimation step for a model. Methods currently supported are: FO, FOCE, ITS, LAPLACE, IMPMAP, IMP, SAEM, BAYES

**Usage**

```
set_evaluation_step(model, idx = -1)
```

**Arguments**

model            (Model) Pharmpy model  
idx              (numeric) index of estimation step, default is -1 (last estimation step)

**Value**

(Model) Pharmpy model object

**See Also**

```
set_estimation_step
add_estimation_step
remove_estimation_step
append_estimation_step_options
add_covariance_step
remove_covariance_step
```

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- set_evaluation_step(model)
model$estimation_steps[1]

## End(Not run)
```

---

```
set_first_order_absorption
      set_first_order_absorption
```

---

**Description**

Set or change to first order absorption rate.

Initial estimate for absorption rate is set to the previous rate if available, otherwise it is set to the time of first observation/2.

**Usage**

```
set_first_order_absorption(model)
```

**Arguments**

model (Model) Model to set or change to use first order absorption rate

**Value**

(Model) PharmPy model object

**See Also**

```
set_bolus_order_absorption
set_zero_order_absorption
```

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- set_first_order_absorption(model)
model$statements$ode_system

## End(Not run)
```

---

```
set_first_order_elimination
      set_first_order_elimination
```

---

**Description**

Sets elimination to first order

**Usage**

```
set_first_order_elimination(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(Model) PharmPy model object

**See Also**

```
set_zero_order_elimination
set_michaelis_menten_elimination
```

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- set_first_order_elimination(model)
model$statements$ode_system

## End(Not run)
```

---

```
set_iiv_on_ruv      set_iiv_on_ruv
```

---

### Description

Multiplies epsilons with exponential (new) etas.

Initial variance for new etas is 0.09.

### Usage

```
set_iiv_on_ruv(model, list_of_eps = NULL, same_eta = TRUE, eta_names = NULL)
```

### Arguments

model	(Model) Pharmpy model to apply IIV on epsilons.
list_of_eps	(array(str) or str (optional)) Name/names of epsilons to multiply with exponential etas. If NULL, all epsilons will be chosen. NULL is default.
same_eta	(logical) Boolean of whether all RUVs from input should use the same new ETA or if one ETA should be created for each RUV. TRUE is default.
eta_names	(array(str) or str (optional)) Custom names of new etas. Must be equal to the number epsilons or 1 if same eta.

### Value

(Model) Pharmpy model object

### See Also

set\_power\_on\_ruv

### Examples

```
## Not run:
model <- load_example_model("pheno")
model <- set_iiv_on_ruv(model)
model$statements$find_assignment("Y")

## End(Not run)
```



---

set\_initial\_condition *set\_initial\_condition*

---

### Description

Set an initial condition for the ode system

If the initial condition is already set it will be updated. If the initial condition is set to zero at time zero it will be removed (since the default is 0).

### Usage

```
set_initial_condition(model, compartment, expression, time = 0)
```

### Arguments

model	(Model) Pharmpy model
compartment	(str) Name of the compartment
expression	(str or numeric) The expression of the initial condition
time	(str or numeric) Time point. Default 0

### Value

(model) Pharmpy model object

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
model <- set_initial_condition(model, "CENTRAL", 10)  
get_initial_conditions(model)  
  
## End(Not run)
```

---

set\_initial\_estimates *set\_initial\_estimates*

---

### Description

Set initial estimates

### Usage

```
set_initial_estimates(model, inits)
```

**Arguments**

model (Model) Pharmpy model  
inits (list(str=numeric)) A list of parameter init for parameters to change

**Value**

(Model) Pharmpy model object

**See Also**

fix\_parameters\_to : Fixing and setting parameter initial estimates in the same function  
unfix\_paramaters\_to : Unfixing parameters and setting a new initial estimate in the same

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- set_initial_estimates(model, list('PTVCL'=2))  
model$parameters['PTVCL']  
  
## End(Not run)
```

---

set\_lower\_bounds      *set\_lower\_bounds*

---

**Description**

Set parameter lower bounds

**Usage**

```
set_lower_bounds(model, bounds)
```

**Arguments**

model (Model) Pharmpy model  
bounds (list(str=numeric)) A list of parameter bounds for parameters to change

**Value**

(Model) Pharmpy model object

**See Also**

set\_upper\_bounds : Set parameter upper bounds  
unconstrain\_parameters : Remove all constraints of parameters

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- set_lower_bounds(model, {'PTVCL': -10})
model$parameters['PTVCL']

## End(Not run)
```

---

```
set_michaelis_menten_elimination
      set_michaelis_menten_elimination
```

---

**Description**

Sets elimination to Michaelis-Menten.

Initial estimate for CLMM is set to CL and KM is set to  $2 \cdot \max(DV)$ .

**Usage**

```
set_michaelis_menten_elimination(model)
```

**Arguments**

model (Model) PharmPy model

**Value**

(Model) PharmPy model object

**See Also**

set\_first\_order\_elimination

set\_zero\_order\_elimination

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- set_michaelis_menten_elimination(model)
model$statements$ode_system

## End(Not run)
```

---

```
set_mixed_mm_fo_elimination  
    set_mixed_mm_fo_elimination
```

---

### Description

Sets elimination to mixed Michaelis-Menten and first order.

Initial estimate for CLMM is set to  $CL/2$  and KM is set to  $2 \cdot \max(DV)$ .

### Usage

```
set_mixed_mm_fo_elimination(model)
```

### Arguments

model (Model) PharmPy model

### Value

(Model) PharmPy model object

### See Also

set\_first\_order\_elimination

set\_zero\_order\_elimination

set\_michaelis\_menten\_elimination

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
model <- set_mixed_mm_fo_elimination(model)  
model$statements$ode_system  
  
## End(Not run)
```

---

set_name	<i>set_name</i>
----------	-----------------

---

**Description**

Set name of model object

**Usage**

```
set_name(model, new_name)
```

**Arguments**

model	(Model) Pharmpy model
new_name	(str) New name of model

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model$name
model <- set_name(model, "run2")
model$name

## End(Not run)
```

---

set_ode_solver	<i>set_ode_solver</i>
----------------	-----------------------

---

**Description**

Sets ODE solver to use for model

Recognized solvers and their corresponding NONMEM advans:

```
+-----+-----+ | Solver | NONMEM ADVAN | +-----+-----+
| CVODES | ADVAN14 | +-----+-----+ | DGEAR | ADVAN8 | +-----+
-----+-----+ | DVERK | ADVAN6 | +-----+-----+
--+ | IDA | ADVAN15 | +-----+-----+ | LSODA | ADVAN13 | +-----+
-----+-----+ | LSODI | ADVAN9 | +-----+-----+
--+
```

**Usage**

```
set_ode_solver(model, solver)
```

**Arguments**

model	(Model) PharmPy model
solver	(str) Solver to use or NULL for no preference

**Value**

(Model) PharmPy model object

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- set_ode_solver(model, 'LSODA')  
  
## End(Not run)
```

---

```
set_peripheral_compartments  
    set_peripheral_compartments
```

---

**Description**

Sets the number of peripheral compartments to a specified number.

**Usage**

```
set_peripheral_compartments(model, n)
```

**Arguments**

model	(Model) PharmPy model
n	(numeric) Number of transit compartments

**Value**

(Model) PharmPy model object

**See Also**

```
add_peripheral_compartment  
remove_peripheral_compartment
```

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- set_peripheral_compartments(model, 2)
model$statements$ode_system

## End(Not run)
```

---

```
set_power_on_ruv      set_power_on_ruv
```

---

**Description**

Applies a power effect to provided epsilons.

Initial estimates for new thetas are 1 if the error model is proportional, otherwise they are 0.1.

**Usage**

```
set_power_on_ruv(
  model,
  list_of_eps = NULL,
  lower_limit = 0.01,
  ipred = NULL,
  zero_protection = FALSE
)
```

**Arguments**

model	(Model) PharmPy model to create block effect on.
list_of_eps	(str or array (optional)) Name/names of epsilons to apply power effect. If NULL, all epsilons will be used. NULL is default.
lower_limit	(numeric (optional)) Lower limit of power (theta). NULL for no limit.
ipred	(str (optional)) Symbol to use as IPRED. Default is to autodetect expression for IPRED.
zero_protection	(logical) Set to TRUE to add code protecting from IPRED=0

**Value**

(Model) PharmPy model object

**See Also**

set\_iiv\_on\_ruv

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- set_power_on_ruv(model)
model$statements$find_assignment("Y")

## End(Not run)
```

---

```
set_proportional_error_model
      set_proportional_error_model
```

---

**Description**

Set a proportional error model. Initial estimate for new sigma is 0.09.

The error function being applied depends on the data transformation.

+	+	+	Data transformation	Proportional error
	+	+	+	:math:y
:math:f + f \epsilon_1	+	+	+	:math:\log(y)
:math:\log(f) + \epsilon_1	+	+	+	+

**Usage**

```
set_proportional_error_model(
  model,
  dv = NULL,
  data_trans = NULL,
  zero_protection = TRUE
)
```

**Arguments**

model	(Model) Set error model for this model
dv	(str or numeric (optional)) Name or DVID of dependent variable. NULL for the default (first or only)
data_trans	(str (optional)) A data transformation expression or NULL (default) to use the transformation specified by the model.
zero_protection	(logical) Set to TRUE to add code protecting from IPRED=0

**Value**

(Model) PharmPy model object



**See Also**

set\_additive\_error\_model : Additive error model  
set\_combined\_error\_model : Combined error model

**Examples**

```
## Not run:  
model <- remove_error_model(load_example_model("pheno"))  
model <- set_proportional_error_model(model)  
model$statements$after_odes  
model <- remove_error_model(load_example_model("pheno"))  
model <- set_proportional_error_model(  
  model,  
  data_trans="log(Y)"  
model$statements$after_odes  
  
## End(Not run)
```

---

```
set_seq_zo_fo_absorption  
      set_seq_zo_fo_absorption
```

---

**Description**

Set or change to sequential zero order first order absorption rate.

Initial estimate for absorption rate is set the previous rate if available, otherwise it is set to the time of first observation/2.

Currently lagtime together with sequential zero order first order absorption is not supported.

**Usage**

```
set_seq_zo_fo_absorption(model)
```

**Arguments**

model (Model) Model to set or change absorption rate

**Value**

(Model) PharmPy model object

**See Also**

set\_bolus\_order\_absorption  
set\_zero\_order\_absorption  
set\_first\_order\_absorption

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- set_seq_zo_fo_absorption(model)
model$statements$ode_system

## End(Not run)
```

---

```
set_time_varying_error_model
      set_time_varying_error_model
```

---

**Description**

Set a time varying error model per time cutoff

**Usage**

```
set_time_varying_error_model(model, cutoff, idv = "TIME")
```

**Arguments**

model	(Model) PharmPy model
cutoff	(numeric) A value at the given quantile over idv column
idv	(str) Time or time after dose, default is Time

**Value**

(Model) PharmPy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- set_time_varying_error_model(model, cutoff=1.0)
model$statements$find_assignment("Y")

## End(Not run)
```

---

set_tmdd	<i>set_tmdd</i>
----------	-----------------

---

**Description**

Sets target mediated drug disposition  
Sets target mediated drug disposition to a PK model.

**Usage**

```
set_tmdd(model, type)
```

**Arguments**

model	(Model) Pharnpy model
type	(str) Type of TMDD model

**Value**

(Model) Pharnpy model object

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- set_tmdd(model, "full")  
  
## End(Not run)
```

---

set_transit_compartments	<i>set_transit_compartments</i>
--------------------------	---------------------------------

---

**Description**

Set the number of transit compartments of model.  
Initial estimate for absorption rate is set the previous rate if available, otherwise it is set to the time of first observation/2.

**Usage**

```
set_transit_compartments(model, n, keep_depot = TRUE)
```

**Arguments**

model (Model) Pharmpy model  
n (numeric) Number of transit compartments  
keep\_depot (logical) FALSE to convert depot compartment into a transit compartment

**Value**

(Model) Pharmpy model object

**See Also**

add\_lag\_time

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- set_transit_compartments(model, 3)  
model$statements$ode_system  
  
## End(Not run)
```

---

set\_upper\_bounds      *set\_upper\_bounds*

---

**Description**

Set parameter upper bounds

**Usage**

```
set_upper_bounds(model, bounds)
```

**Arguments**

model (Model) Pharmpy model  
bounds (list(str=numeric)) A list of parameter bounds for parameters to change

**Value**

(Model) Pharmpy model object

**See Also**

set\_lower\_bounds : Set parameter lower bounds  
unconstrain\_parameters : Remove all constraints of parameters

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- set_upper_bounds(model, list('PTVCL'=10))  
model$parameters['PTVCL']  
  
## End(Not run)
```

---

```
set_weighted_error_model  
      set_weighted_error_model
```

---

**Description**

Encode error model with one epsilon and W as weight

**Usage**

```
set_weighted_error_model(model)
```

**Arguments**

model            (Model) PharmPy model

**Value**

(Model) PharmPy model object

**See Also**

use\_thetas\_for\_error\_stdev : Use thetas to estimate error

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- set_weighted_error_model(model)  
  
## End(Not run)
```

---

```
set_zero_order_absorption  
    set_zero_order_absorption
```

---

**Description**

Set or change to zero order absorption rate.

Initial estimate for absorption rate is set the previous rate if available, otherwise it is set to the time of first observation/2.

**Usage**

```
set_zero_order_absorption(model)
```

**Arguments**

model (Model) Model to set or change to first order absorption rate

**Value**

(Model) Pharnpy model object

**See Also**

```
set_bolus_order_absorption  
set_first_order_absorption
```

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- set_zero_order_absorption(model)  
model$statements$ode_system  
  
## End(Not run)
```

---

```
set_zero_order_elimination  
    set_zero_order_elimination
```

---

**Description**

Sets elimination to zero order.

Initial estimate for KM is set to 1% of smallest observation.

**Usage**

```
set_zero_order_elimination(model)
```

**Arguments**

model (Model) Pharmpy model

**Value**

(Model) Pharmpy model object

**See Also**

set\_first\_order\_elimination  
set\_michaelis\_menten\_elimination

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- set_zero_order_elimination(model)  
model$statements$ode_system  
  
## End(Not run)
```

---

set\_zero\_order\_input *set\_zero\_order\_input*

---

**Description**

Set a zero order input for the ode system  
If the zero order input is already set it will be updated.

**Usage**

```
set_zero_order_input(model, compartment, expression)
```

**Arguments**

model (Model) Pharmpy model  
compartment (str) Name of the compartment  
expression (str or numeric) The expression of the zero order input

**Value**

(model) Pharmpy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- set_zero_order_input(model, "CENTRAL", 10)
get_zero_order_inputs(model)

## End(Not run)
```

---

simplify\_expression    *simplify\_expression*

---

**Description**

Simplify expression given constraints in model

**Usage**

```
simplify_expression(model, expr)
```

**Arguments**

model	(Model) PharmPy model object
expr	(numeric or str) Expression to simplify

**Value**

(Expression) Simplified expression

**Examples**

```
## Not run:
model <- load_example_model("pheno")
simplify_expression(model, "Abs(PTVCL)")

## End(Not run)
```



---

solve_ode_system	<i>solve_ode_system</i>
------------------	-------------------------

---

**Description**

Replace ODE system with analytical solution if possible

Warnings This function can currently only handle the most simple of ODE systems.

**Usage**

```
solve_ode_system(model)
```

**Arguments**

model	(Model) Pharmpy model object
-------	------------------------------

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model$statements$ode_system
model <- solve_ode_system(model)

## End(Not run)
```

---

split_joint_distribution	<i>split_joint_distribution</i>
--------------------------	---------------------------------

---

**Description**

Splits etas following a joint distribution into separate distributions.

**Usage**

```
split_joint_distribution(model, rvs = NULL)
```

**Arguments**

model	(Model) Pharmpy model
rvs	(array(str) or str (optional)) Name/names of etas to separate. If NULL, all etas that are IIVs and non-fixed will become single. NULL is default.

**Value**

(Model) Pharmpy model object

**See Also**

`create_joint_distribution` : combine etas into a join distribution

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- create_joint_distribution(model, c('ETA_1', 'ETA_2'))
model$random_variables$etas
model <- split_joint_distribution(model, c('ETA_1', 'ETA_2'))
model$random_variables$etas

## End(Not run)
```

---

<code>summarize_errors</code>	<i>summarize_errors</i>
-------------------------------	-------------------------

---

**Description**

Summarize errors and warnings from one or multiple model runs.

Summarize the errors and warnings found after running the model/models.

**Usage**

```
summarize_errors(results)
```

**Arguments**

`results` (ModelfitResults or array(ModelfitResults)) List of ModelfitResults or single ModelfitResults

**Value**

(data.frame) A DataFrame of errors with model name, category (error or warning), and an integer as index, an empty DataFrame if there were no errors or warnings found.

**Examples**

```
## Not run:
model <- load_example_model("pheno")
summarize_errors(model)

## End(Not run)
```

summarize\_individuals summarize\_individuals

**Description**

Creates a summary dataframe keyed by model-individual pairs for an input vector of models.

Content of the various columns:

Column	Description
outlier_count	Number of observations with CWRES > 5
ofv	Individual OFV
dofv_vs_parent	Difference in individual OFV between this model and its parent model
predicted_dofv	Predicted dOFV if this individual was excluded
predicted_residual	Predicted residual

**Usage**

```
summarize_individuals(models)
```

**Arguments**

models (array(Model)) Input models

**Value**

(data.frame | NULL) The summary as a dataframe

**Examples**

```
## Not run:
model <- load_example_model("pheno")
fit_results <- fit(model)
results <- run_tool(
  model=model,
  mfl='ABSORPTION(ZO);PERIPHERALS(c(1, 2))',
  algorithm='reduced_stepwise'
)
summarize_individuals([results$start_model, *results$models])

## End(Not run)
```

---

```
summarize_individuals_count_table
      summarize_individuals_count_table
```

---

### Description

Create a count table for individual data

Content of the various columns:

Column	Description
<code>inf_selection</code>	Number of subjects influential on model selection. $ \mathbf{OFV}_{parent} - \mathbf{OFV}  > 3.84$
<code>inf_params</code>	Number of subjects influential on parameters. $\text{predicted\_dofv} > 3.84$
<code>out_obs</code>	Number of subjects having at least one outlying observation ( $\text{CWRES} > 5$ )
<code>out_ind</code>	Number of outlying subjects. $\text{predicted\_residual} > 3.0$
<code>inf_outlier</code>	Number of subjects both influential by any criteria and outlier by any criteria

### Usage

```
summarize_individuals_count_table(models = NULL, df = NULL)
```

### Arguments

`models` (array(Model) (optional)) List of models to summarize.

`df` (data.frame) Output from a previous call to `summarize_individuals`.

### Value

(data.frame) Table with one row per model.

### See Also

`summarize_individuals` : Get raw individual data

---

```
summarize_modelfit_results  
      summarize_modelfit_results
```

---

## Description

Summarize results of model runs

Summarize different results after fitting a model, includes runtime, ofv, and parameter estimates (with errors). If `include_all_estimation_steps` is `FALSE`, only the last estimation step will be included (note that in that case, the `minimization_successful` value will be referring to the last estimation step, if last step is evaluation it will go backwards until it finds an estimation step that wasn't an evaluation).

## Usage

```
summarize_modelfit_results(results, include_all_estimation_steps = FALSE)
```

## Arguments

`results` (ModelfitResults or array(ModelfitResults)) List of ModelfitResults or single ModelfitResults

`include_all_estimation_steps` (logical) Whether to include all estimation steps, default is `FALSE`

## Value

(data.frame) A DataFrame of modelfit results with model name and estimation step as index.

## Examples

```
## Not run:  
model <- load_example_model("pheno")  
summarize_modelfit_results(model$modelfit_results)  
  
## End(Not run)
```

---

```
transform_blq      transform_blq
```

---

## Description

Transform for BLQ data

Transform a given model, methods available are m1, m3, and m4 (1). Current limits of the m3 and m4 method:

- Does not support covariance between epsilons
- Only supports additive, proportional, and combined error model

(1) Beal SL. Ways to fit a PK model with some data below the quantification limit. *J Pharmacokinetic Pharmacodyn*. 2001 Oct;28(5):481-504. doi: 10.1023/a:1012299115260. Erratum in: *J Pharmacokinetic Pharmacodyn* 2002 Jun;29(3):309. PMID: 11768292.

## Usage

```
transform_blq(model, method = "m4", lloq = NULL)
```

## Arguments

model	(Model) PharmPy model
method	(str) Which BLQ method to use
lloq	(numeric (optional)) LLOQ limit to use, if NULL PharmPy will use the BLQ/LLOQ column in the dataset

## Value

(Model) PharmPy model object

## See Also

`remove_loq_data`

## Examples

```
## Not run:  
model <- load_example_model("pheno")  
model <- transform_blq(model, method='m4', lloq=0.1)  
model$statements$find_assignment("Y")  
  
## End(Not run)
```

---

transform\_etas\_boxcox *transform\_etas\_boxcox*

---

### Description

Applies a boxcox transformation to selected etas

Initial estimate for lambda is 0.1 with bounds (-3, 3).

### Usage

```
transform_etas_boxcox(model, list_of_etas = NULL)
```

### Arguments

`model` (Model) Pharnpy model to apply boxcox transformation to.

`list_of_etas` (array(str) or str (optional)) Name/names of etas to transform. If NULL, all etas will be transformed (default).

### Value

(Model) Pharnpy model object

### See Also

`transform_etas_tdist`

`transform_etas_john_draper`

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
model <- transform_etas_boxcox(model, c("ETA_1"))  
model$statements$before_odes$full_expression("CL")  
  
## End(Not run)
```

---

transform\_etas\_john\_draper  
*transform\_etas\_john\_draper*

---

### Description

Applies a John Draper transformation (1) to spelected etas

Initial estimate for lambda is 0.1 with bounds (-3, 3).

(1) John, J., Draper, N. (1980). An Alternative Family of Transformations. Journal of the Royal Statistical Society. Series C (Applied Statistics), 29(2), 190-197. doi:10.2307/2986305

### Usage

```
transform_etas_john_draper(model, list_of_etas = NULL)
```

### Arguments

model	(Model) Pharmpy model to apply John Draper transformation to.
list_of_etas	(array(str) or str (optional)) Name/names of etas to transform. If NULL, all etas will be transformed (default).

### Value

(Model) Pharmpy model object

### See Also

transform\_etas\_boxcox  
transform\_etas\_tdist

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
model <- transform_etas_john_draper(model, c("ETA_1"))  
model$statements$before_odes$full_expression("CL")  
  
## End(Not run)
```



---

transform\_etas\_tdist    *transform\_etas\_tdist*

---

### Description

Applies a t-distribution transformation to selected etas

Initial estimate for degrees of freedom is 80 with bounds (3, 100).

### Usage

```
transform_etas_tdist(model, list_of_etas = NULL)
```

### Arguments

**model**            (Model) Pharnpy model to apply t distribution transformation to.

**list\_of\_etas**    (array(str) or str (optional)) Name/names of etas to transform. If NULL, all etas will be transformed (default).

### Value

(Model) Pharnpy model object

### See Also

transform\_etas\_boxcox

transform\_etas\_john\_draper

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
model <- transform_etas_tdist(model, c("ETA_1"))  
model$statements$before_odes$full_expression("CL")  
  
## End(Not run)
```

---

translate\_nmtran\_time *translate\_nmtran\_time*

---

**Description**

Translate NM-TRAN TIME and DATE column into one TIME column

If dataset of model have special NM-TRAN TIME and DATE columns these will be translated into one single time column with time in hours.

Warnings Use this function with caution. For example reset events are currently not taken into account.

**Usage**

```
translate_nmtran_time(model)
```

**Arguments**

model (Model) PharmPy model object

**Value**

(Model) PharmPy model object

---

unconstrain\_parameters  
*unconstrain\_parameters*

---

**Description**

Remove all constraints from parameters

**Usage**

```
unconstrain_parameters(model, parameter_names)
```

**Arguments**

model (Model) PharmPy model

parameter\_names

(array(str)) Remove all constraints for the listed parameters

**Value**

(Model) PharmPy model object

**See Also**

set\_lower\_bounds : Set parameter lower bounds  
set\_upper\_bounds : Set parameter upper bounds  
unfix\_parameters : Unfix parameters

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model$parameters['PTVCL']  
model <- unconstrain_parameters(model, c('PTVCL'))  
model$parameters['PTVCL']  
  
## End(Not run)
```

---

undrop_columns	<i>undrop_columns</i>
----------------	-----------------------

---

**Description**

Undrop columns of model

**Usage**

```
undrop_columns(model, column_names)
```

**Arguments**

model (Model) Pharmpy model object  
column\_names (array(str) or str) List of column names or one column name to undrop

**Value**

(Model) Pharmpy model object

**See Also**

drop\_dropped\_columns : Drop all columns marked as drop  
drop\_columns : Drop or mark columns as dropped

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
model <- drop_columns(model, c('WGT', 'APGR'), mark=TRUE)  
model <- undrop_columns(model, 'WGT')  
  
## End(Not run)
```

---

unfix_parameters	<i>unfix_parameters</i>
------------------	-------------------------

---

### Description

Unfix parameters

Unfix all listed parameters

### Usage

```
unfix_parameters(model, parameter_names)
```

### Arguments

model (Model) PharmPy model

parameter\_names (array(str) or str) one parameter name or a vector of parameter names

### Value

(Model) PharmPy model object

### See Also

unfix\_parameters\_to : Unfixing parameters and setting a new initial estimate in the same function

fix\_parameters : Fix parameters

fix\_or\_unfix\_parameters : Fix or unfix parameters (given boolean)

fix\_parameters\_to : Fixing and setting parameter initial estimates in the same function

unconstrain\_parameters : Remove all constraints of parameters

### Examples

```
## Not run:
model <- load_example_model("pheno")
model <- fix_parameters(model, c('PTVCL', 'PTVV', 'THETA_3'))
model$parameters$fix
model <- unfix_parameters(model, 'PTVCL')
model$parameters$fix

## End(Not run)
```

---

unfix\_parameters\_to    *unfix\_parameters\_to*

---

### Description

Unfix parameters to  
Unfix all listed parameters to specified value/values

### Usage

```
unfix_parameters_to(model, inits)
```

### Arguments

model	(Model) PharmPy model
inits	(list(str=numeric)) Inits for all parameters to unfix and change init

### Value

(Model) PharmPy model object

### See Also

fix\_parameters : Fix parameters  
fix\_or\_unfix\_parameters : Fix or unfix parameters (given boolean)  
unfix\_paramaters : Unfixing parameters  
fix\_paramaters\_to : Fixing parameters and setting a new initial estimate in the same function

### Examples

```
## Not run:  
model <- load_example_model("pheno")  
model <- fix_parameters(model, c('PTVCL', 'PTVV', 'THETA_3'))  
model$parameters$fix  
model <- unfix_parameters_to(model, {'PTVCL': 0.5})  
model$parameters$fix  
model$parameters['PTVCL']  
  
## End(Not run)
```

---

```
update_initial_individual_estimates
      update_initial_individual_estimates
```

---

**Description**

Update initial individual estimates for a model  
 Updates initial individual estimates for a model.

**Usage**

```
update_initial_individual_estimates(model, individual_estimates, force = TRUE)
```

**Arguments**

`model` (Model) Pharmpy model to update initial estimates  
`individual_estimates` (array) Individual estimates to use  
`force` (logical) Set to FALSE to only update if the model had initial individual estimates before

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
results <- load_example_model_fit_results("pheno")
ie <- results$individual_estimates
model <- update_initial_individual_estimates(model, ie)

## End(Not run)
```

---

```
update_inits      update_inits
```

---

**Description**

Update initial parameter estimate for a model  
 Updates initial estimates of population parameters for a model. If the new initial estimates are out of bounds or NaN this function will raise.

**Usage**

```
update_inits(model, parameter_estimates, move_est_close_to_bounds = FALSE)
```

**Arguments**

`model` (Model) Pharmpy model to update initial estimates  
`parameter_estimates` (array) Parameter estimates to update  
`move_est_close_to_bounds` (logical) Move estimates that are close to bounds. If correlation >0.99 the correlation will be set to 0.9, if variance is <0.001 the variance will be set to 0.01.

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
results <- load_example_model_fit_results("pheno")  
model$parameters$inits  
model <- update_inits(model, results$parameter_estimates)  
model$parameters$inits  
  
## End(Not run)
```

---

use\_thetas\_for\_error\_stdev  
*use\_thetas\_for\_error\_stdev*

---

**Description**

Use thetas to estimate standard deviation of error

**Usage**

```
use_thetas_for_error_stdev(model)
```

**Arguments**

`model` (Model) Pharmpy model

**Value**

(Model) Pharmpy model object

**See Also**

set\_weighted\_error\_model : Encode error model with one epsilon and weight

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- use_thetas_for_error_stdev(model)
model$statements$find_assignment("W")
model$statements$find_assignment("Y")

## End(Not run)
```

---

write\_csv

*write\_csv*

---

**Description**

Write dataset to a csv file and updates the datainfo path

**Usage**

```
write_csv(model, path = NULL, force = FALSE)
```

**Arguments**

model (Model) Model whose dataset to write to file  
path (str (optional)) Destination path. Default is to use original path with .csv suffix.  
force (logical) Overwrite file with same path. Default is FALSE.

**Value**

(Model) Updated model object

**Examples**

```
## Not run:
model <- load_example_model("pheno")
model <- write_csv(model, path="newdataset$csv")

## End(Not run)
```



---

write_model	<i>write_model</i>
-------------	--------------------

---

**Description**

Write model code to file

**Usage**

```
write_model(model, path = "", force = TRUE)
```

**Arguments**

model	(Model) Pharmpy model
path	(str) Destination path
force	(logical) Force overwrite, default is TRUE

**Value**

(Model) Pharmpy model object

**Examples**

```
## Not run:  
model <- load_example_model("pheno")  
write_model(model)  
  
## End(Not run)
```

---

write_results	<i>write_results</i>
---------------	----------------------

---

**Description**

Write results object to json (or csv) file

Note that the csv-file cannot be read into a results object again.

**Usage**

```
write_results(results, path, lzma = FALSE, csv = FALSE)
```

**Arguments**

<code>results</code>	(Results) PharmPy results object
<code>path</code>	(str) Path to results file
<code>lzma</code>	(logical) TRUE for lzma compression. Not applicable to csv file
<code>csv</code>	(logical) Save as csv file

# Index

add\_allometry, 6  
add\_covariance\_step, 7  
add\_covariate\_effect, 8  
add\_estimation\_step, 11  
add\_iiv, 12  
add\_individual\_parameter, 14  
add\_iov, 14  
add\_lag\_time, 15  
add\_metabolite, 16  
add\_peripheral\_compartment, 17  
add\_pk\_iiv, 18  
add\_population\_parameter, 19  
add\_time\_after\_dose, 20  
append\_estimation\_step\_options, 20  
  
bump\_model\_number, 21  
  
calculate\_aic, 22  
calculate\_bic, 22  
calculate\_corr\_from\_cov, 23  
calculate\_corr\_from\_prec, 24  
calculate\_cov\_from\_corrse, 25  
calculate\_cov\_from\_prec, 26  
calculate\_epsilon\_gradient\_expression, 27  
calculate\_eta\_gradient\_expression, 28  
calculate\_eta\_shrinkage, 29  
calculate\_individual\_parameter\_statistics, 30  
calculate\_individual\_shrinkage, 31  
calculate\_parameters\_from\_ucp, 32  
calculate\_pk\_parameters\_statistics, 33  
calculate\_prec\_from\_corrse, 34  
calculate\_prec\_from\_cov, 35  
calculate\_se\_from\_cov, 36  
calculate\_se\_from\_prec, 37  
calculate\_ucp\_scale, 38  
check\_dataset, 38  
check\_high\_correlations, 39  
check\_parameters\_near\_bounds, 40  
  
check\_pharmpy, 41  
cleanup\_model, 41  
convert\_model, 42  
create\_basic\_pk\_model, 43  
create\_config\_template, 44  
create\_joint\_distribution, 44  
create\_report, 45  
create\_results, 45  
create\_rng, 46  
create\_symbol, 47  
  
deidentify\_data, 47  
display\_odes, 48  
drop\_columns, 49  
drop\_dropped\_columns, 50  
  
evaluate\_epsilon\_gradient, 50  
evaluate\_eta\_gradient, 51  
evaluate\_expression, 52  
evaluate\_individual\_prediction, 53  
evaluate\_population\_prediction, 54  
evaluate\_weighted\_residuals, 55  
expand\_additional\_doses, 56  
  
find\_clearance\_parameters, 56  
find\_volume\_parameters, 57  
fit, 58  
fix\_or\_unfix\_parameters, 58  
fix\_parameters, 59  
fix\_parameters\_to, 60  
  
get\_baselines, 61  
get\_bioavailability, 62  
get\_cmt, 62  
get\_concentration\_parameters\_from\_data, 63  
get\_config\_path, 63  
get\_covariate\_baselines, 64  
get\_doseid, 65  
get\_doses, 65

get\_dv\_symbol, 66  
get\_evid, 67  
get\_ids, 67  
get\_individual\_parameters, 68  
get\_individual\_prediction\_expression,  
69  
get\_initial\_conditions, 69  
get\_lag\_times, 70  
get\_mdv, 71  
get\_model\_code, 71  
get\_model\_covariates, 72  
get\_number\_of\_individuals, 72  
get\_number\_of\_observations, 73  
get\_number\_of\_observations\_per\_individual,  
74  
get\_observation\_expression, 76  
get\_observations, 75  
get\_omegas, 76  
get\_pk\_parameters, 77  
get\_population\_prediction\_expression,  
78  
get\_rv\_parameters, 79  
get\_sigmas, 79  
get\_thetas, 80  
get\_unit\_of, 81  
get\_zero\_order\_inputs, 82  
greekify\_model, 82  
  
has\_additive\_error\_model, 83  
has\_combined\_error\_model, 84  
has\_covariate\_effect, 85  
has\_first\_order\_elimination, 85  
has\_linear\_odes, 86  
has\_linear\_odes\_with\_real\_eigenvalues,  
87  
has\_michaelis\_menten\_elimination, 87  
has\_mixed\_mm\_fo\_elimination, 88  
has\_odes, 89  
has\_proportional\_error\_model, 89  
has\_random\_effect, 90  
has\_weighted\_error\_model, 91  
has\_zero\_order\_absorption, 92  
has\_zero\_order\_elimination, 92  
  
install\_pharmpy, 93  
install\_pharmpy\_devel, 94  
is\_linearized, 94  
is\_real, 95  
  
list\_time\_varying\_covariates, 96  
load\_example\_model, 96  
load\_example\_modelfit\_results, 97  
  
make\_declarative, 98  
mu\_reference\_model, 98  
  
omit\_data, 99  
  
plot\_individual\_predictions, 100  
plot\_iofv\_vs\_iofv, 100  
plot\_transformed\_eta\_distributions,  
101  
predict\_influential\_individuals, 101  
predict\_influential\_outliers, 102  
predict\_outliers, 103  
print\_fit\_summary, 104  
print\_model\_code, 104  
print\_model\_symbols, 105  
print\_pharmpy\_version, 105  
  
rank\_models, 106  
read\_dataset\_from\_datainfo, 107  
read\_model, 107  
read\_model\_from\_string, 108  
read\_modelfit\_results, 108  
read\_results, 109  
remove\_covariance\_step, 110  
remove\_covariate\_effect, 111  
remove\_error\_model, 111  
remove\_estimation\_step, 112  
remove\_iiv, 113  
remove\_iov, 114  
remove\_lag\_time, 115  
remove\_loq\_data, 115  
remove\_peripheral\_compartment, 116  
remove\_unused\_parameters\_and\_rvs, 117  
rename\_symbols, 117  
resample\_data, 118  
reset\_index, 119  
reset\_indices\_results, 119  
resume\_tool, 120  
retrieve\_final\_model, 120  
retrieve\_models, 121  
run\_allometry, 122  
run\_amd, 123  
run\_bootstrap, 124  
run\_covsearch, 125  
run\_estmethod, 126

run\_iivsearch, 127  
run\_iovsearch, 128  
run\_modelfit, 129  
run\_modelsearch, 130  
run\_ruvsearch, 131  
run\_tool, 132

sample\_individual\_estimates, 132  
sample\_parameters\_from\_covariance\_matrix, 133  
sample\_parameters\_uniformly, 135  
set\_additive\_error\_model, 136  
set\_bolus\_absorption, 137  
set\_combined\_error\_model, 138  
set\_covariates, 139  
set\_dtbs\_error\_model, 139  
set\_dvid, 140  
set\_estimation\_step, 140  
set\_evaluation\_step, 141  
set\_first\_order\_absorption, 142  
set\_first\_order\_elimination, 143  
set\_iiv\_on\_ruv, 144  
set\_initial\_condition, 145  
set\_initial\_estimates, 145  
set\_lower\_bounds, 146  
set\_michaelis\_menten\_elimination, 147  
set\_mixed\_mm\_fo\_elimination, 148  
set\_name, 149  
set\_ode\_solver, 149  
set\_peripheral\_compartments, 150  
set\_power\_on\_ruv, 151  
set\_proportional\_error\_model, 152  
set\_seq\_zo\_fo\_absorption, 153  
set\_time\_varying\_error\_model, 154  
set\_tmdd, 155  
set\_transit\_compartments, 155  
set\_upper\_bounds, 156  
set\_weighted\_error\_model, 157  
set\_zero\_order\_absorption, 158  
set\_zero\_order\_elimination, 158  
set\_zero\_order\_input, 159  
simplify\_expression, 160  
solve\_ode\_system, 161  
split\_joint\_distribution, 161  
str, 68, 77, 79  
summarize\_errors, 162  
summarize\_individuals, 163  
summarize\_individuals\_count\_table, 164  
summarize\_modelfit\_results, 165  
transform\_blk, 165  
transform\_etas\_boxcox, 167  
transform\_etas\_john\_draper, 168  
transform\_etas\_tdist, 169  
translate\_nmtran\_time, 170

unconstrain\_parameters, 170  
undrop\_columns, 171  
unfix\_parameters, 172  
unfix\_parameters\_to, 173  
update\_initial\_individual\_estimates, 174  
update\_inits, 174  
use\_thetas\_for\_error\_stddev, 175

write\_csv, 176  
write\_model, 177  
write\_results, 177