

# Package ‘atlantistools’

August 16, 2017

**Type** Package

**Title** Process and Visualise Output from Atlantis Models

**Version** 0.4.3

**Description** Atlantis is an end-to-end marine ecosystem modelling framework. It was originally developed in Australia by E.A. Fulton, A.D.M. Smith and D.C. Smith (2007) and has since been adopted in many marine ecosystems around the world (<<http://atlantis.cmar.csiro.au>>). The output of an Atlantis simulation is stored in various file formats like .netcdf and .txt and different output structures are used for the output variables like e.g. productivity or biomass. This package is used to convert the different output types to a unified format according to the “tidy-data” approach by H. Wickham (2014) <DOI:10.18637/jss.v059.i10>. Additionally, ecological metrics like for example spatial overlap of predator and prey or consumption can be calculated and visualised with this package. Due to the unified data structure it is very easy to share model output with each other and perform model comparisons.

**URL** <https://github.com/alketh/atlantistools>

**BugReports** <https://github.com/alketh/atlantistools/issues>

**Depends** R (>= 3.2.3)

**License** GPL-3

**LazyData** true

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**VignetteBuilder** knitr

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agg\_data

*Aggregate data using dplyr functionality.***Description**

This function is a 'wrapper' for the `group_by` and `summarize` procedure used in `dplyr` to aggregate dataframes.

**Usage**

```
agg_data(data, col = "atoutput", groups, out = "atoutput", fun)
agg_perc(data, col = "atoutput", groups, out = "atoutput")
group_data(data, groups)
```

**Arguments**

data	Dataframe the aggregation is applied to.
col	Column of the dataframe the summarise function is applied. Default is atoutput.
groups	Vector of character strings giving the grouping variables.
out	Character string specifying the name of the output column. Default is atoutput.
fun	Aggregation function to apply.

**Value**

grouped datarame with the aggregated data.

**Examples**

```
agg_ref_nums <- agg_data(data = ref_nums, groups = c("species", "agecl"), fun = mean)
```

---

calculate\_biomass\_spatial

*Calculate spatially explicit biomass (in [t]) for each group and age-class per timestep.*

---

**Description**

Calculate spatially explicit biomass time series for each group and ageclass within our model. Data is read in from 'output[...].nc'. Biomass for age-based groups is calculated as (StructN [mgN] + ResN [mgN]) \* Numbers [individuals]. Biomass for non age-based groups is calculated as N [mgN] \* volume [m<sup>3</sup>] (sediment-dz [m] / volume [m<sup>3</sup>] for epibenthic groups). mgN is converted to t based on the settings in the biol.prm file. Simulation time steps are converted to time in years based on output timesteps given in run.prm.

**Usage**

```
calculate_biomass_spatial(nums, sn, rn, n, vol_dz, bio_conv, bps)
```

**Arguments**

nums	Dataframe with information about numbers for age-based groups. Should be generated with <code>load_nc</code> .
sn	Dataframe with information about structural nitrogen for age-based groups. Should be generated with <code>load_nc</code> .
rn	Dataframe with information about reserve nitrogen for age-based groups. Should be generated with <code>load_nc</code> .
n	Dataframe with information about nitrogen for non-age-based groups. Should be generated with <code>load_nc</code> .
vol_dz	Dataframe with information about volume and layer height per polygon. Should be generated with <code>load_nc_physics</code> .
bio_conv	Numeric value to transform weight in mg N to tonnes. Should be generated with <code>get_conv_mgnbiot</code> .
bps	Vector of character strings giving the complete list of epibenthic functional groups (Only present in the sediment layer). The names have to match the column 'Name' in the 'functionalGroups.csv' file. Should be generated with <code>load_bps</code> .

**Value**

Dataframe with columns 'species', 'agecl', 'polygon', 'layer', 'time'. Biomass in [t] is stored in column 'atoutput'.

**Examples**

```
# 1. Using built in datasets.
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
prm_biol <- file.path(d, "VMPA_setas_biol_fishing_Trunk.prm")
fgs <- file.path(d, "SETasGroupsDem_NoCep.csv")
init <- file.path(d, "INIT_VMPA_Jan2015.nc")
bps <- load_bps(fgs = fgs, init = init)

bio_conv <- get_conv_mgnbiot(prm_biol = prm_biol)

df <- calculate_biomass_spatial(nums = ref_nums, sn = ref_structn, rn = ref_resn, n = ref_n,
                               vol_dz = ref_vol_dz, bio_conv = bio_conv, bps = bps)

# 2. Read in dataframes from existing Atlantis simulation.
bboxes <- get_boundary(boxinfo = load_box(file.path(d, "VMPA_setas.bgm")))
nc_gen <- file.path(d, "outputSETAS.nc")
prm_run <- file.path(d, "VMPA_setas_run_fishing_F_Trunk.prm")

groups_age <- c("Planktiv_S_Fish", "Pisciv_S_Fish")
groups_rest <- c("Cephalopod", "Megazoobenthos", "Diatom", "Lab_Det", "Ref_Det")

nums <- load_nc(nc = nc_gen, bps = bps, fgs = fgs,
               select_groups = groups_age, select_variable = "Nums",
               prm_run = prm_run, bboxes = bboxes)
sn <- load_nc(nc = nc_gen, bps = bps, fgs = fgs,
```

```

      select_groups = groups_age, select_variable = "StructN",
      prm_run = prm_run, bboxes = bboxes)
rn <- load_nc(nc = nc_gen, bps = bps, fgs = fgs,
      select_groups = groups_age, select_variable = "ResN",
      prm_run = prm_run, bboxes = bboxes)
n <- load_nc(nc = nc_gen, bps = bps, fgs = fgs,
      select_groups = groups_rest, select_variable = "N",
      prm_run = prm_run, bboxes = bboxes)
vol <- load_nc_physics(nc = nc_gen, select_physics = c("volume", "dz"),
      prm_run = prm_run, bboxes = bboxes, aggregate_layers = FALSE)

df <- calculate_biomass_spatial(nums = nums, sn = sn, rn = rn, n = n, vol_dz = vol,
      bio_conv = bio_conv, bps = bps)

# 3. Read in dataframes from existing Atlantis simulation with Map().
vars <- list("Nums", "StructN", "ResN", "N")
grps <- list(groups_age, groups_age, groups_age, groups_rest)
dfs <- Map(load_nc, select_variable = vars, select_groups = grps,
      MoreArgs = list(nc = nc_gen, bps = bps, fgs = fgs,
      prm_run = prm_run, bboxes = bboxes))

df <- calculate_biomass_spatial(nums = dfs[[1]], sn = dfs[[2]], rn = dfs[[3]], n = dfs[[4]],
      vol_dz = vol, bio_conv = bio_conv, bps = bps)

```

---

calculate\_consumed\_biomass

*Calculate the consumed biomass in [t] of prey j by predator i.*

---

### Description

Consumption data is extracted from output[...]`PROD.nc`. Age based groups are stored as "Eat\_" non age based groups as "Grazing\_". Units are  $\text{mg N m}^{-3} \text{d}^{-1}$ . Factors are species, time, box and agecl (if present). We will refer to species as pred from here on to indicate the predator perspective. Diet contribution data is extracted from `DietCheck.txt`. Currently this only works for models based on the trunk code. Units are The consumed biomass is calculated as follows: - Calculate consumed biomass as Eat (or Grazing) \* boxvolume per time, pred, agecl, box. - Convert to biomass in [t]. - Combine with diet contributions and calculate consumed biomass of prey species.

### Usage

```
calculate_consumed_biomass(eat, grazing, dm, vol, bio_conv)
```

### Arguments

eat	Dataframe with information about consumption for age-based groups. Should be generated with <code>load_nc</code> .
grazing	Dataframe with information about consumption for non-age-based groups. Should be generated with <code>load_nc</code> .







---

change_avail	<i>Change the availability matrix to simplify automated ATLANTIS calibrations.</i>
--------------	--

---

### Description

Change the availability of predator XXX on specific preygroups.

### Usage

```
change_avail(dietmatrix, pred = NULL, pred_stanza = NULL, prey = NULL,
             roc, relative = TRUE, consecutive = FALSE)
```

### Arguments

dietmatrix	Dataframe in 'long' format containing information about availabilities with columns 'pred', 'prey', 'pred_stanza', 'prey_stanza', 'code', 'prey_id' and 'avail'. The dataframe should be generated with <code>load_dietmatrix()</code> .
pred	Character vector of predator Acronyms (see <code>get_acronyms()</code> ). Selecting NULL as pred results in all predators being selected. This can be helpful if you want to increase the feeding pressure on a specific prey item by all groups. Default is NULL.
pred_stanza	Integer vector indicating if the predator is juvenile (= 1) or adult (= 2). pred and pred_stanza need to be of the same length. In rare instances, e.g. pred_stanza being NULL or one single integer either all pred_stanzas are selected or the single pred_stanza is applied to all predators. Default is NULL.
prey	List of character vectors of prey Acronyms (see <code>get_acronyms()</code> ). pred and prey need to be of the same length. Selecting NULL as prey results in all prey groups being selected. This can be helpful if you want to increase the available prey for a specific predator overall. Default is NULL.
roc	Vector of multiplication factors which shall be applied to the old set of parameters. Please supply one value per selected group. In case relative is FALSE the new absolute values can be passed as roc.
relative	Logical if TRUE values are changed relative to base values. If FALSE new values can be passed directly. Default is NULL.
consecutive	Boolean indicating if multiple calls to change_avail are performed one after another TRUE. Default is FALSE.

### Value

parameterfile \*.prm file with the new parameter values.

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
dm <- load_dietmatrix(prm_biol = file.path(d, "VMPA_setas_biol_fishing_Trunk.prm"),
                    fgs = file.path(d, "SETasGroupsDem_NoCep.csv"))

dm1 <- change_avail(dietmatrix = dm,
                  pred = "FPS",
                  pred_stanza = 1,
                  prey = "CEP",
                  roc = 0.1234,
                  relative = FALSE)
# Show only rows with availability of 0.1234
dm1[apply(apply(dm1[, 5:ncol(dm1)], MARGIN = 2, function(x) x == 0.1234), MARGIN = 1, any), ]

dm2 <- change_avail(dietmatrix = dm,
                  pred = c("FPS", "FVS"),
                  pred_stanza = c(1, 2),
                  prey = list(c("FPS", "FVS"), c("FPS", "FVS")),
                  roc = list(c(0.1111, 0.2222), c(0.3333, 0.4444)),
                  relative = FALSE)

# Show only rows with availability of 0.1111, 0.2222, 0.3333 or 0.4444
dm2[apply(apply(dm2[, 5:ncol(dm2)], MARGIN = 2,
              function(x) is.element(x, c(0.1111, 0.2222, 0.3333, 0.4444))), MARGIN = 1, any), ]
```

---

change\_prm

*Change biological parameterfile to simplify automated ATLANTIS calibrations.*

---

**Description**

This function is used to help automate the calibration routine for ATLANTIS models.

**Usage**

```
change_prm(prm_biol, select_acronyms, roc, parameter, relative = TRUE,
          save_to_disc = TRUE, version_flag = 2)
```

**Arguments**

prm_biol	Character string giving the connection to the biological parameterfile. The file-name usually contains biol_fishing and does end in .prm.
select_acronyms	Character vector of functional groups which shall be read in. Names have to match the ones used in the *.prm file. Check column "Code" in "functional-Groups.csv" for clarification.
roc	Vector of multiplication factors which shall be applied to the old set of parameters. Please supply one value per selected group. In case relative is FALSE the new absolute values can be passed as roc.

parameter	Character value of the model parameter which shall be changed. Only one parameter can be selected per function call.
relative	Logical if TRUE values are changed relative to base values. If FALSE new values can be passed directly. Default is TRUE.
save_to_disc	Logical indicating if the resulting prm file should be overwritten (TRUE) or not (FALSE). Defaults to TRUE.
version_flag	The version of ATLANTIS model. 1 for bec_dev, 2 for trunk. default is 2..

### Value

parameterfile \*.prm file with the new parameter values.

### Examples

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
prm_biol <- file.path(d, "VMPA_setas_biol_fishing_Trunk.prm")

new_prm <- change_prm(prm_biol,
  select_acronyms = c("FPS", "FVS"),
  roc = c(2,3),
  parameter = "KWRR",
  save_to_disc = FALSE)
```

---

change_prm_cohort	<i>Change biological parameterfile for parameters which expect multiple values.</i>
-------------------	---

---

### Description

This function is used to help automate the calibration routine for ATLANTIS models.

### Usage

```
change_prm_cohort(prm_biol, select_acronyms, roc, parameter, relative = TRUE,
  save_to_disc = TRUE)
```

### Arguments

prm_biol	Character string giving the connection to the biological parameterfile. The file-name usually contains biol_fishing and does end in .prm.
select_acronyms	Character vector of functional groups which shall be read in. Names have to match the ones used in the *.prm file. Check column "Code" in "functional-Groups.csv" for clarification.

roc	Matrix of multiplication factors which shall be applied to the old set of parameters. Please supply one row per selected group. Each row should have as many entries as the parameter itself. E.g. if you want to change the clearance rate for two fish groups you need to supply a matrix with 2 rows and 10 columns. In case you use different cohort numbers for age-structured groups supply a list of multiplication factors. Each list entry should be group specific.
parameter	Character value of the model parameter which shall be changed. Only one parameter can be selected per function call.
relative	Logical if TRUE values are changed relative to base values. If FALSE new values can be passed directly. Default is TRUE.
save_to_disc	Logical indicating if the resulting prm file should be overwritten (TRUE) or not (FALSE). Defaults to TRUE.

### Value

parameterfile \*.prm file with the new parameter values.

### Examples

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")

new_prm <- change_prm_cohort(prm_biol = file.path(d, "VMPA_setas_biol_fishing_Trunk.prm"),
  select_acronyms = c("FPS", "FVS"),
  roc = matrix(rep(2, times = 20), nrow = 2, ncol = 10),
  parameter = "C",
  save_to_disc = FALSE)
# C_FPS is in line 640. Old values are 0.0002 0.3 0.6 0.6 0.6 0.6 0.5 0.5 0.4 and 0.4.
new_prm[640:641]
# C_FVS is in line 652. Old values are 40.0 40.0 40.0 120.0 150.0 250.0 250.0 300.0 300.0 and 300.0.
new_prm[652:653]

# Also works for lists as argument
new_prm <- change_prm_cohort(prm_biol = file.path(d, "VMPA_setas_biol_fishing_Trunk.prm"),
  select_acronyms = c("FPS", "FVS"),
  roc = list(rep(3, times = 10), rep(2, times = 10)),
  parameter = "C",
  save_to_disc = FALSE)
```

---

check\_df\_names

*Function to check the names of a dataframe.*

---

### Description

This function is used in most plotting routines to check if the correct dataframe is passed.

### Usage

```
check_df_names(data, expect, optional = NULL)
```

**Arguments**

data	Dataframe to check.
expect	Character vector giving the names of the expected columns
optional	Character vector giving the names of optional columns. Default is NULL.

**Examples**

```
## Not run:
check_df_names(preprocess$biomass_age, expect = c("time", "species", "atoutput", "ages"))

## End(Not run)
```

---

check_growth	<i>This function is used to check the individual growth per group over time.</i>
--------------	--

---

**Description**

This function is used to check the individual growth per group over time.

**Usage**

```
check_growth(data, yearly = FALSE)
```

**Arguments**

data	Dataframe with information about individual age based growth over time. This should be generated with preprocess(). You can test either structural or reserve weight.
yearly	Logical specifying if relative change in individual weight shall be calculated on a yearly basis (TRUE) or not (FALSE). Default is FALSE.

**Value**

Dataframe showing the output of the linear model fit (slope & F-statistic) per group and age.

**Examples**

```
check_growth(preprocess$structn_age)
check_growth(preprocess$resn_age)
check_growth(preprocess$resn_age, yearly = TRUE)
```

---

combine_ages	<i>Combine ageclasses to juvenile and adult stanza according to age at maturity.</i>
--------------	--

---

### Description

Combine ageclasses to juvenile and adult stanza according to age at maturity.

### Usage

```
combine_ages(data, grp_col, agemat, value_col = "atoutput")
```

### Arguments

data	Dataframe with ageclass specific information.
grp_col	Character string giving the name of the group column in data. E.g. 'species', 'pred', 'prey' etc.
agemat	First mature age class for age structured groups. This dataframe should be generated with <a href="#">prm_to_df</a> using "age_mat" as parameter.
value_col	Character string giving the name of the column to sum. Default is "atoutput".

### Value

Dataframe with ageclasses combined to stanzas.

### Examples

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
prm_biol <- file.path(d, "VMPA_setas_biol_fishing_Trunk.prm")
fgs <- file.path(d, "SETasGroupsDem_NoCep.csv")

agemat <- prm_to_df(prm_biol = prm_biol, fgs = fgs,
                  group = get_age_acronyms(fgs = fgs),
                  parameter = "age_mat")

combine_ages(ref_nums, grp_col = "species", agemat = agemat)
```

---

combine_groups	<i>Combine values from different groups if specific groups only have a low contribution to the overall value.</i>
----------------	---

---

### Description

Combine values from different groups if specific groups only have a low contribution to the overall value.

**Usage**

```
combine_groups(data, group_col, groups = names(data)[!is.element(names(data),
  c("atoutput", "time", group_col))], combine_thresh = 15)
```

**Arguments**

`data` Dataframe whose groups shall be combined.  
`group_col` Character string giving the name of the group column in 'data'.  
`groups` Vector of character strings giving the grouping variables.  
`combine_thresh` Integer indicating the number of groups to display. Default is 15.

**Value**

dataframe with groups combined to "Rest" if contribution is low.

**See Also**

Other combine functions: [combine\\_runs](#)

**Examples**

```
df <- combine_groups(ref_dm, group_col = "prey")
df <- combine_groups(ref_dm, group_col = "prey", combine_thresh = 2)
```

---

combine_runs	<i>This function is used to combine model output from different simulations!</i>
--------------	--

---

**Description**

This function is used to combine model output from different simulations!

**Usage**

```
combine_runs(outs, runs)
```

**Arguments**

`outs` List of preprocessed Atlantis simulations. Each entry is a list generated with 'preprocess()'.  
`runs` Vector of character strings giving the name of each simulation settings.

**Value**

Named list with the the same format as in 'preprocess()'. Each dataframe has an additional column run.

**See Also**

Other combine functions: [combine\\_groups](#)

**Examples**

```
outs <- list(preprocess, preprocess)
runs <- c("run1", "run2")
test <- combine_runs(outs, runs)
names(test[[1]])
head(test[[1]])
```

---

convert\_bgm

*Transform data from bgm-file to map dataframe.*

---

**Description**

Transform data from bgm-file to map dataframe.

**Usage**

```
convert_bgm(bgm)
```

**Arguments**

bgm                    Character string giving the connection to the atlantis bgm file. The filename ends in .bgm.

**See Also**

Other convert functions: [convert\\_factor](#), [convert\\_time](#)

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
bgm <- file.path(d, "VMPA_setas.bgm")

bgm <- convert_bgm(bgm)
head(bgm)
```



---

convert_factor	<i>Function to convert any column with information about functional groups to a factor whose levels use the LongName of the functional groups file.</i>
----------------	---

---

**Description**

This function is used to match the labels of the plots!

**Usage**

```
convert_factor(data_fgs, col)
```

**Arguments**

data_fgs	Dataframe with information about functionalGroups. Usually loaded with load_fgs().
col	Column of the dataframe which is converted to a factor.

**Value**

Column of a dataframe in factor format.

**See Also**

Other convert functions: [convert\\_bgm](#), [convert\\_time](#)

---

```
convert_relative_initial
```

*Calculate relative timeseries using the initial value as benchmark.*

---

**Description**

Calculate relative timeseries using the initial value as benchmark.

**Usage**

```
convert_relative_initial(data, col = "atoutput")
```

**Arguments**

data	Dataframe to apply the transformation to.
col	Character value giving the name of the column to transform. Default is "atoutput".

**Value**

Dataframe with transformed column 'col'.

**Examples**

```
df <- convert_relative_initial(ref_structn)
head(df[df$layer == 1, ], n = 15)
```

---

convert_time	<i>Convert timestep to actual time!</i>
--------------	---

---

**Description**

Convert timestep to actual time!

**Usage**

```
convert_time(prm_run, col)
```

**Arguments**

prm_run	Character string giving the connection of the run parameterfile. The filename usually contains run_fishing and ends in .prm".
col	Numeric vector. Usually a column in a dataframe with information about time. Either given as timesteps or days.

**Value**

Numeric vector with the time in years.

**See Also**

Other convert functions: [convert\\_bgm](#), [convert\\_factor](#)

---

custom_grid	<i>Utility functions used for various plotting routines within atlantistools.</i>
-------------	---

---

**Description**

Utility functions used for various plotting routines within atlantistools.

**Usage**

```
custom_grid(plot, grid_x, grid_y)
```

**Arguments**

plot	ggplot2 plot.
grid_x	character vector used in facet_grid in horizontal direction.
grid_y	character vector used in facet_grid in vertical direction.

**Value**

ggplot2 plot.

---

extract_prm	<i>Extract values for Atlantis parameters from the biological parameter file.</i>
-------------	---

---

**Description**

Extract values for Atlantis parameters from the biological parameter file.

**Usage**

```
extract_prm(prm_biol, variables)
```

```
extract_prm_cohort(prm_biol, variables)
```

**Arguments**

prm_biol	Character string giving the connection to the biological parameterfile. The file-name usually contains biol_fishing and does end in .prm.
variables	Character string giving the flag to search for. This should be a combination of the parameter name and the group-Code.

**Value**

numeric vector.

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
prm_biol <- file.path(d, "VMPA_setas_biol_fishing_Trunk.prm")

# You can pass a single variable
extract_prm(prm_biol, variables = "KWRR_FVS")

# Or multiple variables
extract_prm(prm_biol, variables = paste("KWRR", c("FVS", "FPS"), sep = "_"))

# Use extract_prm_cohort do extract data for age specific parameters.
# They are usually stored in the next line following the parameter tag.
extract_prm_cohort(prm_biol, variables = "C_FVS")
extract_prm_cohort(prm_biol, variables = paste("C", c("FVS", "FPS"), sep = "_"))
```

---

fishbase_data	<i>fishbase_data</i>
---------------	----------------------

---

**Description**

A table of all the the species found in Fihibase, including taxonomic classification.

**Usage**

fishbase\_data

**Format**

A data frame with 33104 rows and 12 variables:

SpecCode integer. Species code.

Genus character.

Species character.

SpeciesRefNo integer. Reference number.

FBname character. Fishbase name.

SubFamily character.

FamCode integer. Family Code.

GenCode integer. Genetic Code.

SubGenCode integer. Sub Genetic Code.

Family character.

Order character.

Class character.

**Source**

<http://www.fishbase.org/> rfishbase::fishbase

---

flip_layers	<i>Flip layers for visualization.</i>
-------------	---------------------------------------

---

### Description

Within Atlantis the water column id 0 is the water column closest to the sediment. In order to simplify graphical interpretation of vertical plots this order is reversed. The surface layer is 1 by default. The sediment layer id is equal to the number of total layers. Please note that this is only used for graphical display.

### Usage

```
flip_layers(data)
```

### Arguments

data	dataframe with columns polygon and layer. layer id is based on atlantis output (0 = layer closest to the sediment)
------	--

### Value

dataframe with flipped layerids. 1 = surface.

### Examples

```
data <- rbind(expand.grid(species = "sp1", polygon = 0, layer = 0:7),
             expand.grid(species = "sp1", polygon = 1, layer = 0:4),
             expand.grid(species = "sp1", polygon = 2, layer = 0:2),
             expand.grid(species = "sp1", polygon = 3, layer = c(0:3, 7)))
data$output <- runif(nrow(data), min = 0, max = 2)
flip_layers(data)
```

---

get_boundary	<i>Get boundary boxes from Atlantis box information.</i>
--------------	--

---

### Description

Use the output from [load\\_box](#) and obtain a vector specifying which boxes are along the boundary.

### Usage

```
get_boundary(boxinfo)
```

### Arguments

boxinfo	A list as returned from <a href="#">load_box</a> .
---------	--

**Value**

A vector specifying which boxes are on the boundary.

**Author(s)**

Kelli Faye Johnson

**See Also**

[load\\_box](#)

Other get functions: [get\\_colpal](#), [get\\_conv\\_mgnbiot](#), [get\\_groups](#)

Other get functions: [get\\_colpal](#), [get\\_conv\\_mgnbiot](#), [get\\_groups](#)

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
boxes <- load_box(bgm = file.path(d, "VMPA_setas.bgm"))
get_boundary(boxinfo = boxes)
```

---

get\_colpal

*Create discrete color palette used in plots.*

---

**Description**

The colors are derived from <http://colorbrewer2.org/> using the palette 12-class Paired. In addition 9 different gray tones were added in case more than 12 categories are needed (This happens quite often with feeding data).

**Usage**

```
get_colpal()
```

**Value**

Vector of colors in hexadecimal code.

**See Also**

Other get functions: [get\\_boundary](#), [get\\_conv\\_mgnbiot](#), [get\\_groups](#)

---

get_conv_mgnbiot	<i>Extract conversion factor used to transform data from nitrogen in mg to biomass in tonnes.</i>
------------------	---

---

**Description**

Extract conversion factor used to transform data from nitrogen in mg to biomass in tonnes.

**Usage**

```
get_conv_mgnbiot(prm_biol)
```

**Arguments**

prm_biol	Character string giving the connection to the biological parameterfile. The file-name usually contains biol_fishing and does end in .prm.
----------	---

**Value**

Conversion factor as numeric value.

**See Also**

Other get functions: [get\\_boundary](#), [get\\_colpal](#), [get\\_groups](#)

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
prm_biol <- file.path(d, "VMPA_setas_biol_fishing_Trunk.prm")

get_conv_mgnbiot(prm_biol)
```

---

get_diet_fishbase	<i>Extract reference for diet information from <a href="http://www.fishbase.se">http://www.fishbase.se</a></i>
-------------------	--

---

**Description**

This function extracts reference for diet information from <http://www.fishbase.se>

**Usage**

```
get_diet_fishbase(fish, mirror = "se")
```

**Arguments**

fish	Vector of fish species with genus and species information.
mirror	Character string defining the url mirror to use. Defaults to se. In case data extraction is slow use a different mirror. Try to avoid frequently used mirrors like uk or com.

**Value**

Dataframe with species, country, locality, linf and k.

**Examples**

```
## Not run:
# For some reason the examples break with appveyor.
fish <- c("Gadus morhua", "Merlangius merlangus", "Maurolicus muelleri")
diet <- get_diet_fishbase(fish)

fish <- c("Gadus morhua", "Merlangius merlangus", "Ammodytes marinus")
diet <- get_diet_fishbase(fish)

## End(Not run)
```

---

get_groups	<i>Collection of similar functions which get specific columns from the Atlantis functionalGroups.csv</i>
------------	--

---

**Description**

This collection of functions uses the dataframe of functional groups created with [load\\_fgs](#) and creates various character strings of group names or acronym names.

**Usage**

```
get_groups(fgs)

get_age_groups(fgs)

get_acronyms(fgs)

get_age_acronyms(fgs)

get_nonage_acronyms(fgs)

get_fish_acronyms(fgs)
```



**Arguments**

`fgs` Character string giving the connection to the functional groups file. The file-name usually contains Groups and does end in .csv.

**Details**

Currently, the following character strings can be created - `get_groups`: Extract column "Name" - `get_age_groups`: Extract column "Name". Selects groups with 10 ageclasses. - `get_acronym`: Extract column "Code" - `get_age_acronym`: Extract column "Code". Selects groups with 10 ageclasses. - `get_nonage_acronym`: Extracts columns "Code". Only groups with ageclasses different from 10 are selected. - `get_fish_acronyms`: Extract column "Code". Only groups with InvertType equal to "FISH" or "SHARK" are selected.

**Value**

Character string.

**See Also**

Other get functions: [get\\_boundary](#), [get\\_colpal](#), [get\\_conv\\_mgnbiot](#)

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
fgs <- file.path(d, "SETasGroupsDem_NoCep.csv")

get_age_groups(fgs)
get_nonage_acronyms(fgs)
```

---

`get_growth_fishbase` *Extract growth parameters from <http://www.fishbase.se>.*

---

**Description**

This function extracts values for  $L_{inf}$ ,  $k$  and  $t_0$  from <http://www.fishbase.se>

**Usage**

```
get_growth_fishbase(fish, mirror = "se")
```

**Arguments**

`fish` Vector of fish species with genus and species information.

`mirror` Character string defining the url mirror to use. Defaults to se. In case data extraction is slow use a different mirror. Try to avoid frequently used mirrors like uk or com.

**Details**

Before the actual extraction takes place fishbase IDs for every species are extracted using `get_ids_fishbase`. The IDs are needed to generate the urls later on.

**Value**

Dataframe with species, country, locality, linf and k.

**Examples**

```
## Not run:
# For some reason the examples break with appveyor.
fish <- c("Gadus morhua", "Merlangius merlangus")
df <- get_growth_fishbase(fish)
head(df)

df <- get_growth_fishbase(fish, mirror = "de")
head(df)

fish <- c("Sprattus sprattus")
df <- get_growth_fishbase(fish)
head(df)
# Only use for debugging purposes.
fish <- read.csv("Z:/my_data_alex/fish_species_names_from_ibts.csv", stringsAsFactors = FALSE)[, 1]
url <- get_growth_fishbase(fish)
url <- urls$ref_url

## End(Not run)
```

---

<code>get_ids_fishbase</code>	<i>Extract fishbase IDs using the package "rfishbase" to generate species specific fishbase URLs</i>
-------------------------------	--

---

**Description**

This function extracts fishbase IDs using the database provided by the `rfishbase` package.

**Usage**

```
get_ids_fishbase(fish)
```

**Arguments**

`fish`                      Vector of fish species with genus and species information.

**Details**

The function depends on the package "rfishbase" which creates a local copy of the fishbase database. The IDs are needed to generate URLs to scan `www.fishbase.se` for detailed informations about fish growth for example.

**Value**

named vector with species names and fishbase IDs.

**Examples**

```
fish <- c("Gadus morhua", "Merlangius merlangus", "Clupea harengus")
get_ids_fishbase(fish)
```

---

get\_maturity\_fishbase *Extract maturity parameters from <http://www.fishbase.se>.*

---

**Description**

This function extracts values for maturity from <http://www.fishbase.se>

**Usage**

```
get_maturity_fishbase(fish, mirror = "se")
```

**Arguments**

fish	Vector of fish species with genus and species information.
mirror	Character string defining the url mirror to use. Defaults to se. In case data extraction is slow use a different mirror. Try to avoid frequently used mirrors like uk or com.

**Details**

Before the actual extraction takes place fishbase IDs for every species are extracted using [get\\_ids\\_fishbase](#). The IDs are needed to generate the urls later on.

**Value**

Dataframe with species, country, locality, linf and k.

**Examples**

```
## Not run:
# For some reason the examples break with appveyor.
fish <- c("Gadus morhua", "Squalus acanthias")
df <- get_maturity_fishbase(fish)
head(df)

## End(Not run)
```

---

get_ref_fishbase	<i>Extract the bibliographic info from www.fishbase.org.</i>
------------------	--

---

**Description**

Extract bibliographic information for growth parameters (linf, k, t0) from www.fishbase.org

**Usage**

```
get_ref_fishbase(ref_id, mirror = "se")
```

**Arguments**

ref_id	vector of reference ids.
mirror	Character string defining the url mirror to use. Defaults to se. In case data extraction is slow use a different mirror. Try to avoid frequently used mirrors like uk or com.

**Value**

Dataframe

**Examples**

```
## Not run:
df <- get_growth_fishbase("Scyliorhinus canicula")

df$data_ref[df$data_ref == df$main_ref] <- NA
df <- tidyr::gather_(data = df,
                    key_col = "ref_type",
                    value_col = "ref_id",
                    gather_cols = c("main_ref", "data_ref"), na.rm = TRUE)
ref_id <- unique(df$ref_id)
get_ref_fishbase(ref_id)

## End(Not run)
```

---

load_box	<i>Load the box specification file for an Atlantis scenario</i>
----------	---

---

**Description**

Read in the .bgm file that specifies the the box coordinates for an Atlantis scenario.

**Usage**

```
load_box(bgm)
```

**Arguments**

`bgm` Character string giving the connection to the atlantis bgm file. The filename ends in `.bgm`.

**Value**

A list of information regarding boxes for an Atlantis scenario.

**Author(s)**

Kelli Faye Johnson

**See Also**

Other load functions: [load\\_bps](#), [load\\_dietcheck](#), [load\\_fgs](#), [load\\_init\\_age](#), [load\\_init](#), [load\\_nc\\_physics](#), [load\\_nc](#), [load\\_rec](#), [load\\_spec\\_mort](#), [load\\_txt](#)

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
bgm <- file.path(d, "VMPA_setas.bgm")
```

```
boxes <- load_box(bgm)
```

---

<code>load_bps</code>	<i>Extracts the names of the epibenthic biomasspools from the initial conditions file.</i>
-----------------------	--

---

**Description**

Use `fgs` data.frame as read in by [load\\_fgs](#) to get the biomass pool information.

**Usage**

```
load_bps(fgs, init)
```

**Arguments**

`fgs` Character string giving the connection to the functional groups file. The filename usually contains `Groups` and does end in `.csv`.

`init` Character string giving the connection of the initial conditions netcdf file. The filename usually contains `init` and ends in `.nc`.

**Value**

Character vector of epibenthic biomass pools.

**See Also**[load\\_fgs](#)

Other load functions: [load\\_box](#), [load\\_dietcheck](#), [load\\_fgs](#), [load\\_init\\_age](#), [load\\_init](#), [load\\_nc\\_physics](#), [load\\_nc](#), [load\\_rec](#), [load\\_spec\\_mort](#), [load\\_txt](#)

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")

fgs <- file.path(d, "SETasGroupsDem_NoCep.csv")
init <- file.path(d, "INIT_VMPA_Jan2015.nc")

load_bps(fgs, init)
```

---

load_dietcheck	<i>Read in the atlantis dietcheck.txt file and perform some basic data transformations.</i>
----------------	---

---

**Description**

Read in the atlantis dietcheck.txt file and perform some basic data transformations.

**Usage**

```
load_dietcheck(dietcheck, fgs, prm_run, convert_names = FALSE,
  report = FALSE, version_flag = 2)
```

**Arguments**

dietcheck	Character string giving the connection of the dietcheck file. The filename usually contains Dietcheck and ends in .txt".
fgs	Character string giving the connection to the functional groups file. The filename usually contains Groups and does end in .csv.
prm_run	Character string giving the connection of the run parameterfile. The filename usually contains run_fishing and ends in .prm".
convert_names	Logical indicating if group codes are transformed to LongNames (TRUE) or not (default = FALSE).
report	Logical indicating if incomplete DietCheck information shall be printed TRUE or not FALSE.
version_flag	The version of ATLANTIS model. 1 for bec_dev, 2 for trunk. default is 2..

**Value**

A data.frame in long format with the following column names: time, pred, habitat, prey and atoutput (i.e., variable).

**See Also**

Other load functions: [load\\_box](#), [load\\_bps](#), [load\\_fgs](#), [load\\_init\\_age](#), [load\\_init](#), [load\\_nc\\_physics](#), [load\\_nc](#), [load\\_rec](#), [load\\_spec\\_mort](#), [load\\_txt](#)

**Examples**

```
# Apply to bec-dev models.
d <- system.file("extdata", "setas-model-new-becdev", package = "atlantistools")
dietcheck <- file.path(d, "outputSETASDietCheck.txt")
fgs <- file.path(d, "SETasGroups.csv")
prm_run <- file.path(d, "VMPA_setas_run_fishing_F_New.prm")

diet <- load_dietcheck(dietcheck, fgs, prm_run, version_flag = 1)
head(diet, n = 10)

# Apply to trunk models.
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
dietcheck <- file.path(d, "outputSETASDietCheck.txt")
fgs <- file.path(d, "SETasGroupsDem_NoCep.csv")
prm_run <- file.path(d, "VMPA_setas_run_fishing_F_Trunk.prm")
diet <- load_dietcheck(dietcheck, fgs, prm_run)
head(diet, n = 10)
```

---

load\_dietmatrix

---

*Extract the dietmatrix from the biological parameterfile*


---

**Description**

Extracts the diet matrix as long dataframe from the biological parameter file of any ATLANTIS simulation.

**Usage**

```
load_dietmatrix(prm_biol, fgs, transform = TRUE, convert_names = FALSE,
  version_flag = 2)
```

```
write_diet(dietmatrix, prm_biol, save_to_disc = TRUE)
```

**Arguments**

prm_biol	Character string giving the connection to the biological parameterfile. The file-name usually contains biol_fishing and does end in .prm.
fgs	Character string giving the connection to the functional groups file. The file-name usually contains Groups and does end in .csv.
transform	Boolean indicating if the returned dataframe is displayed in "long" (transform = TRUE, default) or "wide" (transform = FALSE) format. You should use the "wide" format in case you aim to change your diet matrix entries.

convert_names	Logical indicating if group codes are transformed to LongNames (TRUE) or not (default = FALSE).
version_flag	The version of ATLANTIS model. 1 for bec_dev, 2 for trunk. default is 2..
dietmatrix	Dataframe of the ATLANTIS dietmatrix generated with load_dietmatrix using transform = FALSE.
save_to_disc	Logical indicating if the resulting prm file should be overwritten (TRUE) or not (FALSE). Defaults to TRUE.

**Value**

dataframe of the availability matrix in long format with columns pred, pred\_stanza (1 = juvenile, 2 = adult), prey\_stanza, prey, avail, code.

**Examples**

```
# Can be applied to trunk models.
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
prm_biol <- file.path(d, "VMPA_setas_biol_fishing_Trunk.prm")
fgs <- file.path(d, "SETasGroupsDem_NoCep.csv")

dm <- load_dietmatrix(prm_biol, fgs)
head(dm, n = 10)

# Use write_diet to update your existing parameterfile.
dietmatrix <- load_dietmatrix(prm_biol, fgs, transform = FALSE)

# Write is set to FALSE here for technical reasons. Make sure to set it to TRUE in case you
# want to update your file.
new_diet <- write_diet(dietmatrix, prm_biol, save_to_disc = FALSE)

# And to bec-dev models.
d <- system.file("extdata", "setas-model-new-becdev", package = "atlantistools")
prm_biol <- file.path(d, "VMPA_setas_biol_fishing_New.prm")
fgs <- file.path(d, "SETasGroups.csv")

dm <- load_dietmatrix(prm_biol, fgs, version_flag = 1)
head(dm, n = 10)
```

---

load\_fgs

*Load the functional group file*


---

**Description**

Read in the functional group file as dataframe.

**Usage**

```
load_fgs(fgs)
```



**Arguments**

`fgs` Character string giving the connection to the functional groups file. The file-name usually contains Groups and does end in .csv.

**Value**

A data.frame of functional group information.

**See Also**

Other load functions: [load\\_box](#), [load\\_bps](#), [load\\_dietcheck](#), [load\\_init\\_age](#), [load\\_init](#), [load\\_nc\\_physics](#), [load\\_nc](#), [load\\_rec](#), [load\\_spec\\_mort](#), [load\\_txt](#)

**Examples**

```
d <- system.file("extdata", "setas-model-new-becdev", package = "atlantistools")
file <- "SETasGroups.csv"
fgs <- load_fgs(file.path(d, file))
head(fgs)
```

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
file <- "SETasGroupsDem_NoCep.csv"
fgs <- load_fgs(file.path(d, file))
head(fgs)
```

---

load_init	<i>This function is used to read in data from the initial conditions file.</i>
-----------	--

---

**Description**

This function is used to read in data from the initial conditions file.

**Usage**

```
load_init(init, vars)
```

**Arguments**

`init` Character string giving the connection of the initial conditions netcdf file. The filename usually contains `init` and ends in .nc.

`vars` Vector of character strings giving the variables to extract from the netcdf file.

**Value**

A list of dataframes with columns `atoutput`, `polygon` and `layer` (if present).

**Author(s)**

Alexander Keth

**See Also**

Other load functions: [load\\_box](#), [load\\_bps](#), [load\\_dietcheck](#), [load\\_fgs](#), [load\\_init\\_age](#), [load\\_nc\\_physics](#), [load\\_nc](#), [load\\_rec](#), [load\\_spec\\_mort](#), [load\\_txt](#)

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
init <- file.path(d, "INIT_VMPA_Jan2015.nc")

load_init(init, vars = "Planktiv_S_Fish1_Nums")
load_init(init, vars = c("Planktiv_S_Fish2_ResN", "Planktiv_S_Fish3_ResN"))
load_init(init, vars = "Megazoobenthos_N")

## Not run:
dir <- "C:/Users/siebo/Documents/Atlantis/BalticAtlantis/run_files_73days_Nej"
init <- file.path(dir, "new_init_Baltic_05Dec2015_v2.nc")
vars <- "Sprat1_ResN"
load_init(init = init, vars = vars)

## End(Not run)
```

---

load_init_age	<i>This function loads weight at age data (in mgN) from the initial conditions file.</i>
---------------	--

---

**Description**

This function loads weight at age data (in mgN) from the initial conditions file.

**Usage**

```
load_init_age(init, fgs, select_variable, select_groups = NULL, bboxes)

load_init_nonage(init, fgs, select_variable = "N", select_groups = NULL,
  bboxes, bps)

load_init_stanza(init, fgs, select_variable = "N", select_groups = NULL,
  bboxes)

load_init_physics(init, select_variable, bboxes)

load_init_weight(init, fgs, bboxes)
```

**Arguments**

init	Character string giving the connection of the initial conditions netcdf file. The filename usually contains init and ends in .nc.
------	---

fgs	Character string giving the connection to the functional groups file. The file-name usually contains Groups and does end in .csv.
select_variable	Character value specifying which variable to load. For load_init_age this can be "Nums", "ResN", "StructN", For load_init_nonage please select "N" (default) For load_init_physics simply pass the names of the physical variables.
select_groups	Character vector of functional groups which shall be read in. Names have to match the ones used in the ncd file. Check column "Name" in "functional-Groups.csv" for clarification. Default is NULL resulting in all available groups.
bboxes	Integer vector giving the box-id of the boundary boxes. Can be created with get_boundary.
bps	Vector of character strings giving the complete list of epibenthic functional groups (Only present in the sediment layer). The names have to match the column 'Name' in the 'functionalGroups.csv' file. Can be created with load_bps.#'

### Value

A dataframes with columns atoutput, polygon, layer (if present), species (if present).

### Author(s)

Alexander Keth

### See Also

Other load functions: [load\\_box](#), [load\\_bps](#), [load\\_dietcheck](#), [load\\_fgs](#), [load\\_init](#), [load\\_nc\\_physics](#), [load\\_nc](#), [load\\_rec](#), [load\\_spec\\_mort](#), [load\\_txt](#)

### Examples

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
init <- file.path(d, "INIT_VMPA_Jan2015.nc")
fgs <- file.path(d, "SETasGroupsDem_NoCep.csv")
bboxes <- get_boundary(load_box(bgm = file.path(d, "VMPA_setas.bgm")))

bps <- load_bps(fgs = fgs, init = init)

# There are no values in the initial conditions file. Therefore atoutput is NA.
load_init_age(init = init, fgs = fgs, bboxes = bboxes,
              select_variable = "ResN",
              select_groups = "Planktiv_S_Fish")

load_init_age(init = init, fgs = fgs, bboxes = bboxes, select_variable = "ResN")
load_init_nonage(init = init, fgs = fgs, bboxes = bboxes, bps = bps,
                 select_groups = "Megazoobenthos")
load_init_nonage(init = init, fgs = fgs, bboxes = bboxes, bps = bps)
load_init_stanza(init = init, fgs = fgs, bboxes = bboxes)
load_init_weight(init = init, fgs = fgs, bboxes = bboxes)
```

---

load\_nc *Load Atlantis outputfiles (netcdf)*

---

### Description

This function loads Atlantis outputfiles (any netcdf file) and converts them to a data.frame.

### Usage

```
load_nc(nc, fgs, bps, select_groups, select_variable, prm_run, bboxes,
        check_acronyms = TRUE, warn_zeros = FALSE, report = TRUE)
```

### Arguments

nc	Character string giving the connection of the netcdf file to read in. The filename usually contains output and ends in .nc".
fgs	Character string giving the connection to the functional groups file. The filename usually contains Groups and does end in .csv.
bps	Vector of character strings giving the complete list of epibenthic functional groups (Only present in the sediment layer). The names have to match the column 'Name' in the 'functionalGroups.csv' file. Can be created with load_bps.#
select_groups	Character vector of functional groups which shall be read in. Names have to match the ones used in the netcdf file. Check column "Name" in "functional-Groups.csv" for clarification.
select_variable	Character value specifying which variable to load. Only one variable of the options available (i.e., c("N", "Nums", "ResN", "StructN", "Eat", "Growth", "Prodn", "Grazing") can be loaded at a time.
prm_run	Character string giving the connection of the run parameterfile. The filename usually contains run_fishing and ends in .prm".
bboxes	Integer vector giving the box-id of the boundary boxes. Can be created with get_boundary.
check_acronyms	Logical testing if functional-groups in select_groups are inactive in the current model run. They will be omitted in the output.
warn_zeros	Logical indicating if check for actual zeros in the data shall be printed or not. Default is FALSE.
report	Logical indicating if progress bars shall be printed (TRUE) or not (FALSE). Default is TRUE.

### Value

A data.frame in long format with the following column names: species, timestep, polygon, agecl, and atoutput (i.e., variable).

**Author(s)**

Alexander Keth

**See Also**

Other load functions: [load\\_box](#), [load\\_bps](#), [load\\_dietcheck](#), [load\\_fgs](#), [load\\_init\\_age](#), [load\\_init](#), [load\\_nc\\_physics](#), [load\\_rec](#), [load\\_spec\\_mort](#), [load\\_txt](#)

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")

nc <- file.path(d, "outputSETAS.nc")
fgs <- file.path(d, "SETasGroupsDem_NoCep.csv")
bps <- load_bps(init = file.path(d, "INIT_VMPA_Jan2015.nc"), fgs = fgs)
bboxes <- get_boundary(boxinfo = load_box(bgm = file.path(d, "VMPA_setas.bgm")))
prm_run <- file.path(d, "VMPA_setas_run_fishing_F_Trunk.prm")

test <- load_nc(nc = nc, bps = bps, fgs = fgs, prm_run = prm_run, bboxes = bboxes,
  select_groups = c("Planktiv_S_Fish", "Cephalopod", "Diatom"),
  select_variable = "ResN")

test <- load_nc(nc = nc, bps = bps, fgs = fgs, prm_run = prm_run, bboxes = bboxes,
  select_groups = c("Planktiv_S_Fish", "Cephalopod", "Diatom"),
  select_variable = "Nums")
```

load\_nc\_physics

*Load Atlantis outputfiles (netcdf)***Description**

This function loads Atlantis outputfiles (netcdf) and converts them to a dataframe.

**Usage**

```
load_nc_physics(nc, select_physics, prm_run, bboxes, aggregate_layers = FALSE,
  warn_zeros = FALSE)
```

**Arguments**

nc	Character string giving the connection of the netcdf file to read in. The filename usually contains output and ends in .nc".
select_physics	Character vector of physical variables which shall be read in. Names have to match the ones used in the netcdf file.
prm_run	Character string giving the connection of the run parameterfile. The filename usually contains run_fishing and ends in .prm".
bboxes	Integer vector giving the box-id of the boundary boxes. Can be created with <code>get_boundary</code> .

aggregate_layers	Logical indicating if values for layers should be aggregated (TRUE) or not (FALSE). Default is FALSE.
warn_zeros	Logical indicating if check for actual zeros in the data shall be printed or not. Default is FALSE.

### Details

This functions converts the ATLANTIS output to a dataframe which can be processed in R.

### Value

A data.frame in long format with the following column names: variable, time, polygon, layer, and atoutput (i.e., variable).

### See Also

Other load functions: [load\\_box](#), [load\\_bps](#), [load\\_dietcheck](#), [load\\_fgs](#), [load\\_init\\_age](#), [load\\_init](#), [load\\_nc](#), [load\\_rec](#), [load\\_spec\\_mort](#), [load\\_txt](#)

### Examples

```
d <- system.file("extdata", "setas-model-new-becdev", package = "atlantistools")
nc <- file.path(d, "outputSETAS.nc")
prm_run <- file.path(d, "VMPA_setas_run_fishing_F_New.prm")
bboxes <- get_boundary(boxinfo = load_box(file.path(d, bgm = "VMPA_setas.bgm")))
select_physics <- c("salt", "NO3", "volume")

test <- load_nc_physics(nc, select_physics, prm_run, bboxes)
str(test)

d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
nc <- file.path(d, "outputSETAS.nc")
prm_run <- file.path(d, "VMPA_setas_run_fishing_F_Trunk.prm")
bboxes <- get_boundary(boxinfo = load_box(file.path(d, bgm = "VMPA_setas.bgm")))

test <- load_nc_physics(nc, select_physics, prm_run, bboxes)
str(test)

test <- load_nc_physics(nc, select_physics = "nominal_dz", prm_run, bboxes)
```

---

load\_rec

*Load information for SSB and Recruits from an Atlantis model run.*

---

### Description

Load information for SSB and Recruits from an Atlantis model run.

**Usage**

```
load_rec(yoy, ssb, prm_biol)
```

**Arguments**

**yoy** Character string giving the connection of the YOY file. The filename usually contains outputYOY and ends in .txt".

**ssb** Character string giving the connection of the YOY file. The filename usually contains outputSSB and ends in .txt".

**prm\_biol** Character string giving the connection to the biological parameterfile. The filename usually contains biol\_fishing and does end in .prm.

**Value**

Dataframe with information about ssb in tonnes and recruits in thousands.

**See Also**

Other load functions: [load\\_box](#), [load\\_bps](#), [load\\_dietcheck](#), [load\\_fgs](#), [load\\_init\\_age](#), [load\\_init](#), [load\\_nc\\_physics](#), [load\\_nc](#), [load\\_spec\\_mort](#), [load\\_txt](#)

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
yoy <- file.path(d, "outputSETASYOY.txt")
ssb <- file.path(d, "outputSETASSB.txt")
prm_biol <- file.path(d, "VMPA_setas_biol_fishing_Trunk.prm")

load_rec(yoy, ssb, prm_biol)
```

---

load\_spec\_mort

*Load mortality information from outputSpecificPredMort.txt*

---

**Description**

Load mortality information from outputSpecificPredMort.txt

**Usage**

```
load_spec_mort(specmort, prm_run, fgs, convert_names = FALSE,
  version_flag = 2)
```

**Arguments**

specmort	Character string giving the connection of the specific mortality file. The file-name usually contains SpecificPredMort and ends in .txt".
prm_run	Character string giving the connection of the run parameterfile. The filename usually contains run_fishing and ends in .prm".
fgs	Character string giving the connection to the functional groups file. The file-name usually contains Groups and does end in .csv.
convert_names	Logical indicating if group codes are transformed to LongNames (TRUE) or not (default = FALSE).
version_flag	The version of ATLANTIS model. 1 for bec_dev, 2 for trunk. default is 2..

**Value**

Dataframe with information about ssb in tonnes and recruits in thousands.

**See Also**

Other load functions: [load\\_box](#), [load\\_bps](#), [load\\_dietcheck](#), [load\\_fgs](#), [load\\_init\\_age](#), [load\\_init](#), [load\\_nc\\_physics](#), [load\\_nc](#), [load\\_rec](#), [load\\_txt](#)

**Examples**

```
d <- system.file("extdata", "setas-model-new-becdev", package = "atlantistools")
specmort <- file.path(d, "outputSETASspecificPredMort.txt")
prm_run <- file.path(d, "VMPA_setas_run_fishing_F_New.prm")
fgs <- file.path(d, "SETasGroups.csv")

df <- load_spec_mort(specmort, prm_run, fgs, version_flag = 1)
head(df)

d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
specmort <- file.path(d, "outputSETASspecificPredMort.txt")
prm_run <- file.path(d, "VMPA_setas_run_fishing_F_Trunk.prm")
fgs <- file.path(d, "SETasGroupsDem_NoCep.csv")

df <- load_spec_mort(specmort, prm_run, fgs)
head(df)
```

---

load\_txt

---

*Function to load various txt files from Atlantis simulations*


---

**Description**

Function to load various txt files from Atlantis simulations

**Usage**

```
load_txt(file, id_col = "Time")
```



**Arguments**

file	Character string giving the connection of the output file. The filename usually contains output and ends in .txt".
id_col	Character strings giving the names of the columns which are not variables. Data from all other columns will be gathered with tidyr. Default is "Time".

**Value**

Dataframe in tidy format!

**See Also**

Other load functions: [load\\_box](#), [load\\_bps](#), [load\\_dietcheck](#), [load\\_fgs](#), [load\\_init\\_age](#), [load\\_init](#), [load\\_nc\\_physics](#), [load\\_nc](#), [load\\_rec](#), [load\\_spec\\_mort](#)

**Examples**

```
d <- system.file("extdata", "setas-model-new-becdev", package = "atlantistools")
file <- file.path(d, "outputSETASSB.txt")
load_txt(file)

file <- file.path(d, "outputSETASYOY.txt")
load_txt(file)
```

---

plot_add_box	<i>Low level plotting function to add sudo confidence range to calibration plots.</i>
--------------	---

---

**Description**

Low level plotting function to add sudo confidence range to calibration plots.

**Usage**

```
plot_add_box(plot, range = c(0.5, 0.2))
```

**Arguments**

plot	ggplot2 object.
range	max and min relative change of data. Default is c(0.5, 0.2).

**Value**

ggplot2 plot.

**See Also**

Other low-level-plot functions: [plot\\_add\\_range](#)

**Examples**

```
# Make sure to use a relative timeseries generated with \code{\link{convert_relative_initial}}.
df <- convert_relative_initial(preprocess$structn_age)

# Create the plot with \code{\link{plot_line}}.
plot <- plot_line(df, col = "agec1")

# Add lower and upper range.
plot_add_box(plot)

# You can set the upper and lower range as you like!
plot_add_box(plot, range = c(0.8, 0.4))
```

---

plot\_add\_polygon\_overview

*Add spatial representation of polygon layout to a ggplot2 object.*

---

**Description**

Add spatial representation of polygon layout to a ggplot2 object.

**Usage**

```
plot_add_polygon_overview(plot, bgm_as_df, polygon_overview = 0.2)
```

**Arguments**

plot	ggplot2 object. Can be a ggplot grob.
bgm_as_df	*.bgm file converted to a dataframe. Please use <a href="#">convert_bgm</a> to convert your bgm-file to a dataframe with columns 'lat', 'long', 'inside_lat', 'inside_long' and 'polygon'.
polygon_overview	numeric value between 0 and 1 indicating the size used to plot the polygon overview in the upper right corner of the plot. Default is 0.2.

**Value**

ggplot grob

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
bgm_as_df <- convert_bgm(bgm = file.path(d, "VMPA_setas.bgm"))

p <- plot_line(preprocess$physics, wrap = NULL)
p <- custom_grid(p, grid_x = "polygon", grid_y = "variable")

grob <- plot_add_polygon_overview(p, bgm_as_df)
gridExtra::grid.arrange(grob)
```

---

plot_add_range	<i>Low level plotting function to add range of observed values to time series plots.</i>
----------------	--

---

## Description

This function can be used to add the range of observed data to a timeseries plot generated with [plot\\_line](#). The density of the color gives an indication of the likelihood of the value.

## Usage

```
plot_add_range(plot, ex_data)
```

## Arguments

plot	ggplot2 object.
ex_data	Dataframe with observed values for the specific timeseries.

## Value

ggplot2 plot.

## See Also

Other low-level-plot functions: [plot\\_add\\_box](#)

## Examples

```
# There is no external data so we need to add some noise first!
ex_data <- preprocess$biomass
ex_data$atoutput <- ex_data$atoutput * runif(n = nrow(ex_data), 0, 1)
ex_data$model <- "test"

# Create the timeseries with \link{plot_line}
plot <- plot_line(preprocess$biomass)

# Add the external data as geom_rug with \link{plot_add_range}
plot_add_range(plot, ex_data)
```

---

plot_bar	<i>Function to plot relative contribution of biomass and numbers per cohort.</i>
----------	--

---

### Description

Function to plot relative contribution of biomass and numbers per cohort.

### Usage

```
plot_bar(data, x = "time", y = "atoutput", fill = "species",
         wrap = NULL, ncol = NULL)
```

### Arguments

data	Dataframe to be plotted.
x	x-variable. Default is 'time'.
y	y-variable. Default is 'atoutput'.
fill	Column to use as filling colour. Default is "species".
wrap	Wrapping column. Default is 'species'
ncol	Number of columns in multipanel plot. Default is 7.

### Value

ggplot2 object

### See Also

Other plot functions: [plot\\_boxes](#), [plot\\_diet\\_bec\\_dev](#), [plot\\_diet](#), [plot\\_line](#), [plot\\_rec](#), [plot\\_species](#)

### Examples

```
plot_bar(preprocess$biomass)

# Most models have a large number of groups. Please make sure to combine groups with a low
# contribution prior to plotting with \link{combine_groups}.
df <- combine_groups(preprocess$biomass, group_col = "species", combine_thresh = 3)
plot_bar(df)

# This function can also be used to plot age-specific data.
plot_bar(preprocess$num_age, fill = "agecl", wrap = "species")

# Please use \link{agg_perc} to visualize the relative cohort structure over time.
df <- agg_perc(preprocess$num_age, groups = c("time", "species"))
plot_bar(df, fill = "agecl", wrap = "species")
```

---

plot_boxes	<i>Plot layout of boxes!</i>
------------	------------------------------

---

**Description**

Plot layout of boxes!

**Usage**

```
plot_boxes(data, color_boxes = TRUE)
```

**Arguments**

data            Dataframe to be plotted.  
color\_boxes    logical indicating if polygons should be color coded or not. Default is TRUE.

**Value**

ggplot2 object

**See Also**

Other plot functions: [plot\\_bar](#), [plot\\_diet\\_bec\\_dev](#), [plot\\_diet](#), [plot\\_line](#), [plot\\_rec](#), [plot\\_species](#)

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
bgm_data <- convert_bgm(file.path(d, "VMPA_setas.bgm"))

# Use color coding for polygons.
plot_boxes(bgm_data)

# Only use text to indicate polygons.
plot_boxes(bgm_data, color_boxes = FALSE)
```

---

plot_consumed_biomass	<i>Circle diagram to visualize the consumed biomass for the whole system.</i>
-----------------------	---

---

**Description**

Circle diagram to visualize the consumed biomass for the whole system.

**Usage**

```
plot_consumed_biomass(bio_consumed, select_time = NULL, show = 0.95)
```

**Arguments**

bio_consumed	Consumed biomass of prey groups by predatorgroup and agecl in tonnes for each timestep and polygon. Dataframe with columns 'pred', 'agecl', 'polygon', 'time', 'prey'. Consumed biomass in [t] is stored in column 'atoutput'. Should be generated with link{calculate_consumed_biomass}.
select_time	Numeric value to control the simulation time in years to visualize. By default the start of the simulation is shown. Default is NULL.
show	Numeric value between 0 - 1 to control the amount of links shown. Default is 0.95. Thus, the most important 95 are grouped together as 'Rest'.

**Value**

plot

**Examples**

```
## Not run:
# Need to fix NAs in init cdf at some point.
plot_consumed_biomass(ref_bio_cons)
plot_consumed_biomass(ref_bio_cons, select_time = 1.8)
plot_consumed_biomass(ref_bio_cons, select_time = 1.8, show = 0.99)

# Add some gns testing.
load("z:/Atlantis_models/Runs/dummy_01_ATLANTIS_NS/preprocess-north-sea.rda", verbose = T)
plot_consumed_biomass(result$biomass_consumed, show = 0.95, select_time = 2.2)

times <- c(0.2, seq(10, 90, 10), 99.8)
par(mfcol = c(3, 4))
for (i in seq_along(times)) {
  plot_consumed_biomass(result$biomass_consumed, show = 0.95, select_time = times[i])
}

times <- seq(0.2, 2.2, 0.2)
# times <- times[times != 1]
par(mfcol = c(3, 4))
for (i in seq_along(times)) {
  plot_consumed_biomass(result$biomass_consumed, show = 0.95, select_time = times[i])
}

## End(Not run)
```

**Description**

Visualize diet proportions from predator and prey perspective. The upper panel plot shows the predator perspective while the lower panel plot shows the prey perspective for a given group. Please note that this function only works with models based on the trunk code. Bec\_dev models should use [plot\\_diet\\_bec\\_dev](#) to get an indication of the feeding interactions.

**Usage**

```
plot_diet(bio_consumed, species = NULL, wrap_col = "agec1",
         combine_thresh = 7)
```

**Arguments**

bio_consumed	Consumed biomass of prey groups by predatorgroup and agec1 in tonnes for each timestep and polygon. Dataframe with columns 'pred', 'agec1', 'polygon', 'time', 'prey'. Consumed biomass in [t] is stored in column 'atoutput'. Should be generated with <code>link{calculate_consumed_biomass}</code> .
species	Character string giving the acronyms of the species you aim to plot. Default is NULL resulting in all available species being plotted.
wrap_col	Character specifying the column of the dataframe to be used as multipanel plot. Default is "agec1".
combine_thresh	Number of different categories to plot. Lets say predator X has eaten 20 different prey items. If you only want to show the 3 most important prey items set combine_thresh to 3. As rule of thumb values < 10 are useful otherwise too many colors are used in the plots. Default is 7.

**Value**

List of grobs composed of ggplot2 objects.

**See Also**

Other plot functions: [plot\\_bar](#), [plot\\_boxes](#), [plot\\_diet\\_bec\\_dev](#), [plot\\_line](#), [plot\\_rec](#), [plot\\_species](#)

**Examples**

```
## Not run:
plots <- plot_diet(ref_bio_cons, wrap_col = "agec1")
gridExtra::grid.arrange(plots[[1]])
gridExtra::grid.arrange(plots[[7]])

# Use names() to get the species names!
names(plots)

## End(Not run)

plot <- plot_diet(ref_bio_cons, species = "Small planktivorous fish", wrap_col = "agec1")
gridExtra::grid.arrange(plot[[1]])
```

---

plot\_diet\_bec\_dev      *Plot contribution of diet contents for each functional group.*

---

### Description

Visualize diet proportions from predator and prey perspective. The upper panel plot shows the predator perspective while the lower panel plot shows the prey perspective for a given group. Please note that this function uses the SpecPredMort.txt file to visualize feeding interactions and therefore is only an indication of the realized true diet within the model. Please use plot\_diet instead.

### Usage

```
plot_diet_bec_dev(data, species = NULL, wrap_col, combine_thresh = 15)
```

### Arguments

data	SpecPredMort.txt read in with load_spec_mort.
species	Character string giving the acronyms of the species you aim to plot. Default is NULL resulting in all available species being plotted.
wrap_col	Character specifying the column of the dataframe to be used as multipanel plot. In case you aim to plot SpecMort.txt data use either "agecl" or "stanza".
combine_thresh	Number of different categories to plot. Lets say predator X has eaten 20 different prey items. If you only want to show the 3 most important prey items set combine_thresh to 3. As rule of thumb values < 10 are useful otherwise too many colors are used in the plots. Default is 15.

### Value

List of ggplot2 objects.

### See Also

Other plot functions: [plot\\_bar](#), [plot\\_boxes](#), [plot\\_diet](#), [plot\\_line](#), [plot\\_rec](#), [plot\\_species](#)

### Examples

```
## Not run:
# Plot SpecMort.txt per ageclass.
plots <- plot_diet_bec_dev(preprocess_setas$diet_specmort, wrap_col = "agecl")
gridExtra::grid.arrange(plots[[1]])

# Only plot specific species
plots <- plot_diet_bec_dev(preprocess_setas$diet_specmort, species = "CEP", wrap_col = "agecl")
gridExtra::grid.arrange(plots[[1]])

# Plot SpecMort.txt per stanza First we need to transform the ageclasses to stanzas.
d <- system.file("extdata", "setas-model-new-becdev", package = "atlantistools")
diet_stanza <- combine_ages(dir = d,
```



```

                                data = preprocess_setas$diet_specmort,
                                col = "pred",
                                prm_biol = "VMPA_setas_biol_fishing_New.prm")
plots <- plot_diet_bec_dev(diet_stanza, wrap_col = "stanza")
gridExtra::grid.arrange(plots[[1]])

## End(Not run)

```

---

plot\_line

*Function to plot time series of atlantis ncdf output.*


---

### Description

Function to plot time series of atlantis ncdf output.

### Usage

```
plot_line(data, x = "time", y = "atoutput", wrap = "species",
          col = NULL, ncol = 7, yexpand = FALSE)
```

### Arguments

data	Dataframe to be plotted.
x	x-variable. Default is 'time'.
y	y-variable. Default is 'atoutput'.
wrap	Wrapping column. Default is 'species'
col	Column to use as colour. Default is NULL.
ncol	Number of columns in multipanel plot. Default is 7.
yexpand	Expands the y axis so it always includes 0. Default is FALSE.

### Value

ggplot2 object

### See Also

Other plot functions: [plot\\_bar](#), [plot\\_boxes](#), [plot\\_diet\\_bec\\_dev](#), [plot\\_diet](#), [plot\\_rec](#), [plot\\_species](#)

### Examples

```

plot_line(preprocess$biomass)
plot_line(preprocess$biomass, col = "species")
plot_line(preprocess$biomass_age, col = "agecl")
plot_line(preprocess$biomass_age, wrap = "agecl", col = "species")

# The function can also be used to compare model outoput with observed data.
d <- system.file("extdata", "setas-model-new-becdev", package = "atlantistools")

```

```

ex_data <- read.csv(file.path(d, "setas-bench.csv"), stringsAsFactors = FALSE)
names(ex_data)[names(ex_data) == "biomass"] <- "atoutput"

data <- preprocess$biomass
data$model <- "atlantis"
comp <- rbind(ex_data, data, stringsAsFactors = FALSE)

# Show atlantis as first factor!
lev_ord <- c("atlantis", sort(unique(comp$model))[sort(unique(comp$model)) != "atlantis"])
comp$model <- factor(comp$model, levels = lev_ord)

# Create plot
plot_line(comp, col = "model")

## Not run:
# Use \link{convert_relative_initial} and \link{plot_add_box}
# with \link{plot_line}. Use \link{convert_relative_initial} to
# generate a relative time series first.
df <- convert_relative_initial(preprocess$structn_age)

# Create the base plot with \link{plot_line}.
plot <- plot_line(df, col = "agecl")

# Add lower and upper range.
plot_add_box(plot)

# Create spatial timeseries plots in conjunction with \link{custom_grid}.
plot <- plot_line(preprocess$physics, wrap = NULL)
custom_grid(plot, grid_x = "polygon", grid_y = "variable")

plot <- plot_line(preprocess$flux, wrap = NULL, col = "variable")
custom_grid(plot, grid_x = "polygon", grid_y = "layer")

## End(Not run)

```

---

plot\_rec

*Plot recruitment.*


---

### Description

Plot recruitment.

### Usage

```
plot_rec(data, ex_data, ncol = 7)
```

### Arguments

data	Dataframe with information about ssb and recruits. This is created from atlantis output files YOY.txt and SSB.txt (Usually output[...]YOY.txt' & 'output[...]SSB.txt') using <a href="#">load_rec</a> .
------	---

ex\_data            Dataframe to compare the atlantis run with.  
 ncol              Number of columns in multipanel plot. Default is 7.

**Value**

ggplot2 object

**See Also**

Other plot functions: [plot\\_bar](#), [plot\\_boxes](#), [plot\\_diet\\_bec\\_dev](#), [plot\\_diet](#), [plot\\_line](#), [plot\\_species](#)

**Examples**

```
## Not run:
d <- system.file("extdata", "setas-model-new-becdev", package = "atlantistools")
ex_data <- read.csv(file.path(d, "setas-ssb-rec.csv"), stringsAsFactors = FALSE)
plot_rec(preprocess_setas$ssb_rec, ex_data)

## End(Not run)
```

---

plot_spatial_box	<i>Visualize the spatial distribution per species and stanza combination.</i>
------------------	---

---

**Description**

Visualize the spatial distribution per species and stanza combination.

**Usage**

```
plot_spatial_box(bio_spatial, bgm_as_df, select_species = NULL,
  timesteps = 2, polygon_overview = 0.2)
```

**Arguments**

bio_spatial	Biomass per group and stanza in tonnes for each timestep, layer and polygon. This dataframe should be generated with <a href="#">calculate_biomass_spatial</a> . The columns of the dataframe have to be 'species', 'species_stanza', 'polygon', 'layer', 'time' and 'atoutput'. Column 'atoutput' is the biomass in tonnes. Please use <a href="#">combine_ages</a> to transform an age-based dataframe to a stanza based dataframe.
bgm_as_df	*.bgm file converted to a dataframe. Please use <a href="#">convert_bgm</a> to convert your bgm-file to a dataframe with columns 'lat', 'long', 'inside_lat', 'inside_long' and 'polygon'.
select_species	Character vector listing the species to plot. If no species are selected NULL (default) all available species are plotted.
timesteps	Integer giving the number of timesteps to visualize. Default is 2. By default the start and end of the simulation is shown. In case timesteps > 2 equally spaced timesteps are added.

polygon\_overview

numeric value between 0 and 1 indicating the size used to plot the polygon overview in the upper right corner of the plot. Default is 0.2.

### Value

grob of 3 ggplot2 plots.

### Examples

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")

bgm_as_df <- convert_bgm(file.path(d, "VMPA_setas.bgm"))

# Spatial distribution in Atlantis is based on adu- and juv stanzas.
# Therefore, we need to aggregate the age-based biomass to
# stanzas with \link{combine_ages}.
bio_spatial <- combine_ages(ref_bio_sp, grp_col = "species", agemat = ref_agemat)

## Not run:
# Apply \link{plot_spatial_box}
grobs <- plot_spatial_box(bio_spatial, bgm_as_df, timesteps = 3)
gridExtra::grid.arrange(grobs[[1]])
gridExtra::grid.arrange(grobs[[9]])

# use names() to select specific plots
names(grobs)

## End(Not run)

# Plot specific species
grobs <- plot_spatial_box(bio_spatial, bgm_as_df,
                          select_species = "Shallow piscivorous fish", timesteps = 3)
gridExtra::grid.arrange(grobs[[1]])
```

---

plot\_spatial\_overlap *Plot spatial overlap.*

---

### Description

Plot spatial overlap.

### Usage

```
plot_spatial_overlap(df_list)
```

### Arguments

df\_list List of dataframes generated with [calculate\\_spatial\\_overlap](#)

**Value**

ggplot2 plot.

**Examples**

```
sp_overlap <- calculate_spatial_overlap(ref_bio_sp, ref_dietmatrix, ref_agemat)
## Not run:
plot_spatial_overlap(sp_overlap)

## End(Not run)
```

---

plot_spatial_ts	<i>Visualize the spatial distribution per species and stanza combination.</i>
-----------------	---

---

**Description**

Visualize the spatial distribution per species and stanza combination.

**Usage**

```
plot_spatial_ts(bio_spatial, bgm_as_df, vol, select_species = NULL,
  ncol = 7, polygon_overview = 0.2)
```

**Arguments**

bio_spatial	Biomass per group and stanza in tonnes for each timestep, layer and polygon. This dataframe should be generated with <a href="#">calculate_biomass_spatial</a> . The columns of the dataframe have to be 'species', 'species_stanza', 'polygon', 'layer', 'time' and 'atoutput'. Column 'atoutput' is the biomass in tonnes. Please use <a href="#">combine_ages</a> to transform an age-based dataframe to a stanza based dataframe.
bgm_as_df	*.bgm file converted to a dataframe. Please use <a href="#">convert_bgm</a> to convert your bgm-file to a dataframe with columns 'lat', 'long', 'inside_lat', 'inside_long' and 'polygon'.
vol	Volume per polygon and timestep. See model-preprocess.Rmd for details.
select_species	Character vector listing the species to plot. If no species are selected NULL (default) all available species are plotted.
ncol	Number of columns in final plot. Default is 7.
polygon_overview	numeric value between 0 and 1 indicating the size used to plot the polygon overview in the upper right corner of the plot. Default is 0.2.

**Value**

grob of 3 ggplot2 plots.

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")

bgm_as_df <- convert_bgm(file.path(d, "VMPA_setas.bgm"))
vol <- agg_data(ref_vol, groups = c("time", "polygon"), fun = sum, out = "volume")

# Spatial distribution in Atlantis is based on adu- and juv stanzas.
# Therefore, we need to aggregate the age-based biomass to
# stanzas with \link{combine_ages}.
bio_spatial <- combine_ages(ref_bio_sp, grp_col = "species", agemat = ref_agemat)

# Apply \link{plot_spatial_ts}
grobs <- plot_spatial_ts(bio_spatial, bgm_as_df, vol)
gridExtra::grid.arrange(grobs[[1]])
gridExtra::grid.arrange(grobs[[7]])
```

---

plot\_species

---

*Create species specific overview plot.*


---

**Description**

This plotting routine is based on Raphael's (Ifremer) plotting routine used during model calibration. Currently 6 plots are created by default: - Biomass over time - Biomass over time per age - StructN over time per age - ResN over time per age - Condition over time per age - Numbers over time per age

**Usage**

```
plot_species(data_pre, species)
```

**Arguments**

data_pre	List of preprocessed Atlantis simulation. The list of dataframes should be created with <code>model-preprocess.Rmd</code> .
species	Character string giving the name of the species to plot. Only age based species are supported.

**Value**

ggplot2 object of class grob

**See Also**

Other plot functions: [plot\\_bar](#), [plot\\_boxes](#), [plot\\_diet\\_bec\\_dev](#), [plot\\_diet](#), [plot\\_line](#), [plot\\_rec](#)

**Examples**

```
plot <- plot_species(preprocess, species = "Shallow piscivorous fish")  
# Use grid.arrange to draw the plot on the current device  
gridExtra::grid.arrange(plot)
```

---

preprocess

*preprocess*

---

**Description**

See "data-vignette-model-preprocess.R" and "model-preprocess.Rmd" for further information.

**Usage**

```
preprocess
```

**Format**

List with 17 dataframes.

1. biomass
2. biomass\_age
3. biomass\_consumed
4. biomass\_spatial\_stanza
5. diet
6. eat\_age
7. flux
8. grazing
9. growth\_age
10. growth\_rel\_init
11. nums
12. nums\_age
13. nums\_box
14. physics
15. resn\_age
16. sink
17. spatial\_overlap
18. ssb\_rec
19. structn\_age
20. vol

---

```
preprocess_txt      Preprocess dataframes loaded in with load_txt()
```

---

### Description

sep\_col is split into multiple columns given by into. If column ageclass is present and values start with 0 one is added to align with agestructure in other functions. Columns without any informations (length(unique()) == 1) are dropped. If the first time step only has zeros as values remove these values. remove zeros overall!

### Usage

```
preprocess_txt(df_txt, sep_col = "code", into)
```

### Arguments

df_txt	Dataframe read in with load_txt().
sep_col	Column to separate into multiple columns. Default is "code".
into	Character vector given the columns to split sep_col in.

### Value

Tidy dataframe.

### Examples

```
d <- system.file("extdata", "setas-model-new-becdev", package = "atlantistools")
df <- load_txt(file = file.path(d, "outputSETASSpecificPredMort.txt"))
df <- preprocess_txt(df_txt = df, into = c("pred", "agecl", "empty_col1", "prey", "empty_col2"))
head(df)

df <- load_txt(file = file.path(d, "outputSETASSpecificMort.txt"))
df <- preprocess_txt(df_txt = df, into = c("species", "agecl", "empty_col", "mort"))
head(df)
```

---

```
prm_to_df      Extract parameters from the biological parameter file and transform them to a dataframe.
```

---

### Description

Extract parameters from the biological parameter file and transform them to a dataframe.



**Usage**

```
prn_to_df(prn_biol, fgs, group, parameter)

prn_to_df_ages(prn_biol, fgs, group, parameter)
```

**Arguments**

prn_biol	Character string giving the connection to the biological parameterfile. The file-name usually contains biol_fishing and does end in .prm.
fgs	Character string giving the connection to the functional groups file. The file-name usually contains Groups and does end in .csv.
group	Character vector giving the functional Groups to extract.
parameter	Character vector giving the parameters to extract.

**Value**

Dataframe with columns 'species' and as many columns as parameters.

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
prn_biol <- file.path(d, "VMPA_setas_biol_fishing_Trunk.prm")
fgs <- file.path(d, "SETasGroupsDem_NoCep.csv")
group <- c("FPS", "FVS")
parameter <- c("mum", "C")

prn_to_df_ages(prn_biol, fgs, group, parameter)
prn_to_df(prn_biol, fgs, group, parameter)
```

---

ref_agemat	<i>agemat.</i>
------------	----------------

---

**Description**

*agemat.*

**Usage**

```
ref_agemat
```

**Format**

**species** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**age\_mat** First mature age class.

---

ref_bio_cons	<i>Consumed biomass.</i>
--------------	--------------------------

---

**Description**

Consumed biomass.

**Usage**

ref\_bio\_cons

**Format**

**pred** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**agecl** Ageclass given as integer from 1 to NumCohorts.

**polygon** Boxid starting from 0 to numboxes - 1.

**time** Simulation time in years. Modeltimestep was converted to actual time based on the settings in the 'run.prm' file.

**prey** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**atoutput** Observation column storing the actual output value. Consumed biomass in tonnes.

---

ref_bio_sp	<i>Spatial biomass.</i>
------------	-------------------------

---

**Description**

Spatial biomass.

**Usage**

ref\_bio\_sp

**Format**

**species** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**agecl** Ageclass given as integer from 1 to NumCohorts.

**polygon** Boxid starting from 0 to numboxes - 1.

**layer** Layerid starting from 0 to numlayers - 1.

**time** Simulation time in years. Modeltimestep was converted to actual time based on the settings in the 'run.prm' file.

**atoutput** Observation column storing the actual output value. Biomass in tonnes.

---

ref_dietmatrix	<i>Dietmatrix.</i>
----------------	--------------------

---

**Description**

See data-raw/data-create-reference-dfs.R for further information.

**Usage**

```
ref_dietmatrix
```

**Format**

**pred** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**pred\_stanza** Predator stanza. 1 = juvenile; 2 = adult.

**prey\_stanza** Prey stanza. 1 = juvenile; 2 = adult.

**code** Flag from the biological parameter file.

**prey** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**avail** Observation column storing the actual output value. Availability ranging from 0 to 1.

**prey\_id** Preyid index based on the functional groups file.

---

ref_dm	<i>Dietcheck.</i>
--------	-------------------

---

**Description**

Dietcheck.

**Usage**

```
ref_dm
```

**Format**

**time** Simulation time in years. Modeltimestep was converted to actual time based on the settings in the 'run.prm' file.

**pred** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**agecl** Ageclass given as integer from 1 to NumCohorts.

**prey** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**atoutput** Observation column storing the actual output value. Diet contribution in percentage.

---

ref_eat	<i>Eat.</i>
---------	-------------

---

**Description**

Eat.

**Usage**

ref\_eat

**Format**

**species** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**agecl** Ageclass given as integer from 1 to NumCohorts.

**polygon** Boxid starting from 0 to numboxes - 1.

**time** Simulation time in years. Modeltimestep was converted to actual time based on the settings in the 'run.prm' file.

**atoutput** Obseravtion column storing the actual output value. Consumption in mg N m<sup>-3</sup> d<sup>-1</sup>.

---

ref_grazing	<i>Grazing.</i>
-------------	-----------------

---

**Description**

Grazing.

**Usage**

ref\_grazing

**Format**

**species** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**agecl** Ageclass given as integer from 1 to NumCohorts.

**polygon** Boxid starting from 0 to numboxes - 1.

**time** Simulation time in years. Modeltimestep was converted to actual time based on the settings in the 'run.prm' file.

**atoutput** Obseravtion column storing the actual output value. Grazing in mg N m<sup>-3</sup> d<sup>-1</sup>.

---

ref_growth	<i>Growth.</i>
------------	----------------

---

**Description**

Growth.

**Usage**

ref\_growth

**Format**

**species** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**agecl** Ageclass given as integer from 1 to NumCohorts.

**polygon** Boxid starting from 0 to numboxes - 1.

**time** Simulation time in years. Modeltimestep was converted to actual time based on the settings in the 'run.prm' file.

**atoutput** Obseravtion column storing the actual output value. Growth in mg N d-1.

---

ref_n	<i>Nitrogen.</i>
-------	------------------

---

**Description**

Nitrogen.

**Usage**

ref\_n

**Format**

**species** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**polygon** Boxid starting from 0 to numboxes - 1.

**layer** Layerid starting from 0 to numlayers - 1.

**time** Simulation time in years. Modeltimestep was converted to actual time based on the settings in the 'run.prm' file.

**atoutput** Obseravtion column storing the actual output value. Nitrogen in mg N m-3.

---

ref_nums	<i>Numbers at age data.</i>
----------	-----------------------------

---

**Description**

Numbers at age data.

**Usage**

ref\_nums

**Format**

**species** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**agecl** Ageclass given as integer from 1 to NumCohorts.

**polygon** Boxid starting from 0 to numboxes - 1.

**layer** Layerid starting from 0 to numlayers - 1.

**time** Simulation time in years. Modeltimestep was converted to actual time based on the settings in the 'run.prm' file.

**atoutput** Obseravtion column storing the actual output value. Numbers.

---

ref_physics	<i>Physical variables.</i>
-------------	----------------------------

---

**Description**

Physical variables.

**Usage**

ref\_physics

**Format**

**variable** Name of the physical variable: "salt", "NO3", "NH3", "Temp", "Chl\_a" and "Denitrification".

**polygon** Boxid starting from 0 to numboxes - 1.

**layer** Layerid starting from 0 to numlayers - 1.

**time** Simulation time in years. Modeltimestep was converted to actual time based on the settings in the 'run.prm' file.

**atoutput** Obseravtion column storing the actual output value. units are salt = PSU; NO3, NH3 = mg N m-3; Temp = degrees Celcius; Chl\_a, Denitrification = ?

---

ref_resn	<i>Reserve nitrogen.</i>
----------	--------------------------

---

**Description**

Reserve nitrogen.

**Usage**

ref\_resn

**Format**

**species** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**agecl** Ageclass given as integer from 1 to NumCohorts.

**polygon** Boxid starting from 0 to numboxes - 1.

**layer** Layerid starting from 0 to numlayers - 1.

**time** Simulation time in years. Modeltimestep was converted to actual time based on the settings in the 'run.prm' file.

**atoutput** Observation column storing the actual output value. Reserve weight in mg N.

---

ref_structn	<i>Structural nitrogen.</i>
-------------	-----------------------------

---

**Description**

Structural nitrogen.

**Usage**

ref\_structn

**Format**

**species** Name of the functional groups given as character string. The names match with the column 'LongName' in the functionalGroups.csv file.

**agecl** Ageclass given as integer from 1 to NumCohorts.

**polygon** Boxid starting from 0 to numboxes - 1.

**layer** Layerid starting from 0 to numlayers - 1.

**time** Simulation time in years. Modeltimestep was converted to actual time based on the settings in the 'run.prm' file.

**atoutput** Observation column storing the actual output value. Structural weight in mg N.

---

ref_to_bibkey	<i>Convert reference to bib-tex-key.</i>
---------------	--

---

**Description**

Convert reference to bib-tex-key.

**Usage**

```
ref_to_bibkey(ref_df, bib)
```

**Arguments**

ref_df	dataframe with columns author, year, title in case title is missing match is performed based on author and year only.
bib	character string giving the name of the .bib file.

**Value**

Character vector.

---

ref_vol	<i>Volume.</i>
---------	----------------

---

**Description**

Volume.

**Usage**

```
ref_vol
```

**Format**

**variable** Name of the physical variable: "volume".

**polygon** Boxid starting from 0 to numboxes - 1.

**layer** Layerid starting from 0 to numlayers - 1.

**time** Simulation time in years. Modeltimestep was converted to actual time based on the settings in the 'run.prm' file.

**atoutput** Obseravtion column storing the actual output value. Volume in m<sup>3</sup>



---

ref_vol_dz	<i>Volume and dz.</i>
------------	-----------------------

---

**Description**

Volume and dz.

**Usage**

ref\_vol\_dz

**Format**

**variable** Name of the physical variable: "volume" and "dz".

**polygon** Boxid starting from 0 to numboxes - 1.

**layer** Layerid starting from 0 to numlayers - 1.

**time** Simulation time in years. Modeltimestep was converted to actual time based on the settings in the 'run.prm' file.

**atoutput** Obseravtion column storing the actual output value. volume in m<sup>3</sup> dz in m

---

scan_prm	<i>Scan character vector for specific flag!</i>
----------	---

---

**Description**

Scan character vector for specific flag!

**Usage**

scan\_prm(chars, variable)

**Arguments**

chars            Vector of character strings

variable        Character string giving the flag to search for.

---

scan\_reference\_fishbase

*Scan list of references for character string for fish species*

---

### Description

Scan list of references for character string for fish species

### Usage

```
scan_reference_fishbase(fish, chr, mirror = "se")
```

### Arguments

fish	Vector of fish species with genus and species information.
chr	Character string to search.
mirror	Character string defining the url mirror to use. Defaults to se. In case data extraction is slow use a different mirror. Try to avoid frequently used mirrors like uk or com.

### Value

Dataframe of potentially relevant references.

### Examples

```
## Not run:
# For some reason the examples break with appveyor.
fish <- c("Gadus morhua", "Merlangius merlangus")
df <- scan_reference_fishbase(fish, chr = "diet")
df <- scan_reference_fishbase(fish, chr = "xxx")
df <- scan_reference_fishbase(fish, chr = "feed")

## End(Not run)
```

---

sc\_init

*Sanity check initial conditions file*

---

### Description

Sanity check initial conditions file

**Usage**

```
sc_init(init, prm_biol, fgs, bboxes, pred = NULL, set_avail = NULL,
        version_flag = 2)

plot_sc_init(df, mult_mum, mult_c, pred = NULL)
```

**Arguments**

init	Character string giving the connection of the initial conditions netcdf file. The filename usually contains init and ends in .nc.
prm_biol	Character string giving the connection to the biological parameterfile. The file-name usually contains biol_fishing and does end in .prm.
fgs	Character string giving the connection to the functional groups file. The file-name usually contains Groups and does end in .csv.
bboxes	Integer vector giving the box-id of the boundary boxes. Can be created with get_boundary.
pred	Vector of predator acronyms to check. If NULL (default) all age based predators are selected.
set_avail	Numeric value. All present availabilities can be set to a specific value. Default value is NULL which results in no changes to the present availability matrix.
version_flag	The version of ATLANTIS model. 1 for bec_dev, 2 for trunk. default is 2..
df	Dataframe to pass to plot_sc_init(). df should be generated with sc_init or read in from *.rda (also generated with sc_init()).
mult_mum	Numeric vector of multiplication factors applied to the initial mum values.
mult_c	Numeric vector of multiplication factors applied to the initial C values.

**Value**

Dataframe/ Plot.

**Examples**

```
d <- system.file("extdata", "setas-model-new-trunk", package = "atlantistools")
init <- file.path(d, "INIT_VMPA_Jan2015.nc")
prm_biol <- file.path(d, "VMPA_setas_biol_fishing_Trunk.prm")
fgs <- file.path(d, "SETasGroupsDem_NoCep.csv")
bboxes <- get_boundary(load_box(bgm = file.path(d, "VMPA_setas.bgm")))

data1 <- sc_init(init, prm_biol, fgs, bboxes)
## Not run:
dir <- system.file("extdata", "gns", package = "atlantistools")
fgs <- "functionalGroups.csv"
init <- "init_simple_NorthSea.nc"
prm_biol <- "NorthSea_biol_fishing.prm"
bboxes <- get_boundary(load_box(dir = dir, bgm = "NorthSea.bgm"))
mult_mum <- seq(0.5, 10, by = 1)
mult_c <- seq(0.5, 10, by = 1)
```

```
no_avail <- FALSE
save_to_disc <- FALSE
data1 <- sc_init(dir, init, prm_biol, fgs, bboxes, save_to_disc = FALSE)
plot_sc_init(df = data1, mult_mum, mult_c)
plot_sc_init(df = data1, mult_mum, mult_c, pred = "Cod")

data2 <- sc_init(dir, init, prm_biol, fgs, bboxes, pred = "Cod", save_to_disc = FALSE)
plot_sc_init(df = data2, mult_mum, mult_c)

## End(Not run)
```

---

str_split_twice	<i>Extract numeric values from string.</i>
-----------------	--

---

### Description

The function splits any character string at each tab and space and returns all (`min_only = FALSE`) or only the first (`min_only = T`) numeric value found in the string.

### Usage

```
str_split_twice(char, min_only = TRUE)
```

### Arguments

char	Character string.
min_only	Logical specifying if only the first numeric value (TRUE) or all numeric values (FALSE) should be returned. Default is TRUE.

### Value

numeric values inside char string.

### Examples

```
str_split_twice(char = "Hello 15")
str_split_twice(char = "flag1 15 16\t15", min_only = FALSE)
```

---

theme_atlantis	<i>Customized theme used in all plots.</i>
----------------	--

---

### Description

This function is a customized theme for ggplot2 plots. It's applied by default to all plots created within atlantistools.

### Usage

```
theme_atlantis(large = 22, medium = 18, small = 14, scale_font = 1,
  rot_xaxis_text = FALSE, rot_strips_y = TRUE)
```

### Arguments

large	Integer giving the size of the font for the main parts of the plot. Default is 22.
medium	Integer giving the size of the font used in the legend and facet labels. Default is 18.
small	Integer giving the size of the font used in the rest of the plot. Default is 14.
scale_font	Numeric used to scale all font sizes. Default is 1.
rot_xaxis_text	Logical indicating if x-axis text should be rotated by 45 degree. Default is FALSE.
rot_strips_y	Logical indicating if facet labels should be rotated by 90 degree. Default is TRUE.

### Examples

```
## Not run: nums_agg <- agg_data(data = ref_nums, groups = c("species", "time"), fun = sum)
ggplot2::ggplot(data = nums_agg, ggplot2::aes(x = time, y = atoutput)) +
  ggplot2::facet_wrap(~species) +
  theme_atlantis()
## End(Not run)
```

---

%>%	<i>Pipeoperator</i>
-----	---------------------

---

### Description

Atlantistools makes heavy use of dplyr data transformations. Therefore it is advisable to import the pipeoperator %>% from magrittr.

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