IBMPopSim package description

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Contents

1 General features and package installation  ........................................ 2
  1.1 Brief overview of Individual-Based-Models (IBMs) ......................... 2
    1.1.1 Events type and associated kernel ........................................ 2
    1.1.2 Events intensities .......................................................... 3
    1.1.3 Simulation algorithm ....................................................... 3
  1.2 Overview of package structure .................................................. 4
  1.3 Installation ........................................................................... 4
    1.3.1 First example to check installation ..................................... 4
  1.4 Quick model creation and simulation ........................................... 5

2 Package description ................................................................. 7
  2.1 Population, individuals, characteristics ...................................... 7
  2.2 Model parameters ................................................................... 8
  2.3 Events creations ....................................................................... 9
    2.3.1 Event kernel code ............................................................... 10
    2.3.2 Event creation with individual intensity ................................. 11
    2.3.3 Event creation with interaction intensity ............................... 12
    2.3.4 Events creation with Poisson and Inhomogeneous Poisson intensity 13
  2.4 Model creation ....................................................................... 14
  2.5 Simulation and outputs ............................................................. 15
    2.5.1 Events bounds .................................................................... 16
    2.5.2 Simulation without swap events .............................................. 16
    2.5.3 Outputs ............................................................................. 18
    2.5.4 Simple multithreading .......................................................... 19
  2.6 Simulation with swap events ..................................................... 19
    2.6.1 Vector of times in \texttt{popsim} ............................................. 19
    2.6.2 Individual life courses .......................................................... 21
3 C++ essentials

3.1 C++ syntax .................................................. 22
3.2 Usual numeric functions .................................. 24
3.3 Individuals characteristics and model parameters: link between R and C++ ................. 24
   3.3.1 Atomic types (characteristics and parameters) ....................................... 24
   3.3.2 Vectors and matrices (model parameters) ........................................... 25
   3.3.3 Predefined functions (model parameters) ........................................... 26
3.4 Random variables ........................................... 29

4 References ...................................................... 30

The IBMPopSim package is conceived to simulate the random evolution of structured population dynamics, called stochastic Individual Based Models (IBMs), based on the description of individual behaviors.

The package IBMPopSim allows the efficient simulation of a wide class IBMs where individuals are marked by their date of birth and a set of (discrete or continuous) characteristics. The scope of applications includes population models in actuarial science, biology, demography, ecology, epidemiology or economy.

An IBM is summarized by the different types events occurring in the population at random dates. These events include the birth/arrival or death/exit of an individual in the population, and individuals changing of characteristics. Once the events modifying the population have been defined, the last step before simulating the population evolution is to specify the so-called events intensities, describing the frequency at which the different types of events occur in the population.

1 General features and package installation

1.1 Brief overview of Individual-Based-Models (IBMs)

Stochastic IBMs can be defined as a general class of population dynamics models, where different types of events can occur to individuals at random dates, depending on their age (or equivalently their date of birth) and characteristics (gender, location, size, socioeconomic class . . . ), and interactions with others.

An IBM can be summarized by the different types events that can occur to individuals and modify the population composition, and how frequently they happen. Let us start by recalling briefly these features before going into detailed on how they can be defined in IBMPopSim.

1.1.1 Events type and associated kernel

In IBMPopSim, the population can evolve according to six main types of events:

- **Birth event**: addition of an individual of age 0 to the population.
- **Death event**: removal of an individual from the population.
- **Entry event**: arrival of an individual in the population.
- **Exit (emigration) event**: exit from the population (other than death).
- **Swap event**: an individual changes of characteristics.
- **Custom event**.

With each event type is associated an **event kernel**, describing how the population is modified following the occurrence of the event. For instance, if an individual \( I \) dies or exit a time \( t \), the individual is removed from the population.
For other types of events, the event kernel has to be specified. Indeed, if an new individual enters the population (entry event), the rule for choosing the age and characteristics of this new individual has to be defined. For instance, the characteristics of a new individual in the population can be chosen uniformly in the space of all characteristics, or can depends on the distribution of his parents or those of the other individuals composing the population.

### 1.1.2 Events intensities

Once the different types or events and their action have been defined, it remains to define how frequently they can occur to define properly an IBM.

An event intensity is a function $\lambda^e(I, t)$ describing the frequency at which an event $e$ can occur to an individual $I$ in the population at a time $t$. Informally, given an history of the population $(\mathcal{F}_t)$, the probability of event $e$ occurring to individual $I$ during a small interval of time $(t, t + dt]$ is proportional to $\lambda^e(I, t)$:

$$\mathbb{P}(\text{event } e \text{ occurring to } I \text{ during } (t, t+dt)|\mathcal{F}_t) = \lambda^e(I, t)dt. \quad (1)$$

The intensity function $\lambda^e$ can depend on:

- The individual’s age and characteristics,
- The time $t$,
- Interactions with other individuals in the population,
- Other model parameters.

When the intensity $\lambda^e$ does not depends on the other individuals living in the population (there are no interactions for this specific event), the intensity function is said to be in the class of individual intensity functions. This is case for instance when the death intensity of individuals only depend on their age and characteristics. Otherwise, the intensity function is said to be the class of interaction functions. In this package, we mainly focus on quadratic interactions, that is of intensity functions of the form

$$\lambda^e(I, t) = \sum_{J \in \text{pop}} U(I, J, t). \quad (2)$$

The interaction function $U(I, J, t)$ represent the interaction between two individuals $I$ and $J$ at time $t$. See for instance the vignette(‘IBMPopSim_interaction’) for examples of intensities with interactions.

Sometimes, the intensity refers to the frequency at which an event can occur in all the population. This is the case for instance when individuals enter the population at a constant rate $C^e$ (intensity of class poisson), or at a rate $(\Lambda^e_t)$ depending on time (intensity of class inhomogeneous_poisson). In this case, the intensity $\Lambda^e_t$ is informally defined as

$$\mathbb{P}(\text{event } e \text{ occurring in the population during } (t, t+dt)|\mathcal{F}_t) = \Lambda^e_t dt. \quad (3)$$

### 1.1.3 Simulation algorithm

The IBM simulation algorithm is based on an acceptance-rejection method for simulating random times called thinning, generalizing the algorithm proposed by (Fournier and Méléard 2004) (see also (Ferrière and Tran 2009), (Bonmezoued 2016)). The main idea of the algorithm is to use candidate event times with constant intensity $\bar{\lambda}$ (which corresponds to exponential variables with parameter $\bar{\lambda}$). The constant $\bar{\lambda}$ must be greater than the true event intensity $\lambda^e(I, t)$ for all individuals $I$:

$$\lambda^e(I, t) \leq \bar{\lambda}, \quad \forall I \in \text{pop}, \ t \in [0, T]. \quad (4)$$
Starting from time $t$, once a candidate event time $t + \tau^e$ has been proposed for a given event $e$ and individual $I$, it remains to accept or reject the candidate event with probability $\frac{\lambda^e(I, t + \tau^e)}{\bar{\lambda}}$. If the candidate event time is accepted, then the event $e$ occurs to the individual $I$ at time $t + \tau^e$. The main idea of the algorithm implemented can be summarized as follows:

**Thinning algorithm**

1. Draw a candidate time $t + \tau^e$ for event $e$, $\tau^e \sim E(\bar{\lambda})$, and an individual $I$.
2. Draw a uniform variable $\theta \sim U([0, 1])$.
3. If $\theta \leq \frac{\lambda^e(I, t + \tau^e)}{\lambda}$ then Apply event kernel to individual $I$, else Do nothing and start again from $t + \tau^e$.

In the presence of interactions, the intensity of an individual $I$ is computed making the sum over all individuals $J$ in the population of the interactions between $I$ and $J$, which can be very time consuming. We introduce in IBMPopSim a “randomized” algorithm, in which this computation is by picking randomly another individual in the population and computing its interaction function with the given individual. In this case the algorithm presented above is simply modified as follows: two individuals are picked, the intensity function $\lambda^e$ in the algorithm is replaced by the interaction function $U^e$, and $\bar{\lambda}$ by a bound of $U^e$.

1.2 Overview of package structure

Emphasis is placed on code efficiency (speed of execution) and ease of implementation, while trying to be as general as possible.

The implementation of an IBM model is based on a few building blocks, easy to manipulate. For code efficiency, we have chosen to let the user write a few lines of C++ code in R interface to define events intensity and action kernel. The user’s code is concatenated with internal code in the package. Compilation is done via the `sourceCpp` call of the Rcpp package. The code produced is usually very fast and can be multithreaded.

Thanks to model parameters and predefined functions, the C++ code to be written is simple and does not require a great knowledge of C++: it is enough to respect the rules of syntax, to do arithmetic operations and tests. Furthermore, various inputs of the model can be modified without having to recompile the code.

1.3 Installation

The latest stable version can be installed from CRAN:

```r
install.packages('IBMPopSim')
```

The latest development version can be installed from github:

```r
# install.packages("devtools")
devtools::install_github("DaphneGiorgi/IBMPopSim")
```

1.3.1 First example to check installation

To illustrate the use of the package and to check the installation, a simple model with Poisson arrivals and exits is implemented.


```r
library(IBMPopSim)

init_size <- 100000
pop <- data.frame(birth = rep(0, init_size), death = NA)
birth = mk_event_poisson(type = 'birth', intensity = 'lambda')
death = mk_event_poisson(type = 'death', intensity = 'mu')
params = list('lambda' = 100, 'mu' = 100)

# mk_model compiles C++ code using sourceCpp from Rcpp
birth_death <- mk_model(events = list(birth, death),
parameters = params)
```

If there are no errors then the C++ built environment is compatible with the package. The model is created and a C++ code has been compiled. The simulation is done using the `popsim` function.

```r
sim_out <- popsim(model = birth_death,
population = pop,
events_bounds = c('birth' = params$lambda, 'death' = params$mu),
parameters = params,
time = 10)

num_births <- length(sim_out$population$birth) - init_size
num_deaths <- sum(!is.na(sim_out$population$death))
print(c("births" = num_births, "deaths" = num_deaths))
## births deaths
## 999 995
```

### 1.4 Quick model creation and simulation

Before going into details on the different steps for defining an IBM in IBMPopSim, let us first go quickly through an example of model creation and simulation. In order to define a model, the user must specify three building blocks:

- An initial population (see Section 2.1),
- The model parameters (see Section 2.2),
- The different events that can occur and their intensity (see Section 2.3).

The model can then be created from these three blocks.

- We take here an initial population, stored in a data frame, composed of 100 000 individuals marked by their gender (encoded by a Boolean characteristic):

```r
pop <- EW_pop_14$sample
```

- The second step is to define the model parameters’ list:

```r
params <- list("alpha" = 0.008, "beta" = 0.02, "p_male" = 0.51,
"birth_rate" = stepfun(c(15,40), c(0,0.05,0)))
```
• The last step is to defined the events that can occur in the population, here birth and death events:

```r
death_event <- mk_event_individual(type = "death",
           intensity_code = "result = alpha * exp(beta * age(I, t));")

birth_event <- mk_event_individual(type = "birth",
           intensity_code = "result = birth_rate(age(I,t));",
           kernel_code = "newI.male = CUnif(0, 1) < p_male;")
```

Note that these events contain some C++ statements that depend (implicitly) on the previously declared parameters in variable `params`.

• Finally, the model is created by calling the function `mk_model`. A C++ source code is obtained from the events and parameters, then compiled using the `sourceCpp` function of the `Rcpp` package.

```r
model <- mk_model(characteristics = get_characteristics(pop),
                   events = list(death_event, birth_event),
                   parameters = params)
```

• In order to simulate a random trajectory of the population until a given time $T$ bounds on the events intensities have to be specified:

```r
a_max <- 115
events_bounds = c("death" = params$alpha * exp(params$beta * a_max),
              "birth" = max(params$birth_rate))
```

Then, the function `popsim` can be called:

```r
sim_out <- popsim(model, pop, events_bounds, params,
                   age_max = a_max, time = 30)
## Simulation on [0, 30]
```

• The data frame `sim_out$population` contains the information (date of birth, date of death, gender) on individuals who lived in the population over the period $[0, 30]$. Functions of the package allows to provide aggregated information on the population.

```r
pop_out <- sim_out$population
head(pop_out)
## birth death male
## 1 -84.97524 NA FALSE
## 2 -84.96228 NA TRUE
## 3 -84.96043 NA FALSE
## 4 -84.95083 NA FALSE
## 5 -84.91351 NA TRUE
## 6 -84.91099 NA FALSE
female_pop <- pop_out[pop_out$male==FALSE, ]
age_pyramid(female_pop, ages = 85:90, time = 30)
## age male value
## 1 85 - 86 FALSE 230
## 2 86 - 87 FALSE 235
## 3 87 - 88 FALSE 218
```
Note that parameters of the model can be changed without recompiling the model.

params$beta <- 0.01

# Update death event bound:
events_bounds["death"] <- params$alpha * exp(params$beta * a_max)

sim_out <- popsimsim(model, pop, events_bounds, params,
                      age_max = a_max, time = 30)

2 Package description

The first step to define a model in IBMPopSim is to define the population structure.

2.1 Population, individuals, characteristics

A population is a data frame in which each row corresponds to an individual, and which has at least two columns:

- A column `birth` containing the birth dates of individuals in the population.
- A column `death` containing the birth dates of individuals in the population (NA for alive individuals).

The data frame can contain more than two columns if individuals are described by additional characteristics such as gender, size, spatial location... In the example below, individuals are described by their birth and death dates, as well a Boolean characteristic called `male`. For instance, the first individual is a female whose age at \( t_0 \) = 0 is \( t_0 - (-106.9055) = 106.9055 \).

head(pop)

```r
## birth death male
## 1 -106.9055 NA FALSE
## 2 -106.8303 NA FALSE
## 3 -104.5097 NA TRUE
## 4 -104.2218 NA FALSE
## 5 -103.5225 NA FALSE
## 6 -103.3644 NA FALSE
```

- **Type of a characteristic.** A characteristic must be of atomic type: logical (bool in C++), integer (int), double or character (char). The function \texttt{?get_characteristic} allows to easily get characteristics names and their types (in R and C++) from a population data frame. See also section 3.3.

- **Name of a characteristic.** We draw the attention to the fact that some names for characteristics are forbidden, or are reserved to specific cases. The reserved words include:
  - birth and death: which are of type double, referring to dates of the events of birth and death,
- **male**: can only be used for a Boolean characteristics, usually referring to the individuals' sex/gender,
- **entry**: can only be used with an event of type **entry**. This characteristic is a **double** containing the date of entry in the population of the individual.
- **out**: can only be used with an event of type **exit**. This characteristic is a Boolean which is **TRUE** if the individual came out of the population at time **death**
- **id**: can only be used in order to defined the individuals unique identifier (see Section 2.6),
- **time**, **I**, **J**, **pop**, **newI**, **t** and **k**: forbidden as characteristics or variable name,
- C++ reserved words: forbidden as characteristics or variable name. Note that **char** is a reserved word. See here for a comprehensive list of C++ reserved words.

### An individual in C++

In the C++ compiled model, an individual **I** is an object of the C++ class **individual** containing some attributes:

- **birth_date** and **death_date** which are **double**,
- the characteristics,
- and some methods:
  * **I.age(t)**: a **const** method returning the age of an individual **I** at time **t**,
  * **I.set_age(a, t)**: to set the age **a** at time **t** of an individual **I** (set birth_date at **t**-**a**),
  * **I.is_dead(t)**: a **const** method returning **true** is the individual **I** is dead at time **t**. For convenience, there is also a global function **age** to get the age of an individual **I** so that **age(I, t)** is equivalent to **I.age(t)**.

### Using characteristic in C++ code

If **Chi** is the name of a characteristic, then **I.Chi** returns the value of the characteristic **Chi** for the individual **I**. For instance, in the previous example **I.male** would equal to **true** if **I** is a male or **false** if **I** is a female.

### A population in C++

During the model creation, a population of individuals is automatically created from the initial population data frame. In C++ we denote by **pop** the current population which is an object of a internal class containing some methods:

- an operator [] to access the k-th individual, note that most of time we denote k the current individual such that **pop[k]** is equivalent to **I** (in fact **I** is defined as a reference on **pop[k]**),
- **pop.kill(k, t)** to kill the individual **k** at time **t**.

### 2.2 Model parameters

A list of model parameters can be defined to simplify the events creation (see also section 3.3 for more details). These parameters can be of various types:

- **Atomic type**: likewise characteristics, a parameter may be **logical**, **integer**, **double** or **character** (and internal in C++ there is conversion to **bool**, **int**, **double** and **char** respectively),
- **Vector or matrix**: we call vector a **Numeric** of **double** of length larger than 1 (and dimension 1), we use RcppArmadillo to convert these types in **arma::vec** and **arma::mat** in C++,
- **Predefined function of one variable**: such functions are **stepfun**, **linfun**, **gompertz**, **weibull**, **piecewise_x**, in C++ these functions are **function_x**,
- **Piecewise function of two variable**: **piecewise_xy** which is **function_xy** in C++,
- **List of predefined functions**: of same C++ type, either **function_x** or **function_xy**.

For example,

```r
params <- list("coeff" = 1.1,
               "death_function" = gompertz(0.1, 0.005))
```
creates a numeric parameter \texttt{coeff} and a Gompertz function \texttt{death_function} which can be used when creating the events of the model.

**Predefined IBMPopSim functions** are classes of functions predefined in the package to simplify models creation. For instance, \texttt{gompertz(a,b)} returns the function \( g(x) = a \exp(bx) \) and \texttt{stepfun} is the usual R function for the creation of step functions. Another example is \texttt{piecewise_xy} which allow the creation of age and time functions. See Section 3.3.3 for a description of all IBMPopSim functions, and Section 3.3.2 for more details of Armadillo objects.

## 2.3 Events creations

The last and most important step of the model creation is the events creation. The call to the function creating an event is of form,

\[
\texttt{mk_event_CLASS(type = "TYPE", name ="NAME", kernel_code = "KERNEL_CODE", ...)}
\]

where \texttt{CLASS} is replaced by the class of the event intensity, \texttt{type} is the event type and \texttt{kernel_code} the event kernel.

The other arguments depend on the intensity class. Tables below summarize the different intensity classes and types of events introduced in Section 1.1.1.

\begin{table}[h]
\centering
\begin{tabular}{ll}
Event type & Type \\
Birth & birth \\
Entry & entry \\
Death & death \\
Exit & exit \\
Swap & swap \\
Custom & custom \\
\end{tabular}
\caption{Event types}
\end{table}

\begin{table}[h]
\centering
\begin{tabular}{ll}
Intensity class & Class \\
Poisson & poisson \\
Inhomogeneous Poisson & inhomogeneous_poisson \\
Individual & individual \\
Interaction & interaction \\
\end{tabular}
\caption{Intensity classes}
\end{table}

The intensity function and kernel of an event are defined through arguments of the function \texttt{mk_event_CLASS}. These arguments are strings composed of few lines of code defining an event frequency and its action of the event on individuals. Since the model is compiled using Rcpp, the code should be written in C++. However, thanks to the model parameters and functions/variables of the package, even the non-experienced C++ user can define a model quite easily. Several examples are given in the several vignettes of this package, and basic C++ tools are presented in Section 3.

The optional argument \texttt{name} gives a name to the event. If not specified, the name of the event is its type, for instance \texttt{death}. However, a name must be specified if the model is composed of several events with the same type.
2.3.1 Event kernel code

The argument kernel_code is NULL by default and doesn’t have to be specified for death and exit events. For those types of events, the event kernel is automatically generated during the model creation.

1. Death event

If the user defines a death event, the death date of the current individual I is set to the current time t. No kernel has to be specified by the user.

2. Exit event

If an exit event is defined, a characteristic out is automatically added to individuals in the population. When an individual I exits the population, I.out is set to TRUE and his exit time is recorded as a “death” date.

3. Birth event

For a birth event, the default generated event kernel is that an individual I gives birth to a new individual newI of age 0 at the current time t, which has the same characteristics than his parent I. If no kernel is specified, the default generated C++ code for a birth event is:

```cpp
individual newI = I;
newI.birth_date = t;
pop.add(newI);
```

The user can modify the birth kernel, by specify the argument kernel_code of mk_event_CLASS. In this case, the generated code is

```cpp
individual newI = I;
newI.birth_date = t;
_KERNEL_CODE_
pop.add(newI);
```

where _KERNEL_CODE_ is replaced by the content of the kernel_code argument. For instance, in a population where individuals are characterized by their gender, the kernel code

```cpp
birth_kernel_code <- "newI.male = (CUnif(0, 1) < p_male);"
```

creates new individuals which are males with probability p_male, or females otherwise. Here, p_male should be included in the list of model parameters.

4. Entry event

If an entry event is defined, a characteristic entry is automatically added to individuals in the population. When an individual I enters the population, I.entry is set as the date at which the individual enters the population. When an entry occurs the individual entering the population is not of age 0. In this case, the user must specify the kernel_code argument indicating how the age and characteristics of the new individual are chosen. For instance, with
entry_kernel_code <- "double a = CUnif(20,40);
    newI.set_age(a,t);
    newI.male = (CUnif(0, 1) < p_male);"

An individual who enters the population has an age $a$ taken uniformly in [20, 40], and is a male with probability $p\_male$. His characteristic $I\_entry$ is equal to $t$.

The available variables for the birth and entry events C++ kernel codes are:

<table>
<thead>
<tr>
<th>Variable</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>$I$</td>
<td>Current individual</td>
</tr>
<tr>
<td>$t$</td>
<td>Current time</td>
</tr>
<tr>
<td>$pop$</td>
<td>Current population (vector)</td>
</tr>
<tr>
<td>$newI$</td>
<td>New individual. By default for birth events newI = I with newI.birth = t</td>
</tr>
<tr>
<td>Model parameters</td>
<td>Depends on the model</td>
</tr>
</tbody>
</table>

When there are several entry events, the user can identify which events generated the entry of an individual by adding a characteristic to the population recording the event name/id when it occurs. The same holds for the other types of events. See e.g. vignette('IBMPopSim_human_pop') for an example with different death events.

5. **Swap event**

Finally, the C++ kernel code for a swap event specifies how the new characteristics of the individual are chosen. The kernel code can depend on the following variables:

<table>
<thead>
<tr>
<th>Variable</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>$I$</td>
<td>Current individual</td>
</tr>
<tr>
<td>$t$</td>
<td>Current time</td>
</tr>
<tr>
<td>$pop$</td>
<td>Current population (vector)</td>
</tr>
<tr>
<td>Model parameters</td>
<td>Depends on the model</td>
</tr>
</tbody>
</table>

We can now describe the different intensity classes.

2.3.2 Event creation with individual intensity

As explained in 1.1.2, events with intensities in the class individual are intensity functions $\lambda^I(I,t)$ which depend only on an individual’s age and characteristics, and on time. They are created using the function

```r
mk_event_individual(type = "TYPE", name ="name",
    intensity_code = "INTENSITY",
    kernel_code = "KERNEL_CODE")
```
The `intensity_code` argument is a character string containing few lines C++ code describing the intensity function. The available variables for the intensity code are given in the following Table.

**Table 5: C++ variables available for individual intensities**

<table>
<thead>
<tr>
<th>Variable</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>I</td>
<td>Current individual</td>
</tr>
<tr>
<td>t</td>
<td>Current time</td>
</tr>
<tr>
<td>Model parameters</td>
<td>Depends on the model</td>
</tr>
</tbody>
</table>

The intensity value has to be stored in a variable called `result`.

**Examples**

1. The intensity code below

   ```
   death_intensity <- "if (I.male)
       result = alpha_1*exp(beta_1*age(I, t));
   else
       result = alpha_2*exp(beta_2*age(I,t));"
   ```

   corresponds to a death intensity equal to \(d_1(a) = \alpha_1 \exp(\beta_1 a)\) for males and \(d_2(a) = \alpha_2 \exp(\beta_2 a)\) for females. In this case, the intensity function depends on the individuals’ age, gender, and on the model parameters \(\alpha = (\alpha_1, \alpha_2)\) and \(\beta = (\beta_1, \beta_2)\).

2. This example creates a death event with intensity depending on age and time, equal to

   \[ d(t,a) = 0.1 \exp(0.005a)1_{0 \leq t < 5} + 0.08 \exp(0.005a)1_{5 \leq t} \]

   This is done by creating the death function \(d\) using the predefined package functions `piecewise_xy` and `gompertz`. The function is then recorded as a model parameter and used in the argument `intensity_code` of `mk_event_individual`.

   ```
   time_dep_function <- piecewise_xy(c(5),
       list(gompertz(0.1,0.005),
              gompertz(0.08,0.005)))
   time_dep_function(0, 65) # death intensity at time 0 and age 65.
   ## [1] 0.1384031
   params <- list("death_function" = time_dep_function)
   death_event <- mk_event_individual(type = "death",
       intensity_code = "result=time_dep_intensity(t,age(I,t));")
   ```

**2.3.3 Event creation with interaction intensity**

Events with intensities in the class `interaction` are events which occur to an individual at a frequency which is the result of interactions with other members of the population (see 1.1.2), and which can be written as

\[
\lambda^e(I,t) = \sum_{J \in \text{pop}} U(I,J,t),
\]

where \(U(I,J,t)\) is the intensity of the interaction between individual \(I\) and \(J\).

An event with intensity in the class `interaction` is created by calling the function
The `interaction_code` argument contains few lines of C++ code describing the interaction function $U$. The interaction function value has to be stored in a variable called `result`. The available variables for the intensity code are given in the following Table:

<table>
<thead>
<tr>
<th>Variable</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>I</td>
<td>Current individual</td>
</tr>
<tr>
<td>J</td>
<td>Another individual in the population</td>
</tr>
<tr>
<td>t</td>
<td>Current time</td>
</tr>
<tr>
<td>Model parameters</td>
<td>Depends on the model</td>
</tr>
</tbody>
</table>

Example

```
dead_interaction_code <- " result = max(J.size - I.size, 0);"
```

In the example above, the death intensity of an individual $I$ is the result of competition between individuals, depending on a characteristic named `size`:

$$U(I, J, t) = (J.size - I.size)^+. $$

If $I$ meets randomly an individual $J$ of size bigger than $I.size$, he can die at the intensity $J.size - I.size$. The bigger is $I.size$, the lower is the death intensity of individual $I$.

The argument `interaction_type`, set by default at `random`, is an algorithm choice for simulating the model. When `interaction_type=full`, the intensity of an individual is computed making the sum over all the individuals in the population of the interactions with the given individual. In IBMPopSim, we introduced a “randomized” algorithm, in which the intensity of an individual is computed by picking randomly another individual in the population and computing its interaction function with the given individual. In most cases, the `random` algorithm is much faster than the `full` algorithm.

Note that events with individual intensities are also much faster to simulate since they only require to observe one individual to be computed.

### 2.3.4 Events creation with Poisson and Inhomogeneous Poisson intensity

When an event occur in the population with an intensity which does not depend on the population, the event intensity is of (inhomogeneous) Poisson class.

**Poisson intensity** When the event intensity is simply a constant, the class is called Poisson in reference to events occurring at jumps time of a Poisson process. Such events are created with the function

```
mk_event_poisson(type="TYPE",
                 name="NAME",
                 intensity="CONSTANT",
                 kernel_code = "KERNEL_CODE")
```
For instance,

```r
mk_event_poisson(type = "entry", intensity = "lambda",
    kernel_code = "double a_I= max(CNorm(20,2),0);
    newI.set_age(a_I,t);"
)
```

creates an event of type Entry, where individuals enter the population at a constant intensity \( \lambda \) (which has to be in the list of model parameters). When an individual \( \text{newI} \) enters the population at time \( t \), his age is chosen as a normally distributed random variable, with mean 20 and variance 4, using the function \( \text{CNorm()} \) (see Section 3.4).

**Inhomogeneous Poisson intensity** When the intensity depends on time (but not on the population), the event can be created similarly by using the function

```r
mk_event_inhomogeneous_poisson(type = "TYPE", name="NAME",
    intensity_code = "INTENSITY",
    kernel_code = "KERNEL_CODE"
)
```

For instance,

```r
mk_event_inhomogeneous_poisson(type = "entry",
    intensity = "result = lambda*(1+ cos(t));",
    kernel_code = "double a_I= CNorm(20,2);
    newI.set_age(a_I,t);"
)
```

creates the same event than before, but now individuals enter the population at the rate \( \lambda(1 + \cos(t)) \) depending on the current time \( t \).

### 2.4 Model creation

Finally, the IBM model is created using the function `mk_model`. The model is composed of:

- The characteristics names and types, which can be obtained from a population data frame with the function `get_characteristics`,
- The events list,
- The model parameters.

```r
model <- mk_model(characteristics = get_characteristics(pop),
    event = events_list,
    parameters = model_params)
```

During this step which can take a few seconds, the model is created and compiled using the Rcpp package. One of the advantages of the model structure in IBMPopSim is that the model depends only on the population characteristics and parameters names and types, rather than their values. This means that once the model has been created, various simulations can be done with different initial populations and parameters values. Thus, only one model has to be created to simulation a class of IBMs with varying parameters (see Section 2.5.2 for an example).

**Example** Here is an example of model with a population structured by age and gender, with birth and death events. The death intensity of an individual of age \( a \) is

\[
    d(a) = 0.008 \exp(0.02a),
\]

and females between 15 and 40 can give birth with birth intensity 0.05. The newborn is a male with probability \( p_{\text{male}} = 0.51 \).
params <- list("p_male"= 0.51,
    "birth_rate" = stepfun(c(15,40),c(0,0.05,0)),
    "death_rate" = gompertz(0.008,0.02))

death_event <- mk_event_individual(type = "death", name= "my_death_event",
    intensity_code = "result = death_rate(age(I,t));")

birth_event <- mk_event_individual( type = "birth",
    intensity_code = "if (I.male)
        result = 0;
    else
        result=birth_rate(age(I,t));",
    kernel_code = "newI.male = CUnif(0, 1) < p_male;"
)

model <- mk_model(characteristics = get_characteristics(pop),
    events = list(death_event,birth_event),
    parameters = params)

summary(model)
## Events:
## #1: individual event of type death
## #2: individual event of type birth
## ---------------------------------------
## Individual description:
## names: birth death male
## R types: double double logical
## C types: double double bool
## ---------------------------------------
## R parameters available in C++ code:
## names: p_male birth_rate death_rate
## R types: double closure closure
## C types: double function_x function_x

2.5 Simulation and outputs

Once the model has been created, the random evolution of the population can be simulated over a period of time $[0,T]$ by calling the function

popsim(model,population, events_bounds, parameters, age_max=Inf, time,...)

Where:

- **model** is the model created in the previous step,
- **population** is a population data frame representing the initial population,
- **parameters** is the list of parameters value,
- **age_max** is the maximum age of individuals in the population (set by default to Inf),
- **time** is the final simulation time $T$ or a vector of times (see @ref(simulation_swap)),
- **events_bounds** is a named vector of bounds for the intensity or interaction function of each event.
2.5.1 Events bounds

Since the IBM simulation algorithm is based on an acceptance-rejection method for simulating random times (see 1.1.3), the user has to specify bounds for the intensity (or interaction) function of each event before simulating a random path of the population evolution. Let us consider the model with a birth and death built in the previous section.

In the model example, the birth intensity of an individual of age $a$ is 0 if he is a male, and $b(a) = 0.0051_{[15,40]}$, if the individual is a female. Thus, the intensity bound for birth events is

$$\bar{\lambda}_b = \sup_{a \geq 0} \text{birth\_rate}(a) = 0.05$$

Since the death intensity function is not bounded, the user will have to specify a maximum age $a_{\text{max}}$ in popsim (all individuals above $a_{\text{max}}$ die automatically). Then, the bound for death events is

$$\bar{\lambda}_d = 0.008 \exp(0.02a_{\text{max}}).$$

The events bounds are defined as a named R vector, with components names corresponding to the events names. In our example, the death event has been named "my_death_event". No name has been specified for the birth event which thus has the default name "birth". Then,

```
a_max <- 120  # maximum age
events_bounds <- c("my_death_event" = 0.008*exp(0.02*a_max),
                  "birth" = max(params$birth_rate))
events_bounds
## my_death_event birth
## 0.08818541 0.05000000
```

Note that the ?max operator has been overloaded for some predefined functions of the package such as ?stepfun.

2.5.2 Simulation without swap events

Once the model and events bounds have been defined, a random trajectory of the population can be simulated by calling

```
sim_out <- popsim(model, pop, events_bounds, params,
                   age_max = a_max, time = 30)
sim_out
## Simulation on [0, 30]
str(sim_out)
```

sim_out is a list composed of

- A list arguments of the simulation inputs, including the initial population, parameters and event bounds.
- A named numeric vector logs of variables related to the simulation algorithm.
- The simulation output called population.
When there are no swap events (individuals don’t change of characteristics), the evolution of the population over the period \([0, 30]\) is recorded in a single data frame \(\text{sim\_out}\$\text{population}\). Each line of \(\text{sim\_out}\$\text{population}\) contains the information of an individual who lived in the population over the period \([0, 30]\). This includes individuals who were initially in \(\text{pop}\), as well as individuals who were born or entered the population.

```r
define str(sim_out$population)
## 'data.frame': 117963 obs. of 3 variables:
## $ birth: num -90 -89.9 -89.9 -89.9 -89.9 ... 
## $ death: num NA NA NA NA NA NA NA NA NA NA ... 
## $ male : logi TRUE TRUE FALSE FALSE TRUE FALSE ... 
```

The following Table describes the elements of the vector \(\text{sim\_out}\$\text{logs}\) containing information on simulation algorithm:

<table>
<thead>
<tr>
<th>Elements</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>proposed_events</td>
<td>Number of candidate event times proposed during the simulation</td>
</tr>
<tr>
<td>effective_events</td>
<td>Number of events which occurred during the simulation</td>
</tr>
<tr>
<td>cleanall_counter</td>
<td>Number of population cleans</td>
</tr>
<tr>
<td>duration_main_algorithm</td>
<td>Simulation time</td>
</tr>
</tbody>
</table>

For instance, the acceptance rate in the toy model is

```r
define 
#### effective_events 
#### 0.2177633
```

and the simulation time is

```r
define 
#### duration_main_algorithm 
#### 95215907
```

Parameters modification and event removal

As explained in Section 2.4 the structure of the compiled model allows the parameters’ values to be changed without recompiling the model. For instance, the parameter of the Gompertz death function can be modified to study the impact of an increase in mortality. Before running the simulation, the events bounds should be updated accordingly accordingly.

```r
define 
params$death_rate <- gompertz(0.01,0.02) # New death rate 

events_bounds[“my_death_event”] <- 0.01*exp(0.02*a_max) # Death event bound update 

new_sim_out <- popsim(model, pop, events_bounds, 
params, age_max = a_max, time = 30) # Population simulation
```

An event can also be remove by setting the event bound to 0:
```r
sim_out_no_birth <- popsim(model, pop,
  events_bounds = c("birth" = 0, "my_death_event" = 0.01*exp(0.02*a_max)),
  params, age_max = a_max, time = 30)
## [1] "event birth is deactivated"
## Simulation on [0, 30]
```

### 2.5.3 Outputs

The R data frame format for the output of the simulation allows all data manipulation functionalities of tidyverse or data.table packages to be used for analyzing the population. We also provide base functions to study the simulation outputs.

For instance, the population age pyramid can be computed at a given time (resp. at multiple dates) with the function `?age_pyramid` (resp. `?age_pyramids`). We refer to the other vignettes for more details on age pyramids computation and visualization.

```r
pop_out <- sim_out$population
# Population age-pyramid at time 30:
pyr <- age_pyramid(pop_out, ages = 0:a_max, time = 30)
plot_pyramid(pyr)
```

![Age Pyramid](image)

- **Group**
  - Females
  - Males

---

Number of individuals

**Age**

- 100 - 101
- 80 - 81
- 60 - 61
- 40 - 41
- 20 - 21
- 0 - 1

---

18
Mortality tables with compatibles with packages such as StMoMo can also be computed by with the functions ?death_table and ?exposure_table.

```r
female_pop <- pop_out[pop_out$male==FALSE, ]
Dxt <- death_table(female_pop, ages = 85:90, period = 20:30)  # Death table
Ext <- exposure_table(female_pop, ages = 85:90, period = 20:30)
```

### 2.5.4 Simple multithreading

If there are no interactions between individuals, i.e. if there are no events with intensity of class interaction, then the simulation can be parallelized easily by setting the optional parameter multithreading (FALSE by default) to TRUE:

```r
sim_out <- popsim(model, pop, events_bounds, params,
                   age_max = a_max, time = 30, multithreading = TRUE)
```

## duration_main_algorithm
## 18101526

By default, the number of threads is the number of concurrent threads supported by the available hardware implementation. The number of thread can be set manually with the optional argument num_threads of ?popsim.

### 2.6 Simulation with swap events

When there are swap events (individuals can change of characteristics), the dates of swap events and the changes of characteristics following each swap event should be recorded for each individual in the population, which is a memory intensive and computationally costly process. To maintain efficient simulations in the presence of swap events, we propose the following solution.

#### 2.6.1 Vector of times in ?popsim.

In the presence of swap events, the argument `time` of ?popsim should a vector of dates \((t_0, \ldots, t_n)\). In this case, popsim returns in the object `population` a list of \(n\) population data frames representing the population at time \(t_1, \ldots, t_n\), simulated from the initial time \(t_0\). For \(i = 1 \ldots n\), the \(i\)th data frame describes individuals who lived in the population during the period \([t_0, t_i]\), with their characteristics at time \(t_i\).

**Examples** A population 100 thousand individuals (or particles) is generated, all born at time 0 and divided into two subgroups. Individuals can swap from subgroup 1 (resp. 2) to subgroup 2 (resp. 1) at rate 0.1 (resp. 0.3).

```r
pop <- data.frame("birth" = rep(0,1e5), "death" = rep(NA,1e5),
                 "sub_grp" = sample(1:2, 1e5, replace = TRUE))

rates <- list( k12 = 0.1, k21=0.3)
#Only swap events occur in the population
swap_event <- mk_event_individual(type = "swap",
                                   intensity_code = "if (I.sub_grp == 1) result = k12;
                                                          else result = k21;",
                                   kernel_code = "I.sub_grp = 3 - I.sub_grp;")
```
Then, the population is simulated from $t_0 = 0$ to $t_n = 20$, and `popsim` returns a list of 20 data frames composed of the population at times $t = 1 \ldots 20$.

```r
time_vec <- 0:20
sim_out <- popsim(model = model_swap, population = pop,
                  events_bounds = c("swap"=max(unlist(rates))),
                  parameters = rates,
                  time = time_vec,
                  multithreading = TRUE)

pop_size <- nrow(pop)
# Mean number of individuals in subgroup 1 at each time:
p_1_t <- lapply(sim_out$population, function(pop_df){
    return(nrow(subset(pop_df, subgrp==1))/pop_size)
})
```

The model is an ergodic two states continuous time Markov chain with stationary distribution $(p_1, p_2) = (0, 75, 0.25)$. The figure below illustrates the convergence of the probability to be in subgroup 1 to $p_1$. 

![Graph showing the convergence of the probability to be in subgroup 1 to $p_1$.]
This example shows that IBMPopSIm can also be used to simulated continuous time Markov Chain with finite state space.

2.6.2 Individual life courses

It is possible to isolate the individuals' life course, by setting the optional argument with_id of `?mk_model` to `TRUE`. In this case, a new characteristic called `id` is automatically added to the population (if not already defined), identifying each individual with a unique integer.

```r
model_swap_id <- mk_model(characteristics = get_characteristics(pop),
                           events = list(swap_event),
                           parameters = rates,
                           with_id = TRUE)
## [1] "add 'id' as individual attributes"

sim_out_id <- popsim(model = model_swap_id,
                      population = pop,
                      parameters = rates,
                      events_bounds = c("swap"=0.3),
                      time = seq(0,5,by=1),
                      multithreading = TRUE)
## [1] "Add 'id' attributes to the population."
## Simulation on [0, 1] [1, 2] [2, 3] [3, 4] [4, 5]

head(sim_out_id$population[[1]])
## id birth death sub_grp
## 1  1    0 NA   1
## 2  9    0 NA   2
## 3 17    0 NA   1
## 4 25    0 NA   2
## 5 33    0 NA   2
## 6 41    0 NA   1
```

These data frames can be merged into a single data frame summarizing the life course of each individual, by calling the function `merge_pop_withid`.

```r
pop_list <- sim_out_id$population
pop_merge <- merge_pop_withid(pop_list, chars_tracked='sub_grp')
head(pop_merge)
## id birth death sub_grp_1 sub_grp_2 sub_grp_3 sub_grp_4 sub_grp_5
## 1  1    0 NA   1   1   1   1   1
## 2 32769  0 NA   1   1   1   1   1
## 3 65537  0 NA   1   1   1   1   1
## 4 98305  0 NA   1   1   1   1   1
## 5 31074  0 NA   1   1   1   1   1
## 6 63842  0 NA   1   1   1   1   1
```

Each line of `pop_merge` corresponds to the life course of one individual. However, the population is only represented at time $t = 0, 1, \ldots, 5$, and this data frame doesn’t account for exact swap event times or multiple swap events that occurred between between two time steps.
3 C++ essentials

The arguments intensity_code, interaction_code and kernel_code of the events functions must contain some C++ code given by the user. You don’t need to be a C++ guru to write the few instructions needed. These functions should use very little C++ syntax. They are essentially arithmetic or logical operations, tests and calls to predefined functions, which we give an overview below.

For code efficiency, you should not allocate memory in these functions (no new) or use type containers (std::vector, std::list, ...). If you think you need to allocate memory, consider as parameter an R vector that will be mapped via arma::vector (see below), or declare the variable as static. Also, it should not be necessary to make a loop. Keep in mind that these functions should be fast.

There are no C++ language restrictions so you can use all the functions of the standard C++11 library. However we detail in this section some functions that should be sufficient. For more details on C++ and Rcpp we recommend:

- C++ tutorial
- Rcpp and RcppArmadillo
- Rcpp for everyone

3.1 C++ syntax

- Each statement must be ended by a semicolon.
- Single-line comments start with two forward slashes //.
- To create a variable, you must specify the type and assign it a value (type variable = value;). Here some examples:

```cpp
int myNum = 5;     // Integer (whole number without decimals)
double myFloatNum = 5.99; // Floating point number (with decimals)
char myLetter = 'D';  // Character
bool myBoolean = true; // Boolean (true or false)
```

- bool data type can take the values true (1) or false (0).
- C++ supports the usual logical conditions from mathematics:
  - Less than: a < b
  - Less than or equal to: a <= b
  - Greater than: a > b
  - Greater than or equal to: a >= b
  - Equal to: a == b
  - Not Equal to: a != b
- The logical operators are: !, &&, ||
- The arithmetic operators are: +, -, *, /, %
- Compound assignment: +=, -=, *=, /=, %=
- Increment and decrement: ++, --
- Conditional ternary operator: ? : The conditional operator evaluates an expression, returning one value if that expression evaluates to true, and a different one if the expression evaluates as false. Its syntax is:

```cpp
condition ? result1; : result2;
```

- Use the if, else if, else statements to specify a block of C++ code to be executed if one or more conditions are or not true.
if (condition1) {
    // block of code to be executed if condition1 is true
} else if (condition2) {
    // block of code to be executed if the condition1 is false and condition2 is true
} else {
    // block of code to be executed if the condition1 is false and condition2 is false
}

• The syntax of the switch statement is a bit peculiar. Its purpose is to check for a value among a number of possible constant expressions. It is something similar to concatenating if-else statements, but limited to constant expressions. Its most typical syntax is:

switch (expression)
{
    case constant1:
        group-of-statements-1;
        break;
    case constant2:
        group-of-statements-2;
        break;
    ...
    default:
        default-group-of-statements
}

It works in the following way: switch evaluates expression and checks if it is equivalent to constant1; if it is, it executes group-of-statements-1 until it finds the break statement. When it finds this break statement, the program jumps to the end of the entire switch statement (the closing brace).

• When you know exactly how many times you want to loop through a block of code, use the for loop.

for (statement 1; statement 2; statement 3) {
    // code block to be executed
}

• The while loop loops through a block of code as long as a specified condition is true.

while (condition) {
    // code block to be executed
}

For more details we recommend a few pages of www.cplusplus.com about:

• Variables
• Operators
• Statements
3.2 Usual numeric functions

The most popular functions of the \texttt{cmath} library, which is included in the package, are the following:

- Exponential and logarithm functions: \texttt{exp}(x), \texttt{log}(x) (natural logarithm)
- Trigonometric functions: \texttt{cos}(x), \texttt{sin}(x), \texttt{tan}(x)
- Power functions: \texttt{pow}(x, a) meaning \(x^a\) and \texttt{sqrt}(x) meaning \(\sqrt{x}\)
- Absolute value: \texttt{abs}(x)
- Truncation functions: \texttt{ceil}(x) meaning \(\lceil x \rceil\) and \texttt{floor}(x) meaning \(\lfloor x \rfloor\)
- Bivariate functions: \texttt{max}(x, y) and \texttt{min}(x, y)

Note that these functions are not vectorial, the arguments \(x\) and \(y\) must be scalar. If the user wants to call some other functions of \texttt{cmath} not listed in the table, this is possible by adding the prefix \texttt{std::} to the name of the function (scope resolution operator :: to access to functions declared in namespace standard \texttt{std}).

3.3 Individuals characteristics and model parameters: link between R and C++

To facilitate the model creation, the individuals’ characteristics and a list model parameters can be declared in the R environment and used in the C++ code.

The data shared between the R environment and the C++ code are:

- The characteristics of the individuals which must be atomic (Boolean, scalar or character).
- The model parameters: a list of variables of type:
  - Atomic, vector or matrix.
  - Predefined real functions of one variable, or list of such functions.
  - Piecewise real function of two variables, of list of such.

3.3.1 Atomic types (characteristics and parameters)

Here is the conversion table used between the atomic types of R and C++.

<table>
<thead>
<tr>
<th>C++ type</th>
<th>R type</th>
</tr>
</thead>
<tbody>
<tr>
<td>bool</td>
<td>logical</td>
</tr>
<tr>
<td>int</td>
<td>integer</td>
</tr>
<tr>
<td>double</td>
<td>double</td>
</tr>
<tr>
<td>char</td>
<td>character</td>
</tr>
</tbody>
</table>

Individuals characteristics

The characteristics of an individual are defined by a named character vector containing the name of the characteristic and the associated C++ type. The function \texttt{get_characteristics} provides a way to extract the characteristics from a population data frame.

```r
library(IBMPopSim)
pop <- IBMPopSim::EW_popIMD_14$sample
get_characteristics(pop)
#>  male   IMD
#> "bool" "int"
```
The requires **birth** and **death** characteristics are of type **double**.

### Atomic model parameters

The model parameters are given by a named list of R objects. We recall that the type of an R object can be determined by calling the ?typeof function. Atomic objects declared as model parameters can be directly used in the C++ code.

In the example below, the variable `code` contains simple C++ instructions depending on the parameters defined in `params`. The `summary` of the model `mod` gives useful information on the types of characteristics and parameters used in C++ code.

```r
params <- list("lambda" = 0.02, "alpha" = 0.4, "mu" = as.integer(2))

code <- "result = lambda + alpha * (age(I, t) + mu);"

event_birth <- mk_event_individual("birth", intensity_code = code)

mod <- mk_model(get_characteristics(pop), events = list("birth" = event_birth),

parameters = params, with_compilation = FALSE)

summary(mod)
```

### 3.3.2 Vectors and matrices (model parameters)

Two of the possible model parameters types given in the argument parameters of the ?mk_model function are R vectors and R matrices. We call R vector a numeric of length at least 2. These types are converted, using the RcppArmadillo library, in C++ Armadillo types, arma::vec and arma::matrix respectively.

The classes arma::vec and arma::matrix are rich and easy-to-use implementations of one-dimensional and two-dimensional arrays. To access to individual elements of an array, use the operator () (or [] in dimension 1).

- (n) or [n] for arma::vec, access the n-th element.
- (i,j) for arma::matrix, access the element stored at the i-th row and j-th column.

**Warning:** The first element of the array is indexed by subscript of 0 (in each dimension).

Another standard way (in C++) to access elements is to used iterators. An iterator is an object that, pointing to some element in a range of elements, has the ability to iterate through the elements of that range using a set of operators (see more details on iterators).

Let v be a an object of type arma::vec and A be a an object of type arma::matrix. Here we show how to get the begin and the end iterators of these objects.

- v.begin(): iterator pointing to the begin of v
- v.end(): iterator pointing to the end of v
- A.begin_row(i): iterator pointing to the first element of row i
3.3.3 Predefined functions (model parameters)

To facilitate the implementation of intensity functions and kernel code, R functions have been predefined in IBMPopSim, which can be defined as model parameters and then called in the C++ code. The goal is to make their use as transparent as possible.

3.3.3.1 Real functions of one variable

Here is a list of such functions that can be defined as a R object and called from R and C++:

- `stepfun`: Step function.
- `linfun`: Linear interpolation function.
- `gompertz`: Gompertz–Makeham intensity function.
- `weibull`: Weibull density function.
- `piecewise_x`: Piecewise real function.

See the reference manual for mathematical definitions of these functions (?stepfun).
Once the model is created, these predefined functions are transformed into C++ functions, identified as `function_x`.

We illustrate below some examples of the use of these functions:

1. We define `dr` a `stepfun` depending on some values in `EW_pop_14$rates`. Note that this function applies to an age `a`.

```r
dr <- with(EW_pop_14$rates, 
           stepfun(x = death_male[, "age"], y = c(0, death_male[, "value"])))
plot(dr, xlab="age", ylab="dr", main="Example of step function")
```
We can also initialize a piecewise real function $f$ defined by the function $dr$ before age 80, and by a Gompertz-Makeham intensity function for ages after 80.

```r
f <- piecewise_x(80, list(dr, gompertz(0.00006, 0.085)))
x <- seq(40, 110)
plot(x, sapply(x, f), xlab="age", ylab="f", main="Example of piecewise function")
```
Example of piecewise function

2. Once a function has been created and defined as model parameter, a model depending on the function can be defined. For example we use in the model below the `stepfun dr` as a model parameter named `rate`. The function is used to defined the intensity of death events in the model.

   ```r
   params <- list("rate" = dr)
   event <- mk_event_individual("death", intensity_code = "result = rate(age(I, t));")
   mod <- mk_model(get_characteristics(pop), events = list("death" = event),
                  parameters = params, with_compilation = FALSE)
   summary(mod)
   # Events:
   # #1: individual event of type death
   # # Individual description:
   # # names: birth death male IMD
   # # R types: double double logical integer
   # # C types: double double bool int
   # # R parameters available in C++ code:
   # # names: rate
   # # R types: closure
   # # C types: function_x
   
   3. After compilation, the parameter `rate` can actually be replaced by any function of type `function_x`. For example you can call

   ```r
   ```
popsim(mod, pop, params = list("rate" = dr), age_max = 120,
    events_bounds = c("death" = dr(age_max)), time = 10)

as well as

popsim(mod, pop, params = list("rate" = f), age_max = 120,
    events_bounds = c("death" = f(age_max)), time = 10)

3.3.3.2 Piecewise real functions of two variables
In the C++ code these R functions declared with
?piecewise_xy are identified as function_xy functions. See ?piecewise_xy for mathematical definition.
This function allows to easily define a step function that depend on age and time.

3.3.3.3 List of functions
As parameter you can use a list of functions. All R functions in the list must
be of the same C++ type: either function_x or function_xy. In C++ code the list of functions is replaced
by a std::vector of function_x or function_xy (with first element indexed by 0).

3.4 Random variables

We use the following notations to describe the available C++ random distributions, which can be used in
the C++ intensity and kernel codes.

- \( U(a, b) \) : Uniform distribution on \([a, b]\) with \( a < b \)
- \( \mathcal{E}(\lambda) \) : Exponential distribution, \( \lambda > 0 \)
- \( \mathcal{N}(\mu, \sigma) \) : Gaussian distribution, \( \mu, \sigma \in \mathbb{R} \)
- Pois(\( \lambda \)) : Poisson distribution, \( \lambda > 0 \)
- \( \Gamma(\alpha, \beta) \) : Gamma distribution, \( \alpha, \beta > 0 \)
- Weib(\( a, b \)) : Weibull distribution, \( a > 0, b > 0 \)
- \( \mathcal{U}\{a, b\} \) : Discrete uniform distribution on \(\{a, a + 1, \ldots, b\}\) with \( a < b \)
- \( \mathcal{B}(p) \) : Bernoulli distribution, the probability of success is \( p \in (0,1) \)
- \( \mathcal{B}(n, p) \) : Binomial distribution \( n \geq 1, p \in (0, 1) \)
- \( \mathcal{D}_n \) : Discrete distribution with values in \(\{0, \ldots, n - 1\}\) and with probabilities \(\{p_0, \ldots, p_{n-1}\}\).

In the table below we show how to call them, which means how to make independent realizations of these
random variables, and we give the reference to the C++ corresponding function of the random library hidden
in this call.

<table>
<thead>
<tr>
<th>Function call</th>
<th>Meaning</th>
<th>C++ random internal function</th>
</tr>
</thead>
<tbody>
<tr>
<td>CUnif(a = 0, b = 1)</td>
<td>( \mathcal{U}(a, b) )</td>
<td>uniform_real_distribution&lt;double&gt;</td>
</tr>
<tr>
<td>CExp(( \lambda = 1 ))</td>
<td>( \mathcal{E}(\lambda) )</td>
<td>exponential_distribution&lt;double&gt;</td>
</tr>
<tr>
<td>CNorm(( \mu = 0, \sigma = 1 ))</td>
<td>( \mathcal{N}(\mu, \sigma) )</td>
<td>normal_distribution&lt;double&gt;</td>
</tr>
<tr>
<td>CPoisson(( \lambda = 1 ))</td>
<td>Pois(( \lambda ))</td>
<td>poisson_distribution&lt;unsigned&gt;</td>
</tr>
<tr>
<td>CGamma(( \alpha = 1, \beta = 1 ))</td>
<td>( \Gamma(\alpha, \beta) )</td>
<td>gamma_distribution&lt;double&gt;</td>
</tr>
<tr>
<td>CWeibull(( a = 1, b = 1 ))</td>
<td>Weib(( a, b ))</td>
<td>weibull_distribution&lt;double&gt;</td>
</tr>
<tr>
<td>CUnifInt(( a = 0, b = 2^{31} - 1 ))</td>
<td>( \mathcal{U}{a, b} )</td>
<td>uniform_int_distribution&lt;int&gt;</td>
</tr>
<tr>
<td>Function call</td>
<td>Meaning</td>
<td>C++ random internal function</td>
</tr>
<tr>
<td>-----------------------</td>
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<td>--------------------------------------------</td>
</tr>
<tr>
<td>CBern($p = 0.5$)</td>
<td>$B(p)$</td>
<td>bernoulli_distribution</td>
</tr>
<tr>
<td>CBinom($n = 1, p = 0.5$)</td>
<td>$B(n, p)$</td>
<td>binomial_distribution&lt;int&gt;</td>
</tr>
<tr>
<td>CDiscrete(p_begin, p_end)</td>
<td>$D_n$</td>
<td>discrete_distribution&lt;int&gt;</td>
</tr>
</tbody>
</table>

In the discrete distribution call `CDiscrete(p_begin, p_end)`, the arguments `p_begin` and `p_end` represent the iterators to the begin and to the end of an array which contains $\{p_0, \ldots, p_{n-1}\}$. Note that the use of iterators is a convenient and fast way to access a column or row of a matrix `arma::mat`.

## 4 References

