**First Steps to R Packages** 

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# Workshop - SSC Student Conference

May 28, 2022



We will go through the basic steps of constructing a package, via an example where 4 functions and 1 data set are to be combined into a single package and prepared for submission to CRAN.

The functions and data set will relate to the two-component mixture distribution having density function

$$f(x) = pf_1(x) + (1-p)f_2(x)$$

where  $p \in (0, 1)$ ,  $f_1(x)$  and  $f_2(x)$  are pdfs of non-central t random variables.

The t random variables are governed by a non-centrality parameter ncp which locates their center and by the number of degrees of freedom df.



The functions that will make up the package are designed to

- simulate the random variables (rtmix())
- calculate quantiles (qtmix())
- calculate probability densities (dtmix()) and
- calculate cumulative probabilities (ptmix())

### The random variate generator: rtmix()

**rtmix**(5, df = c(4, 7), ncp = c(-1, 5), PI = .24)

## [1] -0.7637152 3.8608394 5.7320305 7.2167542 5.8635<mark>216</mark>

5 randomly generated values from the mixture distribution

- with 4 and 7 degrees of freedom
- with component means of -1 and 5
- where the first component has probability 0.24, i.e. p = 0.24

#### The functions

#### The quantile function: qtmix()

qtmix(c(.25, .75, .9), df = c(4, 7), ncp = c(-1, 5),
PI = .24)

## [1] 2.453922 6.096464 7.806849

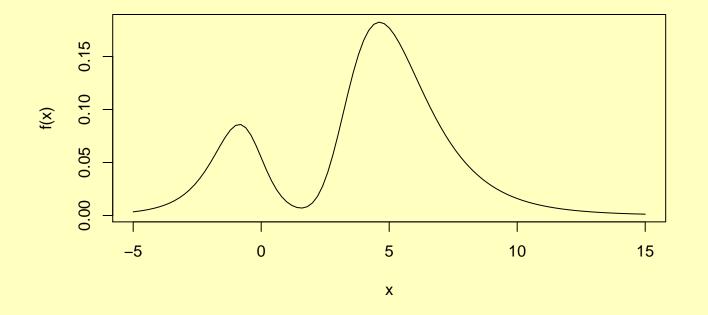
## The 25th, 75th and 90th percentiles of the mixture distribution.

# **The functions**



### The probability density function calculator: dtmix()

curve(dtmix(x, df = c(4, 7), ncp = c(-1, 5), PI = .24),
from = -5, to = 15, ylab = "f(x)")



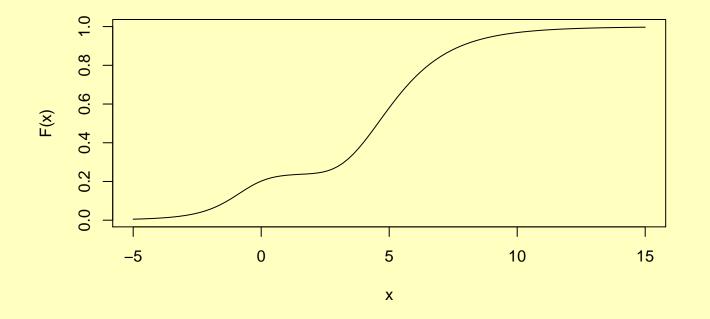
## A plot of the probability density curve of the mixture distribution.

### **The functions**



### The cumulative distribution function calculator: ptmix()

curve(ptmix(x, df = c(4, 7), ncp = c(-1, 5), PI = .24),
from = -5, to = 15, ylab = "F(x)")



## A plot of the cdf curve for the mixture distribution.



### The data file is called simdata:

str(simdata)

## 'data.frame': 200 obs. of 2 variables: ## \$ x: num -0.043 -1.211 -1.079 -0.383 -1.381 ... ## \$ y: num -0.748 -2.713 -2.518 -3.78 -2.62 ...



Since the functions concern a 2-component mixture of trandom variables, we name the package *tmix*.

Our first step is to create a package directory which has that name.

File tolic View Search Terminal Help
johbraun@braun7:~\$ mkdir tmix
johbraun@braun7:~\$ \_\_\_\_\_\_



Within the *tmix* directory, we require the following:

- A *man* directory. We need at least one, usually both, of the following
  - An *R* directory.
  - A data directory.

johbraun@braun7:~\$ mkdir tmix johbraun@braun7:~\$ cd tmix johbraun@braun7:~/tmix\$ mkdir R johbraun@braun7:~/tmix\$ mkdir data johbraun@braun7:~/tmix\$ mkdir man johbraun@braun7:~/tmix\$



Within the *tmix* directory, we also minimally require the following text files:

- A DESCRIPTION file.
- A NAMESPACE file.



You can also construct the package directories from within R or RStudio using the package.skeleton function. This function is built into the default R distribution.

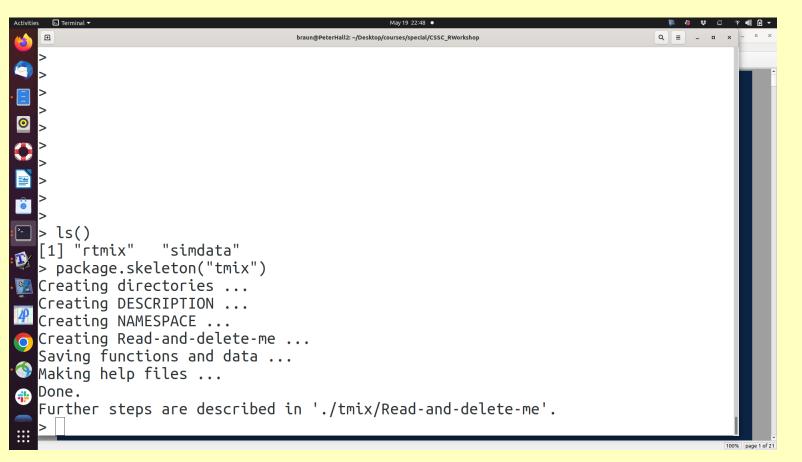
You need at least one function or data object in your workspace in order to use this function.

Start with only one or two objects and complete the build and check process on a package that only contains them. Then gradually add functions and data to your package, rebuilding and rechecking at each stage. This will make debugging easier and faster.



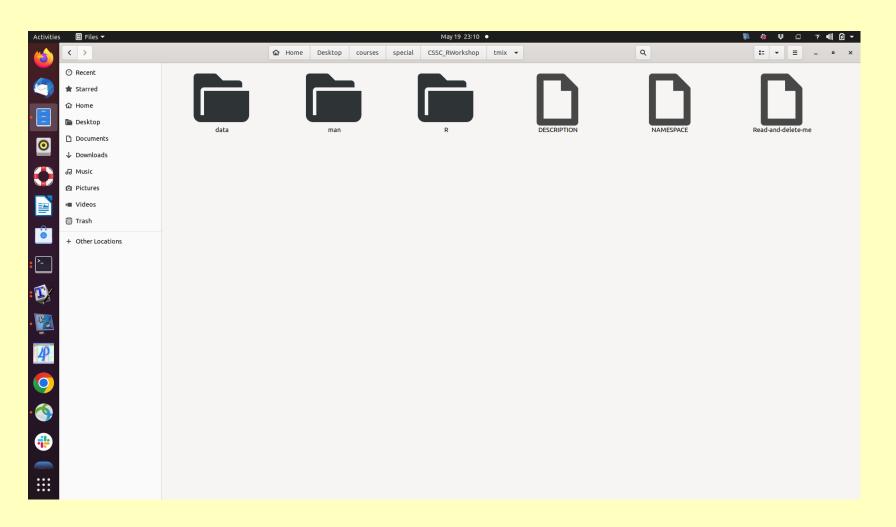
# Here, I am starting with one function and one data object in the workspace.

uble of mine

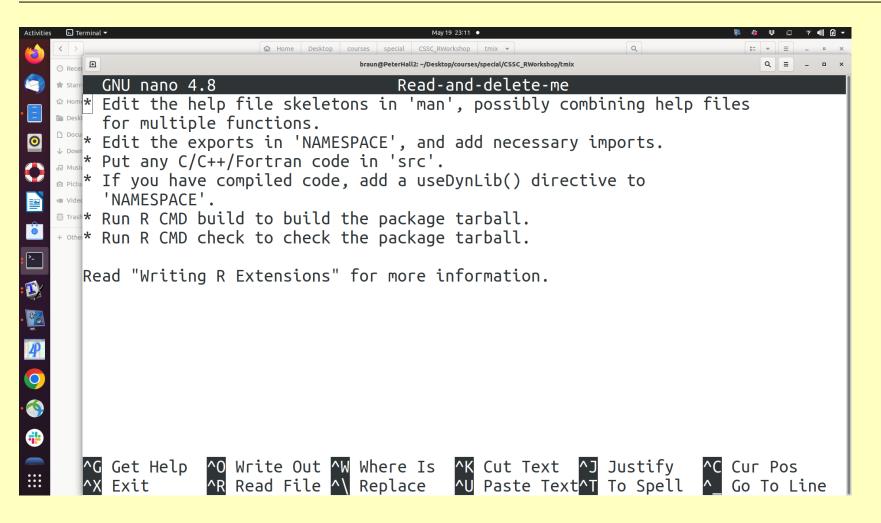


The command package.skeleton("tmix") creates the directory structure with the beginnings of the files but just for the two objects. I need to add in the other three objects, one at a time, later.\*

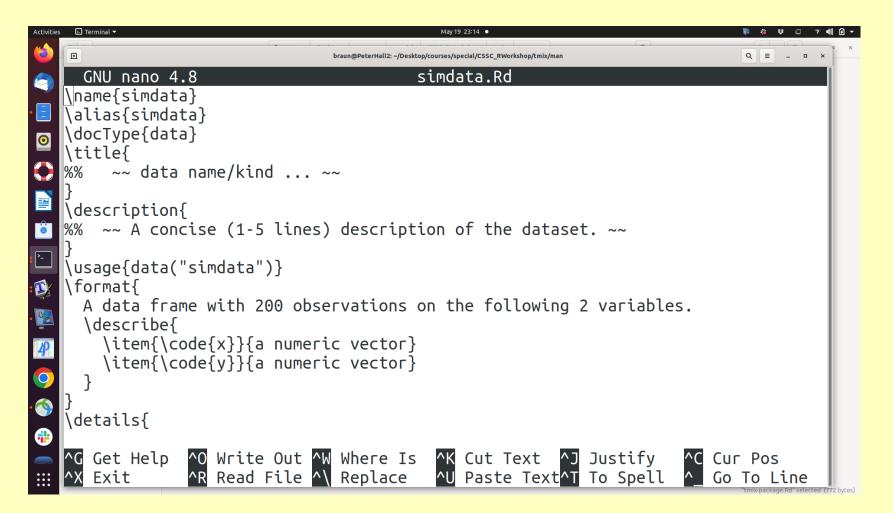
\*Watch the award-winning 19-second silent movie packageSkeleton.mp4 to see the commands in action.



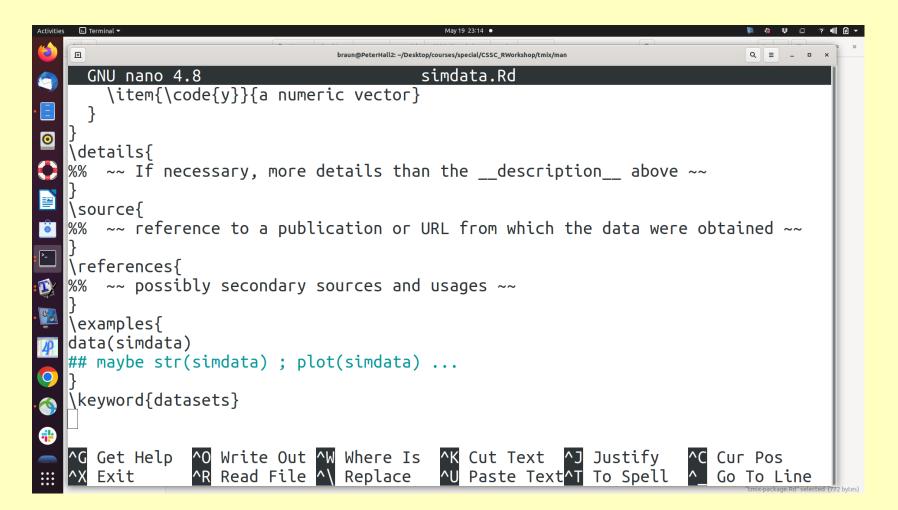
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The third and fourth \* points can be ignored unless you are doing something more sophisticated. Writing R Extensions can be found on the CRAN site. You do need to consult this.



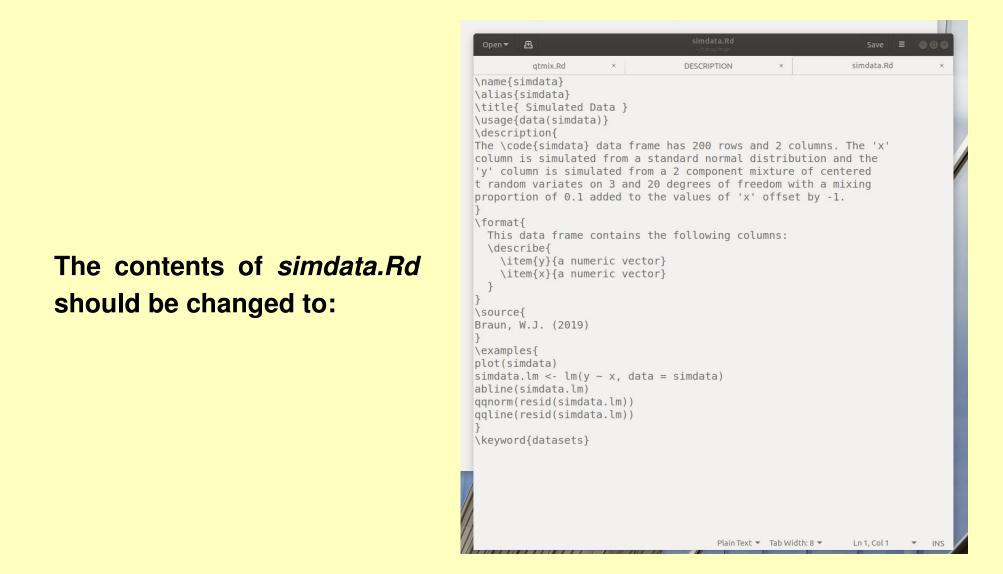
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# **Cleaning up the man file - automatic method**



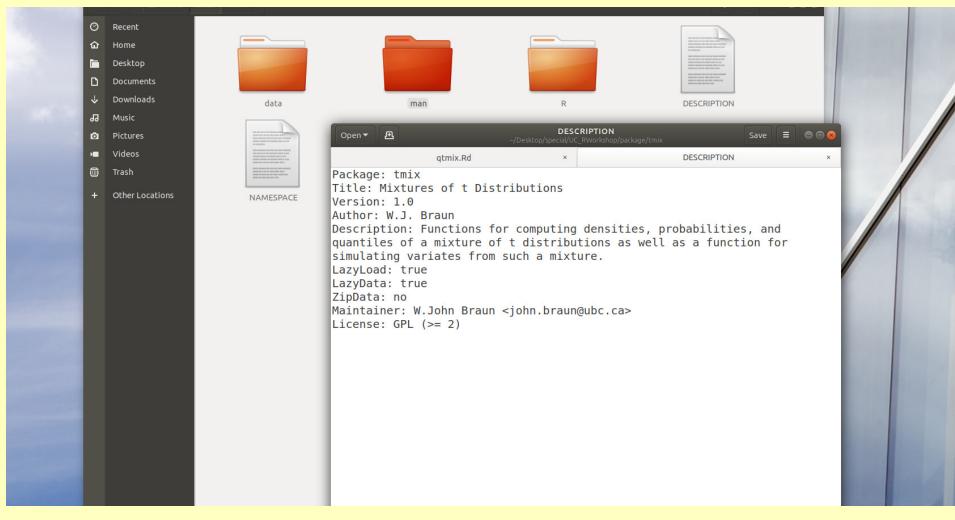


But we're getting ahead of ourselves.

# **The DESCRIPTION file**



### **Possible contents for the DESCRIPTION are as follows:**



The contents of the Description should all be on one line. Multiple lines are used here for display purposes only.



This file is needed so that you don't have to worry that the names you choose for your functions or data sets will conflict with names of functions and data sets from other packages.

The contents of NAMESPACE for this example:

```
importFrom("stats", "qt", "dt", "pt", "rt", "rbinom")
exportPattern(".")
export("dtmix", "ptmix", "qtmix", "rtmix")
```

We have not included our data file in the export list, only the functions.

We have included all 4 of our functions in the export list.

If we had held one back, the package user would not have direct access to that function.



### The R and data directories

# We create four files in the R subdirectory, calling them

- qtmix.R
- rtmix.R
- dtmix.R
- ptmix.R

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The contents of these files are the corresponding functions.

The data, stored in a file named *simdata.R* will be copied into the *data* subdirectory.



The man directory will contain 5 files:

- qtmix.Rd
- rtmix.Rd
- dtmix.Rd
- ptmix.Rd
- simdata.Rd

Here, we will construct those files manually. Optionally, you can include a help file for the package itself called *tmix-package.Rd*. This gives an overview of the package.

You can avoid some of this work by using the *roxygen2* package (Wickham et al, 2019).

# The help directory: man



| The contents of <i>qtmix.Rd</i> are: | <pre>Que defined<br/>//renderly control of the control</pre> |
|--------------------------------------|---|
|                                      | Plain Text ▼ Tab Width: 8 ▼ Ln 9, Col 12 ▼ INS  |



Most of the contents of qtmix.Rd are required, with the exception of the following: details, seealso, references, author and examples.

These options are strongly recommended, since the more detail that is provided, the better.

For a list of keywords that you might use, try

RShowDoc("KEYWORDS")



The qtmix() function invoked a Newton-Raphson iteration to solve the required inverse problem.

Since this iteration requires a loop, it may have been better to write the code into a *C* or *Fortran* program that would be called from R.

If this had been done, the external code would be written in the form of a subroutine or collection of subroutines and stored in an additional directory called *src*.

Special commands are required within the R functions (e.g. qtmix()) which would need to access the external code. The R manual has more information on this.

Documentation for the package itself is also recommended, as is the creation of vignettes, which describe the package in action.



Once the pieces of the package are assembled, it is necessary to build the package. In RStudio, this is fairly straightforward.\*

At the command line (Mac, Linux or Cygwin in Windows), we would type

R CMD build tmix

To check that the package components are properly constructed, we next type

R CMD check tmix

# It is usually advisable to build and check the package as it is being constructed, instead of waiting until all the pieces have been assembled.

\*See the 99 second silent movie called RStudiobuildCheck.mp4 for checking and building a copy of the package called *tmix2*.



Before submitting the package to CRAN, we need to do a more thorough check as follows:

R CMD check --as-cran tmix

This last check is to be done using the most recent development version of R. Before submitting to CRAN, the submittor should check the R manual to ensure that all requirements have been met.

The final steps in the submission can be carried out at https://cran.r-project.org/submit.html.

Alternatively, and maybe preferably, you could post your package on a github site to encourage collaborative development of it.