# Step by Step Tutorial to creating R Packages

Heng Wang
Michigan State University

#### Introduction

- R is an open source statistical software
- R provides functions to perform statistical operations
- Classical (regression, logistic regression, ANOVA, etc)
- Modern (neural networks, bootstrap, genomic selection, etc)
- Can be easily extended by make new packages

- To install an R package, use function install.packages()
- For example, to fit a mixed model, we could use function mixed.solve(). It requires package "rrBLUP".

#### Steps to Build an R package

- Step 1. Prepare your functions
- Step 2. Build the structure of the package using package.skeleton()
- Step 3. Edit **DESCRIPTION** File
- Step 4. Edit the help File
- Step 5. Preparation for Windows Users (RTools)
- Step 6. Build and install the R package
- Step 7. Check the R package
- Step 8. Add functions and data to a package

- -- Step 1. Prepare your functions
- Before you write your functions, clear the working space using rm(list=ls()).
- Write your function. Load all the data you want to include in the package.
- Set working directory to the position containing the .R file.

## Build an R Package -- Step 2. package.skeleton()

- Run package.skeleton(name, list).
- For example: package.skeleton(name="cum", list=c("my.cumsumprod", "xvec.example", "output.example")
- Or, package.skeleton(name="cum", code\_files="cumsumprod.R")
- A new folder cum is built. If just run
  package.skeleton(), then anRpackage will be
  built.

#### Step 2 (Cont.)

- Inside cum / anRpackage you many find several folders:
- R: contains R code files
- o data: contains data files
- man: contains documentation files (.Rd)
- You may also have src folder, if your function contains C, C++, or FORTRAN source.
- Other files: tests, exec, inst, etc.

#### Step 2 (Cont.)

- ... also some files.
- Read-and-delete-me: contain instructions for following steps.
  - \* Edit the help file skeletons in 'man', possibly combining help files for multiple functions.
  - \* Edit the exports in 'NAMESPACE', and add necessary imports.
  - \* Put any C/C++/Fortran code in 'src'.
  - \* If you have compiled code, add a useDynLib() directive to 'NAMESPACE'.
  - \* Run R CMD build to build the package tarball.
  - \* Run R CMD check to check the package tarball.
  - Read "Writing R Extensions" for more information.
- DESCRIPTION: manual file for the users.
- NAMESPACE

#### -- Step 3. Edit **DESCRIPTION** File

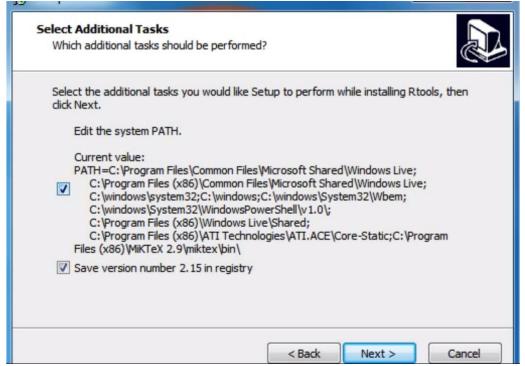
- Package: cum
  - -- name of the package
- Type: Package
- Title: What the package does (short line)
  - -- contains no more than 65 characters
- Version: 1.0
  - -- a sequence of non-negative integers, like: 1.0.2, 1-0-2
- Date: 2013-02-27
  - -- Date that the package was created. Today's date by default
- Author: Who wrote it
  - -- all the authors, no limit
- Maintainer: Who to complain to <u>yourfault@somewhere.net</u>
  - -- one name and an email address
- Description: More about what it does (maybe more than one line)
  - -- Description of the package, no length limit
- License: What license is it under?
  - -- Usually GPL-2 (GNU General Public License Version 2), which is good for CRAN / Bioconductor. Check "Writing R Extensions" for all license abbreviations.

## Build an R Package -- Step 4. Edit the help File

- Fill the content in each category
- Delete the comments or instructions. Change the default content
- Do this for each .Rd file in man folder.

- -- Step 5. Preparation for Windows Users
- Download and install Rtools. <a href="http://cran.r-project.org/bin/windows/Rtools/">http://cran.r-project.org/bin/windows/Rtools/</a>
- Attention! Check the checkbox to update the

current PATH.



#### Step 5 (Cont.)

- Change the PATH in Control Panel.
- Click System, then Advanced system settings.
- Click the Advanced tap in the prompt window. Then click the Environment Variables.
- In **PATH**, click **Edit**...
- C:\Windows\System64\;c:\Rtools\bin;c:\Rtools\gcc-4.6.3\bin;C:\Program Files\R\R-
  - 2.15.1\bin\x64;c:\Rtools\perl\bin;c:\Rtools\MinGW\bin;c:\R\bin;c:\Rtools\MinGW;c:\Perl\bin;c:\Program Files\MiKTex 2.6\miktex\bin;C:\Program Files (x86)\SSH Communications Security\SSH Secure Shell

# Build an R Package -- Step 6. Build and install the R package

- In search box, type command prompt
- In command prompt, change directory to the place that contains the R package
- Build R package using R CMD build pkgName.
   For example I use R CMD build cum. A tar.gz file is built under the working directory.

#### Step 6 (Cont.)

- Install the R package using R CMD INSTALL pkgName. Here I use R CMD INSTALL cum\_1.0.tar.gz.
- If any error occurs, check the .Rd file. Then rerun R CMD build, R CMD INSTALL.

## Build an R Package -- Step 7. Check the R package

- Install Miktex / (Mactex) package inconsolata using mpm --verbose --install inconsolata.
- Check the R package using R CMD check pkgName. I use R CMD check cum.
- In **R** environment, type **library(pkgName)**. For example, **library(cum)**.
- You can type

?cum

?my.cumsumprod

?xvec.example

?output.example

## -- Step 8. Add functions and data to a package

- Change the working directory to the folder that contains your new functions and/or data.
- Copy the functions into working space.
- Run prompt() to the new function, i.e., prompt(cumadd). Now you have a help file for cumadd.
- Edit the .Rd help file.
- Move the .R file and the .Rd file to the package folder. Put the .R file in the R folder. Put the .Rd file in the man folder.

#### Step 8 (Cont.)

- Read the data file into the working space.
- Save the data as an .rda file.
- Create the help file using prompt() function.
- Edit the .Rd help file.
- Move the .rda file and the .Rd file to the package folder. The .rda file goes to the data folder. The .Rd file goes to the man folder.

#### Step 8 (Cont.)

- Build and install the package again.
- R CMD build cum
- R CMD INSTALL cum\_1.0.tar.gz
- R CMD check cum
- In R console, type library(cum)
  - ?cum
  - ?my.cumsumprod
  - ?xvec.example
  - ?output.example
  - ?yvec
  - ?cumadd

### Questions?

## Thank you!