### R Programming

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#### Intro

Understanding the basics of R programming helps to improve analysis/reporting scripts and extend what we can do with R.

Good coding practice follows the DRY principle: Don't Repeat Yourself. Rather than modifying copy-pasted code chunks, we might

- write a custom function
- use loops or iteration functions to perform multiple similar tasks

Custom functions can be used to provide convenient wrappers to complex code chunks as well as implement novel functionality.

#### **Data Structures Revisited**

For basic data analysis, our data is usually imported and we use high-level functions (e.g. from **dplyr**) to handle it.

For programming, we need to work with lower-level data structures and be able to

- create basic objects
- extract components
- coerce one data type to another

#### **Vectors**

numeric(), character() and logical() can be used to initialize vectors of the corresponding type for a given length

```
x <- numeric(3)
x
# [1] 0 0 0</pre>
```

Elements can be assigned by indexing the positions to be filled, e.g.

```
x[1] \leftarrow 4
x[-c(2, 3)] \leftarrow 4
```

This is particularly useful when programming an iterative procedure.

as.logical(), as.numeric() and as.character() coerce to the corresponding type, producing NAs if coercion fails.

### **Logical Vectors**

Logical vectors are commonly used when indexing. The vector might be produced by a logical operator:

```
x <- c(1, 1, 2, 2)
x > 1
# [1] FALSE FALSE TRUE TRUE
x[x > 1]
# [1] 2 2
```

duplicated() is also useful here:

```
duplicated(x)
# [1] FALSE TRUE FALSE TRUE
!duplicated(x)
# [1] TRUE FALSE TRUE FALSE
```

Single element logical vectors (scalars) are useful for flow control, see later.

#### **Numeric Vectors**

The are several convenience function for creating numeric vectors, notably seq() and rep().

As they are so useful there are fast shortcuts for particular cases

```
seq_len(4)
# [1] 1 2 3 4
x <- 3:5
seq_along(x)
# [1] 1 2 3
rep.int(1:2, times = c(2, 3))
# [1] 1 1 2 2 2</pre>
```

### **Character Vectors**

Character vectors may be used for creating names

```
names(x) <- paste0(LETTERS[1:3], 1229:1231)
x
# A1229 B1230 C1231
# 3 4 5
names(x)
# [1] "A1229" "B1230" "C1231"</pre>
```

Names can be used as an alternative to numeric or logical vectors when indexing

```
x["B1230"]
# B1230
# 4
```

### **Matrices**

A matrix is in fact also a vector, with an attribute giving the dimensions of the matrix

```
M <- matrix(1:6, 2, 3)
M
# [,1] [,2] [,3]
# [1,] 1 3 5
# [2,] 2 4 6
str(M)
# int [1:2, 1:3] 1 2 3 4 5 6
attributes(M)
# $dim
# [1] 2 3</pre>
```

Useful functions for matrices include dim(), ncol(), nrow(), colnames() and rownames(). rbind() and cbind() can be used to row-bind or column-bind vectors.

Matrices enable computation via matrix algebra as well as row/column-wise operations.

### Lists

Lists collect together items which may be different types or lengths. Like a vector, elements may be named.

```
results <- list(matrix = M, vector = x)
results

# $\text{matrix}

# [,1] [,2] [,3]

# [1,] 1 3 5

# [2,] 2 4 6

# $\text{vector}

# A1229 B1230 C1231

# 3 4 5
```

Lists are often used to return the results of a function.

Elements can be indexed by [ to return a list or [ [ to return a single element. \$ can be used to extract elements by name, e.g. results\$matrix.

#### **Data Frames**

Data frames are lists of variables of the same length and hence can often be treated as a matrix

```
dat <- data.frame(x = x, id = letters[1:3])
dat[[1]]
# [1] 3 4 5
dat[1, 2]
# [1] a
# Levels: a b c</pre>
```

The tibble() (or as\_data\_frame()) from the tibble package (or dplyr) can be used to produce a tibble instead.

When the data are all numeric, as .matrix() can be used to coerce to a matrix.

#### **Your Turn**

The exercises. Rmd file provides a template for the exercises.

The lm function calls the "workhorse" function lm.fit to actually fit the model. Unlike lm, which works from a formula, lm.fit works from the model matrix and the response vector.

Define a response y containing 10 numeric values. Define an explanatory variable z of the numbers 1 to 10.

Use the function cbind() to create a matrix x with 1s in the first column and z in the second column.

Fit a model using fit1 <- lm.fit(x, y). Use str to explore the structure of the results. Use \$ to extract the coefficients.

#### **Control Structures**

Control structures are the commands that make decisions or execute loops.

Conditional execution: if/else, switch

Loops: for, while, repeat

### if/else

An if statement can stand alone or be combined with an else statement

```
x <- 1:3
if (all(x > 0)) {
    res <- mean(x)
} else {
    res <- mean(abs(x))
}</pre>
```

The condition must evaluate to logical vector of length one. The functions all(), any(), is.null() and other is. functions are useful here.

# **Conditioning on equality**

Using == may not be appropriate as it compares each element; identical() will test the whole object

```
x <- y <- 1:2
x == y
# [1] TRUE TRUE
identical(x, y)
# [1] TRUE</pre>
```

all.equal() will alow for some numerical "fuzz"

```
z <- sqrt(2)
identical(z * z, 2)
# [1] FALSE
all.equal(z * z, 2)
# [1] TRUE</pre>
```

#### switch

The switch() function provides a more readable alternative to nested if statements

The final unnamed argument is the default.

#### for

A for loop repeats a chunk of code, iterating along the values of a vector or list

```
x <- c("banana", "apple")
for (nm in x) print(nm)
# [1] "banana"
# [1] "apple"</pre>
```

Unassigned objects are not automatically printed; hence call to print(). This also applies to ggplot objects, which only display when printed.

```
for (i in seq_along(x)) {
    message("Element ", i, ": ", x[i])
}
# Element 1: banana
# Element 2: apple
```

seq\_along() is used here rather than 1:length(x) as length(x) may be zero. message is used to print messages to the console.

### while and repeat

The while loop repeats while a condition is TRUE

The repeat loop repeats until exited by break

```
repeat {
    x <- x + 1
    if (max(x) > 10) break
}
```

break can be used in for or while loops too.

next can be used to skip to the next iteration.

### **Growing Objects**

Adding to an object in a loop, e.g. via c() or cbind()

```
res <- NULL
for (i in 1:5000) res <- c(res, 1)
```

forces a copy to be made at each iteration.

It is far better to create an object of the necessary size first

```
res <- numeric(5000)
for (i in seq_along(res)) res[i] <- 1</pre>
```

To initialise a list we can use

```
res <- vector(mode = "list", length = 100)
```

### Benchmarking

There will usually be many ways to write code for a given task. To compare alternatives, we can benchmark the expression

#### **Your Turn**

Load the broom package, as we will use glance in this exercise.

Assign to y the variable mtcars\$mpg. Assign to n the number of columns of mtcars minus 1. Create a vector named r.squared with this number of elements.

Write a for loop to iterate through the names of mtcars except mpg, that

- prints a message stating the current variable
- assigns the current variable to x
- fits a linear model using lm, regressing y on x
- uses glance to obtain the  $\mathbb{R}^2$  value and stores it in an element of r.squared corresponding to the current iteraton

Use which.max() to find the best predictor for mpg and the corresponding  $R^2$  value.

### Vectorization

Vectorization is operating on vectors (or vector-like objects) rather than individual elements.

Many operations in R are vectorized, e.g.

```
x <- 1:3
y <- 3:1
x == y
# [1] FALSE TRUE FALSE
res <- list(a = 1:3, b = 1:6)
lengths(res)
# a b
# 3 6</pre>
```

We do not need to loop through each element!

# Recycling

Vectorized functions will recycle shorter vectors to create vectors of the same length

```
1:4 + 0:1
# [1] 1 3 3 5
```

This is particularly useful for single values

```
cbind(1, 3:4)
# [,1] [,2]
# [1,] 1 3
# [2,] 1 4
```

and for generating regular patterns

```
paste0(rep(1:3, each = 2), c("a", "b"))
# [1] "1a" "1b" "2a" "2b" "3a" "3b"
```

### **Vectorization and Matrices**

Vectorizations applies to matices too, not only through matrix algebra

```
M <- matrix(1:4, nrow = 2, ncol = 2)
M + M
# [,1] [,2]
# [1,] 2 6
# [2,] 4 8</pre>
```

but also vectorized functions

```
M <- M + .3
round(M)
# [,1] [,2]
# [1,] 1 3
# [2,] 2 4
```

# **Matrices and Recycling**

Values are recycled down matrix, which is convenient for row-wise operations

To do the same for columns we would need to explicitly replicate, which is not so efficient.

```
M - rep(1:3, each = 2)

# [,1] [,2] [,3]

# [1,] 0 1 2

# [2,] 1 2 3
```

### **Vectorization vs For Loop**

Operations that can be vectorized will be more efficient than a loop in R

### **Row/Column-wise Operations**

Several functions are available implementing efficient row/column-wise operations, e.g. colMeans(), rowMeans(), colSums(), rowSums(), sweep()

```
M <- matrix(1:4, nrow = 2, ncol = 2)
rowMeans(M)
# [1] 2 3</pre>
```

These provide an alternative to iterating though rows and columns in R (the iteration happens in C, which is faster).

The matrixStats provides further "matricised" methods.

### **Iteration Functions**

Iteration functions provide a general alternative to for loops. They are not necessarily faster, but can be more compact.

apply() applies a function over rows/columns of a matrix.

lapply(), sapply() and vapply() iterate over a list or vector. vapply() is recommended for programming as it specifies the type of return value

```
vapply(list(a = 1:3, b = 1:6), FUN = mean, FUN.VALUE = numeric(1))
# a b
# 2.0 3.5
```

mapply() iterates over two or more lists/vectors in parallel.

The **purrr** packages provides alternatives to these that have a simpler, more consistent interface with fixed type of return value.

### Random Sampling

In order to test code, or when running simulations, you may want to generate random numbers. The numbers may be a random sample from a pool

```
pool <- 1:10
sample(pool, 7)
# [1] 3 4 5 7 2 8 9
sample(pool, 7, replace = TRUE)
# [1] 7 7 1 3 2 7 4</pre>
```

#### **Statistical Distributions**

Alternatively you can sample from a statistical distribution such as the standard normal

```
rnorm(7, mean = 0, sd = 1)
# [1] 0.7383 0.5758 -0.3054 1.5118 0.3898 -0.6212 -2.2147
```

or the uniform distribution

```
runif(7, min = 1, max = 100)
# [1] 87.10 34.69 48.73 60.36 49.86 19.44 82.91
```

Common distributions are available in the base package, e.g. rchisq(), rt(), others are provided in packages, e.g. mvrnorm() is provided by MASS.

#### Random Seeds

Sampling relies on generating random numbers.

We can fix the random seed so that all the code that follows will be repeatable

```
set.seed(1746)
sample(1:7)
# [1] 2 3 6 1 4 5 7
set.seed(1746)
sample(1:7)
# [1] 2 3 6 1 4 5 7
```

To repeatedly run the same code with different seeds, we can use replicate().

```
sim <- replicate(100, mean(rexp(10, rate = 10)))</pre>
```

### **Functions**

Functions are defined by two components:

- the arguments of the function
- the body of the function that computes the result

They are created using function()

```
t_statistic <- function(n) {
    x <- rnorm(n)
    y <- rnorm(n)
    t.test(x, y)$statistic
}</pre>
```

#### **Your Turn**

Find the code for the t\_statistic function in exercises.Rmd. Source in the function by running the function code. Try calling the function with different values of n to check it works.

Use replicate to simulate 1000 t-statistics comparing two samples of size 10 drawn from N(0, 1).

Theoretically the statistics should follow a t-distribution with 18 (= 10 + 10 - 2) degrees of freedom. Use rt to simulate 1000 values from this distribution.

Use the ggplot code in exercises. Rmd to plot compare the density of the simulated values with the density of the theoretical values.

### **Specified Arguments**

specified arguments are those named in the function definition, e.g. in rnorm()

```
args(rnorm)
# function (n, mean = 0, sd = 1)
# NULL
```

the arguments are n, mean and sd.

mean and sd have been given default values in the function definition, but n has not, so the function fails if the user does not pass a value to n

```
rnorm()
# Error in rnorm(): argument "n" is missing, with no default
```

### Name and Order of Arguments

The user can pass objects to these arguments using their names or by supplying unnamed values in the right order

```
rnorm(5, 1, 10)
# [1] -14.0073 -0.2515  7.9310 10.1248 -3.4390
rnorm(5, sd = 10)
# [1] 14.36 10.59 -9.35 16.62 -1.95
```

So naming and order is important! Some guidelines

- put compulsory arguments first, e.g. data
- put rarely used arguments last, e.g. tolerance setting
- use short but meaning argument names
- if relevant, use the same argument names as similar functions

### **Using Arguments**

Arguments are used as objects in the function code.

An new environment is created each time the function is called, separate from the global workspace.

```
x <- 1
y <- 3
f <- function(x, y){
    a <- 1
        x <- x + a
        x + y
}
f(x, y)
# [1] 5
x
# [1] 1
a
# Error in eval(expr, envir, enclos): object 'a' not found</pre>
```

# **Lexical Scoping**

If an object is not defined within the function, or passed in as an argument, R looks for it in the *parent environment* where the function was defined

```
x <- 1
y <- 3
f <- function(x){
    x + y
}
f(x)
# [1] 4
rm(y)
f(x)
# Error in f(x): object 'y' not found</pre>
```

It is safest to use arguments rather than depend on global variables!

# **Unspecified Arguments**

... or the *ellipsis* allow unspecified arguments to be passed to the function.

This device is used by functions that work with arbitrary numbers of objects, e.g.

```
args(sum)
# function (..., na.rm = FALSE)
# NULL
sum(1, 4, 10, 2)
# [1] 17
```

It can also be used to pass on arguments to another function, e.g.

```
t_statistic <- function(x, g, ...) {
    t.test(x ~ g, ...)$stat
}</pre>
```

# Using ...

Arguments passed to ... can be collected into a list for further analysis

```
f <- function(...){
    dots <- list(...)
    vapply(dots, mean, numeric(1), na.rm = TRUE)
}
x <- 1
y <- 2:3
f(x, y)
# [1] 1.0 2.5</pre>
```

Similarly the objects could be concatenated using c

### **Return Values**

By default, functions return the object created by the last line of code

```
f <- function(x) {
    x <- x + 1
    log(x)
}</pre>
```

Alternatively return() can be used to terminate the function and return a given object

```
f <- function(x) {
    if (all(x > 0)) return(log(x))
    x[x <= 0] <- 0.1
    log(x)
}</pre>
```

Multiple objects can be returned in a list.

# **Naming Functions**

As with arguments, function names are important:

- use a name that describes what it returns (e.g. t\_statistic) or what it does (e.g. remove\_na)
- try to use one convention for combining words (e.g. snake case t\_statistic or camel case tStatistic)
- avoid using the same name as other functions

### **Side Effects**

A side-effect is a change outside the function that occurs when the function is run, e.g.

- plot to the graphics window or other device
- printing output to the console
- write data to a file

A function can have many side-effects and a return value, but it is best practice to have a separate function for each task, e.g creating a plot or a table.

Writing to file is usually best done outside a function.

### **Your Turn**

In the qq\_norm chunk of exercises.Rmd there is some code to compute the slope and intercept of the line to add to a quantile-quantile plot, comparing sample quantiles against theoretical quantiles of a N(0,1) distribution.

Turn this code into a function named  $qq_norm$  taking the sample data as an argument and returning the slope and intercept in a list.

Run this chunk to source the function, then run the normal-QQ chunk which uses the qq\_norm function to compute parameters for an example plot.

Copy qq\_norm to the qq chunk and rename it qq. Add a new argument fun to specify any quantile function (e.g. qt, qf, etc). Give it the default value qnorm. Inside the function use qfun <- match.fun(fun) to get the quantile function matching fun, then use qfun instead of qnorm to compute qtheory. Use . . . to pass on arguments to qfun.

Save functions.R and run the t-QQ chunk in exercises.Rmd.

# **Using Functions From Other Packages**

In our own functions (outside of packages), it is possible to use library

```
scale_rows <- function(X){
    library(matrixStats)
    X <- X - rowMeans(X)
    X/rowSds(X)
}</pre>
```

But this loads the entire package, potentially leading to clashes with functions from other packages. It is better to use the **import** package:

```
scale_rows <- function(X){
   import::from(matrixStats, rowSds)
   X <- X - rowMeans(X)
   X/rowSds(X)
}
scale_rows(matrix(1:12, nrow = 3))</pre>
```

## **Custom ggplot**

**ggplot2**, like **dplyr** and other tidyverse packages, uses *non-standard evaluation*, that is, it refers to variable names as if they were objects in the current environment

```
ggplot(mtcars, aes(x = mpg, y= disp)) +
   geom_point()
```

To emulate this, we have to use tools from rlang: enquo then!!

```
ggscatter <- function(data, x, y){
   import::from(rlang, enquo, `!!`)
   import::from(ggplot2, ggplot, aes, geom_point)
   nse_x <- enquo(x)
   nse_y <- enquo(y)
   ggplot(data, aes(x = !! nse_x, y = !! nse_y)) +
        geom_point()
}
ggscatter(mtcars, x = mpg, y = disp)</pre>
```

# **Externalizing Function Code**

It is a good idea to separate function code from analysis code.

Put related functions together and source as required

```
source("modelFunctions.R")
source("plotFunctions.R")
```

The **import** package enables only necessary, top-level functions to be imported to the global workspace:

```
import::here(poissonModel, quasiPoissonModel, .from = "modelFunctions.R")
```

In either case, import::from commands can be put outside the function body to make the code easier to read.

# **Documenting Functions**

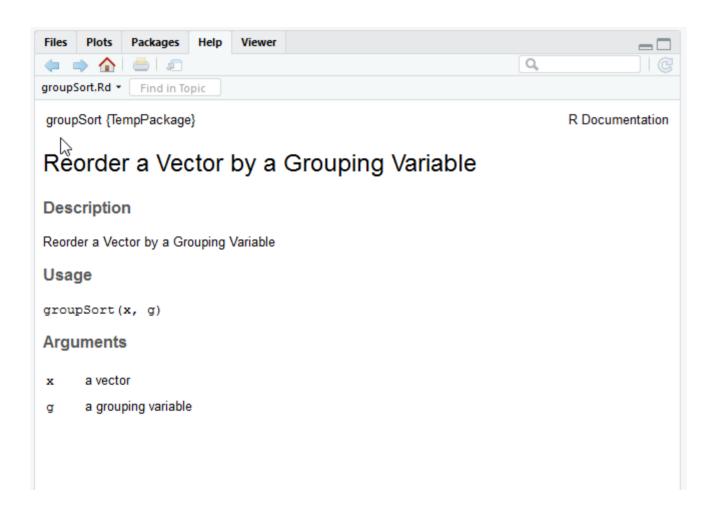
Comments help to record what a function does

```
# reorder x by grouping variable g
groupSort <- function(x, g) {
   ord <- order(g) #indices for ascending order of g
   x[ord]
}</pre>
```

The **docstring** package enables *roxygen* comments to be turned into a help file

```
library(docstring)
groupSort <- function(x, g) {
    #' Reorder a Vector by a Grouping Variable
    #'
    #' @param x a vector
    #' @param g a grouping variable
    ord <- order(g) #indices for ascending order of g
    x[ord]
}</pre>
```

#### ?groupSort



For fuller documentation, see the docstring vignette.

### **Validation**

When developing a function, we will want to validate its output.

A simple approach is to try different inputs

```
log_2 <- function(x){
    log(x, 2)
}
log_2(2^2)
# [1] 2
log_2(2^0)
# [1] 0</pre>
```

Doing this each time we change the function becomes tedious to check and error-prone as we miss important tests.

# Unit testing

The **testthat** packages allows us to create a test suite:

```
context("log_2 works correctly")

test_that("log_2 returns log to base 2", {
    expect_equal(log_2(2^3), 3)
    expect_equal(log_2(2^0), 0)
})

test_that("negative values give error", {
    expect_error(log_2(2^-1))
})
```

## **Running Tests**

If we save the tests in a file, e.g. tests.R, we can use test\_file() to run and check all tests:

```
library(testthat)
test file("tests.R")
# v | OK F W S | Context
# x | 2 1 | log_2 works correctly
 tests.R:9: failure: negative values give error
 \log_2(2^{-1}) did not throw an error.
# OK:
# Failed: 1
# Warnings: 0
# Skipped: 0
```

### **Your Turn**

Copy the qq function to a new R script and save as functions. R. Add roxygen comments at the start of the function body to define a title and parameter documentation.

Run the documentation chunk of exercises. Rmd to view your documentation.

Open the tests.R script. Using expect\_equal add some tests for the following

- a sample of 100,000 from N(0, 1) gives approximately slope 1, intercept 0
- a sample of 100,000 from N(0, 1/2) gives approximately slope 2, intercept 0
- sample of 100,000 from N(2, 1) gives approximately slope 1, intercept -2

Use the tol argument in expect\_equal to set a tolerance of 0.01.

Run the tests chunk of exercises. Rmd to run your tests with testfile. Try changing the expected tolerance to get a test to fail.

# **Sanity Checks**

To avoid mistakes, you may want to add some basic sanity checks

```
logit <- function(p){
    stopifnot(p > 0 & p < 1)
    log(p/(1 - p))
}
logit(2)
# Error in logit(2): p > 0 & p < 1 is not TRUE
logit(0.5)
# [1] 0</pre>
```

## **Error Messages**

Often the R messages can be quite obscure

```
zap <- function(x) if (max(x) < 1e7) 0 else x x <- c(1, 2, NA) zap(x) # Error in if (max(x) < 1e+07) 0 else x: missing value where TRUE/FALSE needed
```

More helpful error message can be implemented using stop

# Warning Messages

Warning messages should be given using warning()

```
safe_log2 <- function(x) {
    if (any(x == 0)) {
        x[x == 0] <- 0.1
        warning("zeros replaced by 0.1")
    }
    log(x, 2)
}
safe_log2(0:1)
# Warning in safe_log2(0:1): zeros replaced by 0.1
# [1] -3.322   0.000</pre>
```

Other messages can be printed using message().

# **Suppressing Warnings**

If a warning is expected, you may wish to suppress it

```
log(c(3, -1))
# Warning in log(c(3, -1)): NaNs produced
# [1] 1.099 NaN
x <- suppressWarnings(log(c(3, -1)))</pre>
```

All warnings will be suppressed however!

Similarly suppressMessages() will suppress messages.

# **Catching Errors/Warnings**

The purrr package has various functions to catch issues.

possibly() lets you modify a function to return a specified value when there is an error

```
log("a")
# Error in log("a"): non-numeric argument to mathematical function
library(purrr)
poss_log <- possibly(log, otherwise = NA)
poss_log("a")
# [1] NA</pre>
```

safely() works in a similar way but returns a list with elements "result" and "error", so you can record the error message(s).

quietly() lets you modify a function to return printed output, warnings and messages along with the result.

### traceback()

When an unexpected error occurs, there are several ways to track down the source of the error, e.g. traceback()

```
f1 <- function(x){ f2(x) }
f2 <- function(x){ x + qqqq }
f1(10)
# Error in f2(x): object 'qqqq' not found
traceback()
# 2: f2(2) at #1
# 1: f1(10)</pre>
```

In RStudio, if Debug > On Error > Error Inspector is checked and the traceback has at least 3 calls, the option to show traceback is presented

```
Error in x + d : non-numeric argument to binary operator

$\text{$L$ Show Traceback}$

Rerun with Debug
```

### debugonce()

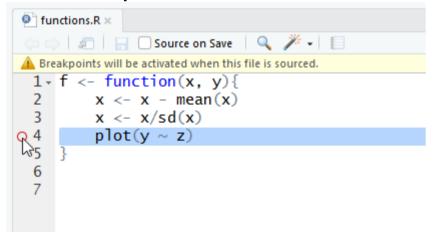
debugonce() flags a function for debugging the next time it is called

When in debug mode type n or  $\[ \leftarrow \]$  to step to the next line and c to continue to the end of a loop or the end of the function.

## **Breakpoints**

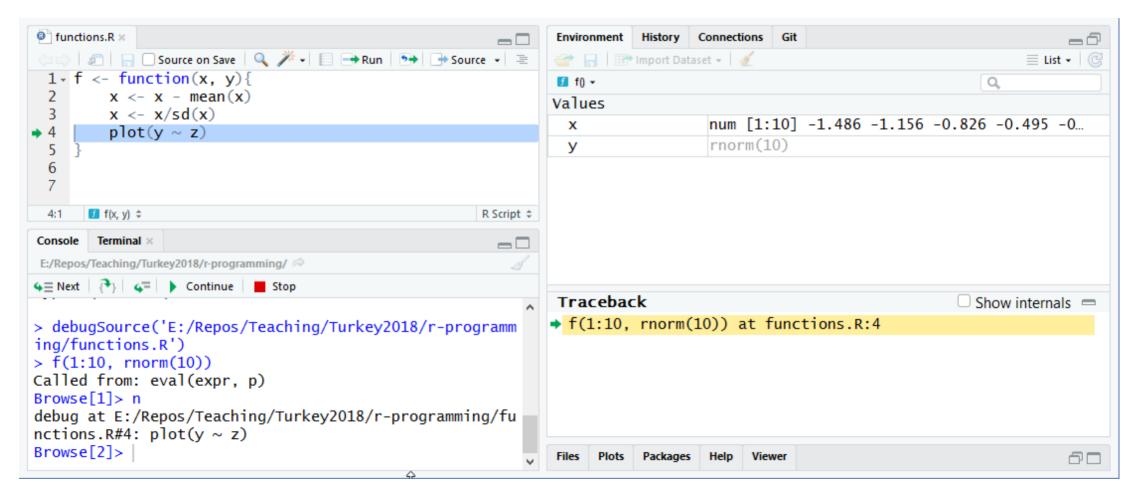
Stepping through a function line by line can be tedious. In RStudio we can set custom breakpoints in the source pane

### Set breakpoint in RStudio



#### Source the code

### Start debugging from breakpoints



# **RStudio's Rerun with Debug**

The Rerun with Debug option will rerun the command that created the error and enter debug mode where the error occurred.

### Good points:

- Easy to enter debug mode (when option shown)
- Can click in Traceback pane to view objects at any point in the call stack

### Bad points:

- May have gone past source of error (use breakpoints instead)
- May enter deeply nested function: use recover() to select an earlier entry point

Outside RStudio use options(error = recover), run code to debug, then set options(error = NULL).

### **Your Turn**

Open the function debug\_practice.R and source the example function f.

Try to run f(10) - there's an error! Use traceback() to see which function call generated the error, then fix the problem.

Run f(10) again - there is another error! Can you fix this directly?

Try running f(1) - is the result what you expected? Use debugonce() to setting debugging on f() and re-run f(1). Step through the function, printing each object as it is created to see what is happening.

Can you think how to improve the function? See if you can modify the function to give a sensible result for any integer.

### **Version Control**

When developing code, we often want to keep old versions.

We might save with different files names

or comment on changes

```
## ED 2018-09-28 use geom_bar instead of geom_histogram
## p <- p + geom_histogram(stat = "identity")
p <- p + geom_bar(stat = "identity")
## HT 2018-09-26 remove legend from plot
p <- p + theme(legend.position = "none")</pre>
```

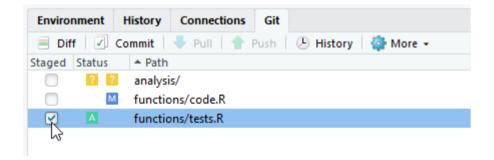
Either way it can get messy and hard to track/revert changes!

## git

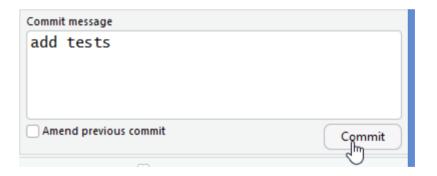
git is a version control system that allows us to record changes made to files in a directory.

When git is set up on an RStudio project (see refs at end) we can

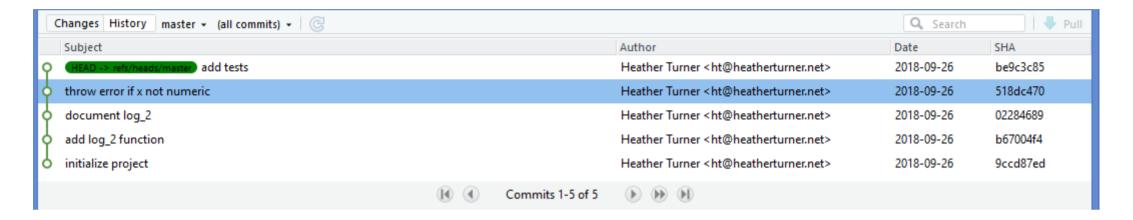
### stage changes to commit

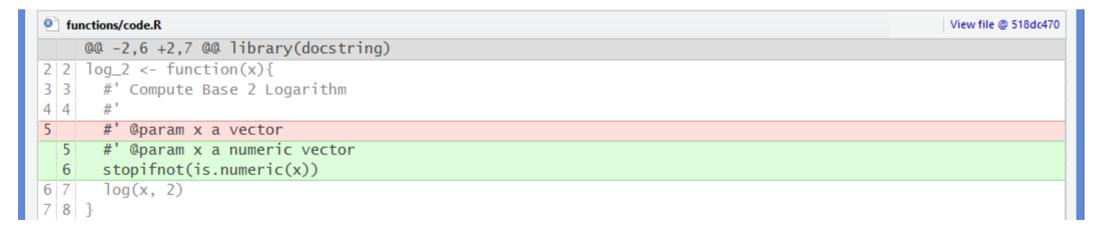


#### commit a set of changes



# **Commit History**





### GitHub/GitLab/Bitbucket

Services such as GitHub enable you to create a remote copy of your respository. You *push* commits to the remote repo (and *pull* if working with others).

### For all projects

provides a backup: you can checkout any version to reset the project

### For long-term/collaborative projects

- open issues: note and discuss changes to make
- collaborate with others: merge changes made by different people
- distribute your code: it can be installed directly from remote repo

# Writing an R Package

If using functions across many projects, or you want to share your functions with the wider world, it's best to put those functions in a package.

A package is built from the package source, which is a directory of the function code, tests, etc organised with a particular structure.

The **usethis** package helps to create the right structure and add components the the package, e.g. with create\_package() and use\_tests().

The **devtools** package helps to develop the package, e.g. with load\_all() to load the functions as if the package were installed and document() to create helpfiles from the roxygen comments.

# Package vs Stand-alone Function

	Package	Standalone function	
Function code	with related function code in R/	anywhere in any .R file	
roxygen comments	above function definition	in function body	
Imports	roxygen comments	in function .R (import::from)	
Exports	roxygen comments	in analysis .R (import::here)	
testthat tests	in tests/testthat/	in separate .R file	
Long-form docs	.Rmd in vignettes/	-	
Shared data	file in data/, roxygen in R/	-	
Package metadata	DESCRIPTION file	-	
Package intro	README.md	-	
Package news	NEWS.md	-	

# **Licensing and Sharing Packages**

You define the license in the package DESCRIPTION. This may be

- Proprietary: LICENSE file, e.g. "Do not distribute outside of Company X".
- Open source: you should adopt a standard license e.g. MIT or GPL-3.

You can share your package in different ways

	Quality standards	Installation	Discoverability
Own/department website	None	Medium	Hard
GitHub	Voluntary checks	Medium	Medium
CRAN	Compulsory checks	Easy	High
Bioconductor/rOpenSci	Code review	Easy	High

# **Further Reading**

### **Programming**

- R for Data Science <a href="http://r4ds.had.co.nz">http://r4ds.had.co.nz</a>
- Advanced R http://adv-r.had.co.nz/.

#### Git/GitHub

- Setup: Version Control with Git and SVN (RStudio support)
- Happy Git and GitHub for the useR <a href="http://happygitwithr.com/">http://happygitwithr.com/</a>

### **Packages**

- Forwards Package Development workshops https://github.com/forwards/workshops/
- R Packages http://r-pkgs.had.co.nz/