Introduction to Functional Programming

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An introduction to writing functions to improve your coding efficiency and optimize performance.

Learning objectives

- 1. Write and use functions in R
- 2. Document functions for easy re-use
- 3. Replace loops with functions optimized for vector calculations

Don't Repeat Yourself

The DRY principle aims to reduce repetition in software engineering. By writing and using functions to accomplish a set of instructions multiple times, you reduce the opportunities for mistakes and often improve performance of the code you write. Functional programming makes it easy to apply the same analyses to different sets of data, without excessive copy-paste-update cycles that can introduce hard-to-detect errors.

Writing functions

Why do we write functions? Usually, we create functions after writing some code with certain variables, then copy/pasting that code and changing the variable names. Then more copy/paste. Then we forget to change one of the variables and spend a day figuring out why our results don't make sense.

For example, consider the iris data set:

head(iris)

```
Sepal.Length Sepal.Width Petal.Length Petal.Width Species
##
## 1
              5.1
                           3.5
                                         1.4
                                                      0.2 setosa
## 2
              4.9
                           3.0
                                         1.4
                                                      0.2 setosa
## 3
              4.7
                           3.2
                                                      0.2
                                         1.3
                                                           setosa
## 4
              4.6
                           3.1
                                         1.5
                                                      0.2
                                                           setosa
## 5
              5.0
                           3.6
                                         1.4
                                                      0.2 setosa
## 6
              5.4
                           3.9
                                         1.7
                                                      0.4 setosa
```

And if we look at the values for each of the four measurements, some are quite different in their ranges.

summary(iris)

```
##
     Sepal.Length
                      Sepal.Width
                                       Petal.Length
                                                        Petal.Width
##
           :4.300
                                              :1.000
    Min.
                     Min.
                             :2.000
                                      Min.
                                                       Min.
                                                               :0.100
##
    1st Qu.:5.100
                     1st Qu.:2.800
                                      1st Qu.:1.600
                                                       1st Qu.:0.300
   Median :5.800
##
                     Median :3.000
                                      Median :4.350
                                                       Median :1.300
   Mean
           :5.843
                     Mean
                            :3.057
                                      Mean
                                              :3.758
                                                       Mean
                                                               :1.199
##
    3rd Qu.:6.400
                     3rd Qu.:3.300
                                      3rd Qu.:5.100
                                                       3rd Qu.:1.800
           :7.900
                                              :6.900
##
    Max.
                     Max.
                             :4.400
                                      Max.
                                                       Max.
                                                               :2.500
##
          Species
    setosa
               :50
```

```
## versicolor:50
## virginica:50
##
```

Want to standardize so each has mean zero and variance of one

New script file, iris-stand.R

```
# Standarize iris values
# Jeff Oliver
# jcoliver@email.arizona.edu
# 2018-08-16

# Standardize petal length
mean.petal.length <- mean(iris$Petal.Length)
sd.petal.length <- sd(iris$Petal.Length)
standard.petal.length <- (iris$Petal.Length - mean.petal.length)/sd.petal.length</pre>
```

Check process

```
summary(standard.petal.length)
```

```
## Min. 1st Qu. Median Mean 3rd Qu. Max.
## -1.5623 -1.2225 0.3354 0.0000 0.7602 1.7799
```

Repeat for petal width, copy/paste from above

```
# Standardize petal length
mean.petal.width <- mean(iris$Petal.Width)
sd.petal.width <- sd(iris$Petal.Width)
standard.petal.width <- (iris$Petal.Width - mean.petal.width)/sd.petal.width
summary(standard.petal.width)</pre>
```

```
## Min. 1st Qu. Median Mean 3rd Qu. Max.
## -1.4422 -1.1799 0.1321 0.0000 0.7880 1.7064
```

Did copy/paste really save time? And how many opportunities for mistakes?

Write a function!

```
standardize <- function(data) {
  mean.value <- mean(data)
  sd.value <- sd(data)
  z.value <- (data - mean.value)/sd.value
  return(z.value)
}</pre>
```

Replace previous work

```
standard.petal.length <- standardize(data = iris$Petal.Length)
standard.petal.width <- standardize(data = iris$Petal.Width)</pre>
```

Document the function

```
# Standardize a vector of numerical data
# data: vector of numerical data
# returns: vector of original data standardized to have mean = 0, sd = 1
standardize <- function(data) {
   mean.value <- mean(data)</pre>
```

```
sd.value <- sd(data)
  z.value <- (data - mean.value)/sd.value
  return(z.value)
}
Use apply to do it to multiple columns
iris.stand <- apply(X = iris, MARGIN = 2, FUN = standardize)</pre>
iris.stand <- apply(X = iris[, 1:4], MARGIN = 2, FUN = standardize)</pre>
iris.stand <- as.data.frame(iris.stand)</pre>
Add the species column
iris.stand$Species <- iris$Species</pre>
Write a function to do it all (but do this in a separate file - move our other function there, too)
# data: data frame to standardize
# ignore: (optional) vector of columns to skip in standardization process
# returns: data frame with all non-ignored columns standardized to have mean = 0, sd = 1
standardize.df <- function(data, ignore = c()) {</pre>
  ignored.cols <- data[, ignore]</pre>
  standardized.df <- apply(X = data[, -ignore], MARGIN = 2, FUN = standardize)
  standardized.df <- cbind(standardized.df, ignored.cols)</pre>
  return(standardized.df)
}
iris.stand <- standardize.df(data = iris, ignore = 5)</pre>
# data: data frame to standardize
# ignore: (optional) vector of columns to skip in standardization process
# returns: data frame with all non-ignored columns standardized to have mean = 0, sd = 1
standardize.df <- function(data, ignore = c()) {</pre>
  ignored.cols <- as.data.frame(data[, ignore])</pre>
  colnames(ignored.cols) <- colnames(data)[ignore]</pre>
  standardized.df <- apply(X = data[, -ignore], MARGIN = 2, FUN = standardize)
  standardized.df <- cbind(standardized.df, ignored.cols)</pre>
  return(standardized.df)
iris.stand <- standardize.df(data = iris, ignore = 5)</pre>
```

Old Lesson

Learning objectives

- 1. Write and use functions in R
- 2. Document functions for easy re-use
- 3. Replace loops with functions optimized for vector calculations

Don't Repeat Yourself

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Writing functions

Why do we write functions? Usually, we create functions after writing some code with certain variables, then copy/pasting that code and changing the variable names. Then more copy/paste. Then we forget to change one of the variables and spend a day figuring out why our results don't make sense.

For example, consider the airquality data set:

```
##
     Ozone Solar.R Wind Temp Month Day
## 1
        41
                190 7.4
                            67
                                        2
## 2
        36
                118 8.0
                            72
                                   5
## 3
        12
                149 12.6
                            74
                                   5
                                        3
## 4
        18
                313 11.5
                            62
                                   5
                                        4
                                    5
                                        5
## 5
        NA
                 NA 14.3
                            56
## 6
        28
                 NA 14.9
                                    5
                                        6
                            66
```

Let's start by writing a script to run linear regression to see if the solar radiation (the Solar.R column) predicts the ozone level (the Ozone column). Create a new file called "airquality-regression.R":

```
# Analyze air quality data
# Jeffrey Oliver
# jcoliver@email.arizona.edu
# 2017-06-22

# Relationship between ozone and solar radiation
simple <- lm(airquality[, "Ozone"] ~ airquality[, "Solar.R"])</pre>
```

Now we are interested in just the correlation coefficient (r^2) and the p-value for this relationship, so we use summary and extract those values:

```
# Extract the model parameters
simple.summary <- summary(simple)
simple.r2 <- simple.summary$r.squared
simple.p <- simple.summary$coefficients[2, 4]</pre>
```

And finally we can print out these two values with the cat command:

```
cat("Solar r^2 =", simple.r2)
cat("Solar p =", simple.p)
```

```
## Solar r^2 = 0.1213419Solar p = 0.0001793109
```

If we want to do the same thing for the relationship between ozone and wind, we can copy/paste this code and just change the predictor column specification in the 1m command from "Solar.R" to "Wind" and update the message in the cat commands:

```
simple <- lm(airquality[, "Ozone"] ~ airquality[, "Wind"])
simple.summary <- summary(simple)
simple.r2 <- simple.summary$r.squared</pre>
```

```
simple.p <- simple.summary$coefficients[2, 4]
cat("Wind r^2 =", simple.r2)
cat("Wind p =", simple.p)</pre>
```

This isn't too great an effort, but as the number of predictors grows, the time and opportunty for mistakes also grows. Since we are doing the same exact process for each predictor variable (run lm, run summary, extract values of interest, then print values of interest), we can encapsulate this in a function.

Behold, a function!

To define a function, create a new file called "regression-functions.R". While functions do not necessarily have to exist outside of the R script in which they are called, it is generally considered good practice. Start this file with the usual header containing information about the contents:

```
# Functions to automate linear regression
# Jeff Oliver
# jcoliver@email.arizona.edu
# 2017-06-15
```

And we define a function pretty much the same way we assign values to a variable, but here we use the function function:

```
RegressSimple <- function() {
}</pre>
```

We have a general idea of what the function should do, so write that as a brief comment above the function:

```
# Run linear regression for all predictors and a response variable in a data frame
RegressSimple <- function() {
}</pre>
```

Now what do we do? How does one actually go about writing functions? Maybe you know exactly what your function should do and all the inputs and outputs. I rarely find myself in that case. Rather, I start by doing the same thing as I did before: I copy paste the portion of code I want to run. So in this case, from the airquality-regression.R file, copy the section for the solar radiation analysis and paste it into the body of the RegressSimple function. That is, paste it between the pair of curly braces { }:

```
# Run linear regression for all predictors and a response variable in a data frame
RegressSimple <- function() {
    simple <- lm(airquality[, "Ozone"] ~ airquality[, "Solar.R"])
    simple.summary <- summary(simple)
    simple.r2 <- simple.summary$r.squared
    simple.p <- simple.summary$coefficients[2, 4]
    cat("Solar r^2 =", simple.r2)
    cat("Solar p =", simple.p)
}</pre>
```

We don't want the function to print out values from the models, so delete the code that extracts the values and prints them with the two cat statements:

```
# Run linear regression for all predictors and a response variable in a data frame
RegressSimple <- function() {
   simple <- lm(airquality[, "Ozone"] ~ airquality[, "Solar.R"])
   simple.summary <- summary(simple)
}</pre>
```

At this point, we can consider input and output of the function. We'll need to give the function two pieces of input: the data to work with and which variable to use as the response. We specify input by declaring the

names of the values (data and response) in the function call:

```
# Run linear regression for all predictors and a response variable in a data frame
RegressSimple <- function(data, response) {
   simple <- lm(airquality[, "Ozone"] ~ airquality[, "Solar.R"])
   simple.summary <- summary(simple)
}</pre>
```

And while we're at it, we need to document these inputs:

```
# Run linear regression for all predictors and a response variable in a data frame
# data: the data frame to analyze
# response: the name of the response variable
RegressSimple <- function(data, response) {
   simple <- lm(airquality[, "Ozone"] ~ airquality[, "Solar.R"])
   simple.summary <- summary(simple)
}</pre>
```

Now that we have inputs, we can update the variables in the code we copied from the regression script. What do we need to update?

- The name of the data frame we used, airquality is replaced by data
- The model specification now uses the abstracted response variable, stored in response instead of the hard coded "Ozone"; note that we *do not* put response in double quotes

```
# Run linear regression for all predictors and a response variable in a data frame
# data: the data frame to analyze
# response: the name of the response variable
RegressSimple <- function(data, response) {
   simple <- lm(data[, response] ~ data[, "Solar.R"])
   simple.summary <- summary(simple)
}</pre>
```

The goal of the function is to run linear regression for *all* the predictors in the data frame, so how can we do this? The simplest way is to use a **for** loop for all the columns in the data frame, updating the **lm** call with the column specification of the predictor variable:

```
# Run linear regression for all predictors and a response variable in a data frame
# data: the data frame to analyze
# response: the name of the response variable
RegressSimple <- function(data, response) {
  for (predictor in 1:ncol(data)) {
    simple <- lm(data[, response] ~ data[, predictor])
    simple.summary <- summary(simple)
  }
}</pre>
```

For output, we probably want the results of the linear model for each of our predictors. In this case, we'll use a list object and assigning the output of summary to an element in that list. Note because we are using a list, we use two-bracket notation [[]] to indicate an element in the list. And finally, we sent back these results with the return function:

```
# Run linear regression for all predictors and a response variable in a data frame
# data: the data frame to analyze
# response: the name of the response variable
RegressSimple <- function(data, response) {
  model.summaries <- list()
  for (predictor in 1:ncol(data)) {
    simple <- lm(data[, response] ~ data[, predictor])</pre>
```

```
element.name <- colnames(data)[predictor]
  model.summaries[[element.name]] <- summary(simple)
}
return(model.summaries)
}</pre>
```

Before we try this out, update the documentation with a description of the output

```
# Run linear regression for all predictors and a response variable in a data frame
# data: the data frame to analyze
# response: the name of the response variable
# returns: a list where each element is the output from summary(lm)
RegressSimple <- function(data, response) {
    model.summaries <- list()
    for (predictor in 1:ncol(data)) {
        simple <- lm(data[, response] ~ data[, predictor])
        element.name <- colnames(data)[predictor]
        model.summaries[[element.name]] <- summary(simple)
    }
    return(model.summaries)
}</pre>
```

Using our function

So how do we use this? We need to load this function into memory so we can use it. Go back to the script with our original regression analyses, "airquality-regression.R". Comment out our previous linear regression code and add a call to source to load our function file. Hint: in RStudio, you can select multiple lines and comment them out with the shortcut Shift-Ctrl-C.

```
# Analyze air quality data
# Jeffrey Oliver
# jcoliver@email.arizona.edu
# 2017-06-22

source(file = "regression-functions.R")

# Relationship between ozone and solar radiation
# simple <- lm(airquality[, "Ozone"] ~ airquality[, "Solar.R"])

# Extract the model parameters
# simple.summary <- summary(simple)
# simple.r2 <- simple.summary$r.squared
# simple.p <- simple.summary$coefficients[2, 4]
# cat("Solar r^2 = ", simple.r2)
# cat("Solar p = ", simple.p)</pre>
```

And we can now use this function, passing airquality as the data and "Ozone" as the response:

```
airquality.models <- RegressSimple(data = airquality, response = "Ozone")</pre>
```

```
## Warning in summary.lm(simple): essentially perfect fit: summary may be
## unreliable
```

Hmmmm...that's an odd warning. Let's add some reporting code to see if we can figure out what it's doing. Update the RegressSimple function to print the name of the predictor using the cat function:

```
# Run linear regression for all predictors and a response variable in a data frame
# data: the data frame to analyze
# response: the name of the response variable
# returns: a list where each element is the output from summary(lm)
RegressSimple <- function(data, response) {
    model.summaries <- list()
    for (predictor in 1:ncol(data)) {
        simple <- lm(data[, response] ~ data[, predictor])
        element.name <- colnames(data)[predictor]
        model.summaries[[element.name]] <- summary(simple)
        cat("Predictor:", element.name, "\n")
    }
    return(model.summaries)
}</pre>
```

Now go back our airquality-regression.R script and run the RegressSimple command again.

```
airquality.models <- RegressSimple(data = airquality, response = "Ozone")</pre>
```

```
## Warning in summary.lm(simple): essentially perfect fit: summary may be
## unreliable
```

Drat. Same warning, but no message printed? Why not? Because we made changes to the regression-functions.R file, but did not load them into memory with the source command, R is using the old version of the RegressSimple function. So we need to run the source command first:

```
source(file = "regression-functions.R")
```

Then the RegressSimple command:

Predictor: Month
Predictor: Day

```
airquality.models <- RegressSimple(data = airquality, response = "Ozone")

## Warning in summary.lm(simple): essentially perfect fit: summary may be
## unreliable

## Predictor: Ozone

## Predictor: Solar.R

## Predictor: Wind

## Predictor: Temp</pre>
```

OK, so there are some problems. First, we aren't really interested in the effect of Day or Month on ozone levels, so when we call RegressSimple, we should only pass it data we want to analyze. Here we drop the fifth and sixth columns, which are the Month and Day columns, respectively:

```
airquality.models <- RegressSimple(data = airquality[, -c(5:6)], response = "Ozone")</pre>
```

But look at the output from function call again. Our function actually ran regression on a model using ozone to predict ozone - that's probably what caused the warning message "summary may be unreliable". But we don't want to exclude ozone from the data we pass to the function, because that is the response variable. We therefore need to update our function definition so we don't run an ozone ~ ozone model. More generally, we need to make sure our response variable is not treated as a predictor. To do this, we:

- $1. \ \, {\rm Find} \,\, {\rm out} \,\, {\rm which} \,\, {\rm column} \,\, {\rm is} \,\, {\rm the} \,\, {\rm response} \,\, {\rm variable}$
- 2. Create a vector of predictor variable names using colnames
- 3. Update our for loop to only use those predictor variables
- 4. Use the predictor variable name for the model.summaries list element name

Open the regression-functions.R file and update RegressSimple:

```
# Run linear regression for all predictors and a response variable in a data frame
# data: the data frame to analyze
# response: the name of the response variable
# returns: a list where each element is the output from summary(lm)
RegressSimple <- function(data, response) {
    response.index <- which(colnames(data) == response)
    predictors <- colnames(data)[-response.index]
    model.summaries <- list()
    for (predictor in predictors) {
        simple <- lm(data[, response] ~ data[, predictor])
        model.summaries[[predictor]] <- summary(simple)
        cat("Predictor:", predictor, "\n")
    }
    return(model.summaries)
}</pre>
```

Now when we run RegressSimple, there are only three predictors used in the models, as we expect:

```
source(file = "regression-functions.R")
airquality.models <- RegressSimple(data = airquality[, -c(5:6)], response = "Ozone")

## Predictor: Solar.R

## Predictor: Wind

## Predictor: Temp

Since it RegressSimple is now only running the models we want, remove the cat command from the function:

# Run linear regression for all predictors and a response variable in a data frame

# data: the data frame to analyze

# response: the name of the response variable

# returns: a list where each element is the output from summary(lm)</pre>
```

```
# data: the data frame to unaryse
# response: the name of the response variable
# returns: a list where each element is the output from summary(lm)
RegressSimple <- function(data, response) {
    response.index <- which(colnames(data) == response)
    predictors <- colnames(data)[-response.index]
    model.summaries <- list()
    for (predictor in predictors) {
        simple <- lm(data[, response] ~ data[, predictor])
        model.summaries[[predictor]] <- summary(simple)
    }
    return(model.summaries)
}</pre>
```

Because airquality.models is a list, we can access the objects using double-bracket notation:

```
solar.model <- airquality.models[["Solar.R"]]
solar.corr <- solar.model$r.squared
solar.p <- solar.model$coefficients[2, 4]
cat("Solar r^2 = ", solar.corr, "\n")
cat("Solar p = ", solar.p, "\n")

## Solar r^2 = 0.1213419
## Solar p = 0.0001793109</pre>
```

Make it Class-y

But now we're back to the copy-paste-update cycle if we want to get all the correlation coefficients and p-values. If we know that's all we want, we can create another function, one that does the work of heavy lifting of extracting coefficients and printing them out to the screen.

In the file with our RegressSimple function, create another function, and call it print.RegressSimple:

```
print.RegressSimple <- function(x, ...) {
}</pre>
```

Briefly, what we are doing is creating a function that specifies the output of anything that is class RegressSimple (we'll get to how we make that happen in a moment). The thing to note now is that in a print.___ function, the first argument, x is the object we wish to print; in this case, a product of the function RegressSimple. Stick with me here, it will become clearer. For the purposes of this lesson, just remember that in the print.RegressSimple function, the variable x is the list that is produced from a call to RegressSimple. We want this function to extract the r² and p-values for each model and display a table of those values for each predictor variable. Something like:

```
## variable r2 p

## Solar.R 0.12 1.7e-04

## Wind 0.36 9.2e-13

## Temp 0.48 2.9e-18
```

Add a brief explanation of what this function does, then add code to get the names of the variables and set up a data frame to hold the values we want to print:

Now extract the values we want to print for each element in the RegressSimple object (which is the variable x in print.RegressSimple):

Finally, we add cat and print statements to output the values.

```
# Print values of interest from each predictor in a RegressSimple object
print.RegressSimple <- function(x, ...) {
    # Get a vector of the elements' names</pre>
```

Now when we run our code, we can just enter the name of the variable to have it print out our nice table of just the r^2 and p-values.

```
source(file = "regression-functions.R")
airquality.models <- RegressSimple(data = airquality[, -c(5:6)], response = "Ozone")
airquality.models
## $Solar.R
##
## lm(formula = data[, response] ~ data[, predictor])
## Residuals:
      Min
               1Q Median
                               3Q
                                      Max
## -48.292 -21.361 -8.864 16.373 119.136
## Coefficients:
                    Estimate Std. Error t value Pr(>|t|)
## (Intercept)
                    18.59873
                                6.74790 2.756 0.006856 **
## data[, predictor] 0.12717
                                0.03278 3.880 0.000179 ***
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
## Residual standard error: 31.33 on 109 degrees of freedom
    (42 observations deleted due to missingness)
## Multiple R-squared: 0.1213, Adjusted R-squared: 0.1133
## F-statistic: 15.05 on 1 and 109 DF, p-value: 0.0001793
##
##
## $Wind
##
## Call:
## lm(formula = data[, response] ~ data[, predictor])
## Residuals:
      Min
               1Q Median
                               3Q
                                      Max
## -51.572 -18.854 -4.868 15.234 90.000
##
## Coefficients:
```

```
##
                     Estimate Std. Error t value Pr(>|t|)
                                 7.2387
                                          13.38 < 2e-16 ***
## (Intercept)
                     96.8729
## data[, predictor]
                                          -8.04 9.27e-13 ***
                     -5.5509
                                 0.6904
## ---
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
##
## Residual standard error: 26.47 on 114 degrees of freedom
##
     (37 observations deleted due to missingness)
## Multiple R-squared: 0.3619, Adjusted R-squared: 0.3563
## F-statistic: 64.64 on 1 and 114 DF, p-value: 9.272e-13
##
##
## $Temp
##
## Call:
## lm(formula = data[, response] ~ data[, predictor])
##
## Residuals:
##
      Min
                1Q Median
                               30
                                      Max
##
  -40.729 -17.409
                   -0.587
                          11.306 118.271
##
## Coefficients:
                     Estimate Std. Error t value Pr(>|t|)
##
                                 18.2872 -8.038 9.37e-13 ***
## (Intercept)
                     -146.9955
## data[, predictor]
                        2.4287
                                  0.2331 10.418 < 2e-16 ***
## Signif. codes: 0 '***' 0.001 '**' 0.05 '.' 0.1 ' ' 1
## Residual standard error: 23.71 on 114 degrees of freedom
     (37 observations deleted due to missingness)
## Multiple R-squared: 0.4877, Adjusted R-squared: 0.4832
## F-statistic: 108.5 on 1 and 114 DF, p-value: < 2.2e-16
```

Well that didn't work. It just printed out the each element of the list. That's because we need to take one final step. The print.RegressSimple only works on objects that are of class RegressSimple. So what is the class of our airquality.models?

```
class(airquality.models)
```

[1] "list"

It's a list, *not* a RegressSimple object. So how do we instruct R to make the output of RegressSimple to be an object of class RegressSimple? Big surprise here, we use the class function! In our RegressSimple function, right before we return the model.summaries list, we set the class to be "RegressSimple":

```
# Run linear regression for all predictors and a response variable in a data frame
# data: the data frame to analyze
# response: the name of the response variable
# returns: a list where each element is the output from summary(lm)
RegressSimple <- function(data, response) {
   response.index <- which(colnames(data) == response)
   predictors <- colnames(data)[-response.index]
   model.summaries <- list()
   for (predictor in predictors) {
      simple <- lm(data[, response] ~ data[, predictor])
      model.summaries[[predictor]] <- summary(simple)</pre>
```

```
# Set class and return results
  class(model.summaries) <- "RegressSimple"</pre>
  return(model.summaries)
Now re-run the lines in our regression script:
source(file = "regression-functions.R")
airquality.models <- RegressSimple(data = airquality[, -c(5:6)], response = "Ozone")
airquality.models
## Regression results:
##
       variable r2
                           р
## [1,] Solar.R 0.1213419 1.793109e-04
                 0.3618582 9.271974e-13
## [2,] Wind
## [3,] Temp
                 0.4877072 2.931897e-18
There it is, our table of values! Our final two files will then be:
airquality-regression.R:
# Analyze air quality data
# Jeffrey Oliver
# jcoliver@email.arizona.edu
# 2017-06-22
source(file = "regression-functions.R")
airquality.models <- RegressSimple(data = airquality[, -c(5:6)], response = "Ozone")
airquality.models
regression-functions.R:
# Functions to automate linear regression
# Jeff Oliver
# jcoliver@email.arizona.edu
# 2017-06-15
# Run linear regression for all predictors and a response variable in a data frame
# data: the data frame to analyze
# response: the name of the response variable
# returns: a list where each element is the output from summary(lm)
RegressSimple <- function(data, response) {</pre>
  response.index <- which(colnames(data) == response)</pre>
  predictors <- colnames(data)[-response.index]</pre>
  model.summaries <- list()</pre>
  for (predictor in predictors) {
    simple <- lm(data[, response] ~ data[, predictor])</pre>
   model.summaries[[predictor]] <- summary(simple)</pre>
  }
  # Set class and return results
  class(model.summaries) <- "RegressSimple"</pre>
  return(model.summaries)
```

}

```
# Print values of interest from each predictor in a RegressSimple object
print.RegressSimple <- function(x, ...) {</pre>
  # Get a vector of the elements' names
 predictors <- names(x)</pre>
  # Set up a dataframe for results of interest
 model.results <- data.frame(variable = predictors,</pre>
                         r2 = 0,
                         p = 0
  # Extract r-squared and p-values
  model.results$r2 <- sapply(x, "[[", "r.squared")</pre>
  model.coeffs <- sapply(x, "[[", "coefficients")</pre>
  model.results$p <- model.coeffs[8, ]</pre>
  # Print values
  cat("Regression results: ", "\n")
  print(as.matrix(model.results), quote = FALSE)
```

Additional resources

- A deeper dive to functional programming
- An even deeper dive into functional programming
- Some opinions and suggestions for naming things like functions (see the 'Object names' section)
- A PDF version of this lesson

Back to learn-r main page

Questions? e-mail me at jcoliver@email.arizona.edu.