Statistical Computing (36-350)

Lecture 11: Refactoring

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Agenda

- Abstraction adjusts programming to human strengths
- Refactoring adjusts code to bring out commonalities
- Ways of refactoring: names, objects, common operations, general operations
- Example: The jack-knife

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Abstraction — hiding details and specifics, dealing in generalities and common patterns — is a way to do this

We have talked about lots of examples of this already

Data structures; Functions; Interfaces; Functions as objects

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Once we have some code, and it (more or less) works, re-write it to emphasize commonalities:

- Parallel and transparent naming
- Grouping related values into objects
- Common or parallel sub-tasks become shared functions
- Common or parallel over-all tasks become general functions



Naming

R puts next to no limits on names of variables and functions :. we should use names that make sense to humans

- Names should indicate purpose or meaning
 E.g., call something plot or predict when, but only when, it plots or predicts
- Similar objects should have similar names.

Example: conventions for functions related to random variables

dnorm probability density of normal r.v.
rnorm random value from normal r.v.
pnorm cumulative probability of normal r.v.
quantile of normal r.v.
quantile of normal r.v.
probability density of gamma r.v.

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qchisq ?
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- people (including you) waste time less puzzling over it
- you are more easily replaced as a programmer

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Create a single data object (data frame, list, ...) that includes them all as parts

Replace mentions of the individual variables with mentions of parts of the unified object

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Makes sure that the right values are always present (pass the object as an argument to functions, rather than the components)
Memorization: if you know you are going to want to do the same calculation many times on these data values, do it once when you create the object, and store the result as a component

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Write one function to do the general operation, with additional arguments (typically including functions)

Call the new general function with appropriate arguments, rather than the old functions

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Old functions provide test cases to check if general function works

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Extended example: the jackknife

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```
Have an estimator \hat{\theta} of parameter \theta want the standard error of our estimate, se_{\hat{\theta}} The jackknife approximation: omit case i, get estimate \hat{\theta}_{(-i)} Take the variance of all the \hat{\theta}_{(-i)} multiply that variance by \frac{(n-1)^2}{n} to get \approx variance of \hat{\theta} then se_{\hat{\theta}} = square root of that variance (Why \frac{(n-1)^2}{n}? Think about just getting the standard error of the mean)
```

Jackknife for gamma parameters

```
gamma.jackknife <- function(data) {
    n <- length(data)
    jackknife.ests <- matrix(NA,nrow=2,ncol=n)
    rownames(jackknife.ests) = c("a","s")
    for (omitted.point in 1:n) {
        fit <- gamma.est(data[-omitted.point])
            jackknife.ests["a",omitted.point] <- fit$a
            jackknife.ests["s",omitted.point] <- fit$s
    }
    variance.of.ests <- apply(jackknife.ests,1,var)
    jackknife.vars <- ((n-1)^2/n)*variance.of.ests
    jackknife.stderrs <- sqrt(jackknife.vars)
    return(jackknife.stderrs)
}</pre>
```

Jackknife for the mean

```
mean.jackknife <- function(data) {
    n <- length(data)
    jackknife.ests <- vector(length=n)
    for (omitted.point in 1:n) {
        new.mean <- mean(data[-omitted.point])
    }
    variance.of.ests <- var(new.mean)
    jackknife.var <- ((n-1)^2/n)*variance.of.ests
    jackknife.stderr <- sqrt(jackknife.vars)
    return(jackknife.stderr)
}</pre>
```

Jackknife for linear regression coefficients

```
jackknife.lm <- function(data,p) {
  n <- nrow(data)
  jackknife.ests <- matrix(0,nrow=p,ncol=n)
  for (omit in 1:n) {
    new.coefs <- lm(YOUR.FORMULA.HERE,data=data[-omit,])$coefficients
    jackknife.ests[,omit] <- new.coefs
}
variance.of.ests <- apply(jackknife.ests,1,var)
  jackknife.var <- ((n-1)^2/n)*variance.of.ests
  jackknife.stderr <- sqrt(jackknife.vars)
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```

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figure out the size of the data
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scale up variances
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Refactor by extracting the common "omit one" operation Refactor by defining a general "jackknife" operation

The Common Operation

Works for vectors, lists, 1D and 2D arrays, matrices, data frames:

```
omit.case <- function(data,i) {
  d <- dim(data)
  if (is.null(d) || (length(d)==1)) {
    return(data[-i])
  } else {
    return(data[-i,])
  }
}</pre>
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EXERCISE: Modify so it also handles higher-dimensional arrays

The General Operation

```
jackknife <- function(estimator,data) {
  if (is.null(dim(data))) { n <- length(data) }
  else { n <- nrow(data) }
  jackknife.ests <- c()
  for (omit in 1:n) {
    reestimate <- estimator(omit.case(data,omit))
    jackknife.ests <- cbind(jackknife.ests,reestimate)
  }
  var.of.reestimates <- apply(jackknife.ests,1,var)
  jackknife.var <- ((n-1)^2/n)* var.of.reestimates
  jackknife.stderr <- sqrt(jackknife.var)
  return(jackknife.stderr)
}</pre>
```

Could allow other arguments to estimator, spin off finding n as its own function, etc.

It works

Summary

Refactoring adjusts code to emphasize patterns

- Names are informative and systematic
- Objects keep related values together
- Common sub-tasks become specialized lower-level functions
- General patterns of operations become high-level general functions

Refactoring makes code look more like top-down design Refactoring usually involves abstraction Abstraction emphasizes human strengths