

Statistical Computing (36-350)

Lecture 9: Functions as Objects

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- Functions are objects, and can be arguments to other functions
 - Example: `curve`
 - Example: `gradient` and `gradient.descent`
- Functions as return values
 - Example: Linear predictor
 - Example: the gradient operator
- Example: `surface`

READING: Sections 7.5, 7.11 and 7.13 of Matloff

OPTIONAL RECOMMENDED READING: Chapter 3 of Chambers

CODE FROM THIS LECTURE: At class website, with comments

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Both ideas can be understood from your experience with calculus

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Definite integral of f over $[a, b]$: takes a function and two points, gives a number

$$\int_a^b f(x) dx \equiv \lim_{n \rightarrow \infty} \sum_{i=0}^{n-1} \left(\frac{b-a}{n} \right) f \left(a + i \frac{b-a}{n} \right)$$

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∇ is an **operator**, not a functional

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Indefinite integration: $\int_{-\infty}^x f(u) du$ takes f and gives a new function

Fourier transform: takes f and gives a new function

$$\tilde{f}(\omega) = \int_{x=-\infty}^{x=\infty} f(x) e^{2i\pi\omega x} dx$$

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We have already seen an example: `apply` takes a function as one of its arguments

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gives argument list of `foo`: names are argument names, values are expressions for defaults (if any)
- `parent environment`: access with `environment(foo)`

User-defined and built-in R functions are both of class `function`

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Example: curve

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(because you did all of the assigned reading, including section 2.3.3 of the textbook)

A call to `curve` looks like this:

```
curve(expr, from = a, to = b, ...)
```

`expr` is some expression involving a variable called `x`

which is swept from the value `a` to the value `b`

`...` are other plot-control arguments

`curve` presumes that the expression can take a vector of `x` values and return a vector of numerical values, e.g.,

```
curve(x^2 * sin(x))
```

is fine

Using curve with your own functions

If we have defined a function already, we can use it in curve:

```
psi <- function(x,c=1) {ifelse(abs(x)>c,2*c*abs(x)-c^2,x^2)}  
curve(psi(x,c=10),from=-20,to=20)
```

Try this! Also try

```
curve(psi(x=10,c=x),from=-20,to=20)
```

and explain it to yourself

If our function doesn't take vectors to vectors, curve becomes unhappy

```
> mse <- function(y0,a,Y=gmp$pcgmp,N=gmp$pop) {  
+   mean((Y - y0*(N^a))^2)  
+ }  
> curve(mse(a=x,y0=6611),from=0.10,to=0.15)  
Error in curve(mse(a = x, y0 = 6611), from = 0.1, to = 0.15) :  
  'expr' did not evaluate to an object of length 'n'  
In addition: Warning message:  
In N^a : longer object length is not a multiple of shorter object length
```

How do we solve this?

apply applies the same function to every row or column of an array
sapply applies the same function to every element of an array or vector, and tries to simplify the result down to an array

```
> sapply(seq(from=0.10,to=0.15,by=0.01),mse,y0=6611)
[1] 154701953 102322975 68755655 64529167 104079528 207057513
> mse(6611,0.10)
[1] 154701953
```

Now (try it!):

```
mse.plottable <- function(a,...){ return(sapply(a,mse,...)) }
curve(mse.plottable(a=x),from=0.10,to=0.15)
curve(mse.plottable(a=x,y0=5100),from=0.10,to=0.20)
```

Later, we will see many more related tricks for splitting up problems and applying the same function repeatedly

Example: gradient

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find the partial derivative of f with respect to each component of x
return the vector of partial derivatives
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return the vector of partial derivatives
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It makes no sense to re-write this every time we change f !

\therefore write code to calculate the gradient of an arbitrary function

```
gradient <- function(f,x,deriv.steps) {  
  # not real code  
  evaluate the function at  $x$  and at  $x+\text{deriv.steps}$   
  take slopes to get partial derivatives  
  return the vector of partial derivatives  
}
```

A naive implementation would use a for loop

```
gradient <- function(f,x,deriv.steps,...) {  
  p <- length(x)  
  stopifnot(length(deriv.steps)==p)  
  f.old <- f(x,...)  
  gradient <- vector(length=p)  
  for (coordinate in 1:p) {  
    x.new <- x  
    x.new[coordinate] <- x.new[coordinate]+deriv.steps[coordinate]  
    f.new <- f(x.new,...)  
    gradient[coordinate] <- (f.new - f.old)/deriv.steps[coordinate]  
  }  
  return(gradient)  
}
```

Works, but it's so repetitive!

Better: use matrix manipulation and apply

```
gradient <- function(f,x,deriv.steps,...) {  
  p <- length(x)  
  stopifnot(length(deriv.steps)==p)  
  x.new <- matrix(rep(x,times=p),nrow=p) + diag(deriv.steps,nrow=p)  
  f.new <- apply(x.new,2,f,...)  
  gradient <- (f.new - f(x,...))/deriv.steps  
  return(gradient)  
}
```

(clearer, and half as long)

Presumes that `f` takes a vector and returns a single number

Any extra arguments to `gradient` will get passed to `f`

Check: Does this work when `f` is a function of a single number?

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If it really matters, use the `grad` function in the `numDeriv` package

Now we can use this as a piece of a larger machine:

```
gradient.descent <- function(f,x,max.iterations,step.scale,
  stopping.deriv,...) {
  for (iteration in 1:max.iterations) {
    grad <- gradient(f,x,...)
    if(all(abs(grad) < stopping.deriv)) { break() }
    x <- x - step.scale*grad
  }
  fit <- list(argmin=x,final.gradient=grad,final.value=f(x,...),
    iterations=iteration)
  return(fit)
}
```

(As written, we need to specify `deriv.steps` when calling this, but that's not an argument.

(How can you tell? Why make this choice?))

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Works equally well whether f is mean squared error of a regression, ψ error of a regression, (negative log) likelihood, cost of a production plan, ...

Wrappers and Anonymous Functions

`gradient.descent` presumes `f` takes a vector
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What to do?

- 1 Put a wrapper around `mse`:

```
mse.for.optimization <- function(param,...) {  
  return(mse(y0=param[1],a=param[2],...))  
}  
gradient.descent(f=mse.for.optimization, blah blah blah)
```

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- 2 Use an anonymous function:

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  a=param[2],...)},blah blah blah)
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(in fact the `f=` is optional here)

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  a=param[2],...)},blah blah blah)
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(in fact the `f=` is optional here)

Anonymous functions work because the return value of `function` is
a function object

Anonymous functions don't clutter your workspace, but they don't
stick around for you to examine later

Scoping `f` takes values for all names which aren't its arguments from the environment where it was defined, not the one where it is called (e.g., not from inside `gradient` or `gradient.descent`)

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Debugging If f and g are both complicated, avoid debugging $g(f)$ as a block; divide the work by writing *very simple* `f.0` to debug/test g , and debug/test the real f separately

Returning Functions: A trivial example

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```
make.noneuclidean <- function(ratio.to.diameter=pi) {  
  circumference <- function(d) { return(ratio.to.diameter*d) }  
  return(circumference)  
}
```

Define `make.noneuclidean` but don't run it yet

```
> circumference(10)  
Error: could not find function "circumference"  
> kings.i <- make.noneuclidean(3)  
> kings.i(10)  
[1] 30  
> formals(kings.i)  
$d  
> body(kings.i)  
{  
  return(ratio.to.diameter * d)  
}  
> environment(kings.i)  
<environment: 0xe43d64>  
> circumference(10)  
Error: could not find function "circumference"
```

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make.linear.predictor <- function(x,y) {  
  linear.fit <- lm(y~x)  
  predictor <- function(x) {  
    return(predict(object=linear.fit,newdata=data.frame(x=x)))  
  }  
  return(predictor)  
}
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    return(predict(object=linear.fit,newdata=data.frame(x=x)))  
  }  
  return(predictor)  
}
```

The predictor function persists and works, even when the data we used to create it is gone

```
> library(MASS); data(cats)
> vet_predictor <- make.linear.predictor(x=cats$Bwt,y=cats$Hwt)
> rm(cats)           # Data set goes away
> vet_predictor(4.0) # My cat's body mass in kilograms
      1
15.77959           # Predicted mass of my cat's heart in grams
```

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}
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}  
  
> mse.gradient <- nabla(mse.for.optimization)  
> mse.gradient(c(6611,0.15),deriv.steps=c(1,1e-6))  
[1] 1.646082e+05 1.428795e+10  
> gradient(mse.for.optimization,c(6611,0.15),c(1,1e-6))  
[1] 1.646082e+05 1.428795e+10  
> gradient(mse.for.optimization,c(6611,0.15),c(1,1e-6),Y=2*gmp$pcgmp)  
[1] -2.908638e+05 -2.486987e+10  
> mse.gradient(c(6611,0.15),deriv.steps=c(1,1e-6),Y=2*gmp$pcgmp)  
[1] -2.908638e+05 -2.486987e+10
```

The simple first-differences method is not so hot, so use the `grad` function from `numDeriv`

```
del <- function(f,...) {  
  require(numDeriv)  
  g <- function(x,...) { grad(func=f,x=x, ...)}  
  return(g)  
}
```

How would you check this?

Example: surface

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Built-in plotting function `contour`:

```
contour(x,y,z, [[other stuff]])
```

x and y are vectors of coordinates, z is a matrix of the corresponding shape

(see `help(contour)` for graphical options)

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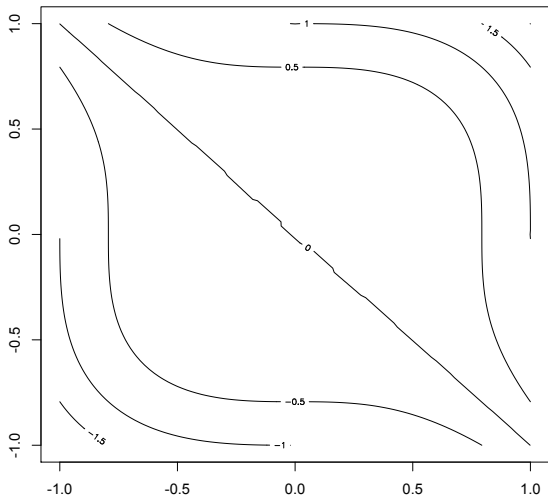
Strategy: `surface` should make x and y sequences, evaluate the expression at each combination to get z , and then call `contour`

First attempt at surface

Only works with vector-to-number functions:

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```
surface.0 <- function(f,from.x=0,to.x=1,from.y=0,to.y=1,n.x=101,  
  n.y=101,...) {  
  x.seq <- seq(from=from.x,to=to.x,length.out=n.x)  
  y.seq <- seq(from=from.y,to=to.y,length.out=n.y)  
  plot.grid <- expand.grid(x=x.seq,y=y.seq)  
  z.values <- apply(plot.grid,1,f)  
  z.matrix <- matrix(z.values,nrow=n.x)  
  contour(x=x.seq,y=y.seq,z=z.matrix,...)  
  invisible(list(x=x.seq,y=y.seq,z=z.matrix))  
}
```



```
surface.0(function(p){return(sum(p^3))},from.x=-1,from.y=-1)
```

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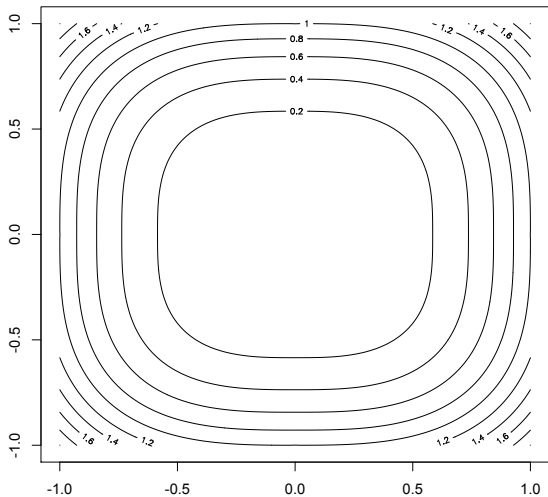
When we type something like `x^2+y^2` as an argument to `surface.0`, R tries to evaluate it prematurely

`substitute` returns the *unevaluated* expression

`curve` uses first `substitute(expr)` and then `eval(expr, envir)`, having made the right `envir`

Second attempt at surface

```
surface.1 <- function(expr,from.x=0,to.x=1,from.y=0,to.y=1,n.x=101,
  n.y=101,...) {
  x.seq <- seq(from=from.x,to=to.x,length.out=n.x)
  y.seq <- seq(from=from.y,to=to.y,length.out=n.y)
  plot.grid <- expand.grid(x=x.seq,y=y.seq)
  unevaluated.expression <- substitute(expr)
  z.values <- eval(unevaluated.expression,envir=plot.grid)
  z.matrix <- matrix(z.values,nrow=n.x)
  contour(x=x.seq,y=y.seq,z=z.matrix,...)
  invisible(list(x=x.seq,y=y.seq,z=z.matrix))
}
```



```
surface.1(abs(x^3)+abs(y^3),from.x=-1,from.y=-1)
```

Evaluating a function at every combination of two arguments is a really common task

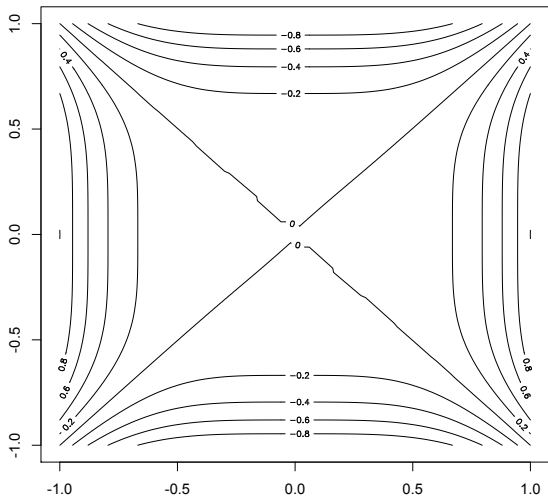
There is a function to do it for us: `outer` (seen in lecture 3)

Evaluating a function at every combination of two arguments is a really common task

There is a function to do it for us: `outer` (seen in lecture 3)

```
surface.2 <- function(expr,from.x=0,to.x=1,from.y=0,to.y=1,n.x=101,
  n.y=101,...) {
  x.seq <- seq(from=from.x,to=to.x,length.out=n.x)
  y.seq <- seq(from=from.y,to=to.y,length.out=n.y)
  unevaluated.expression <- substitute(expr)
  z <- function(x,y) {
    return(eval(unevaluated.expression,envir=list(x=x,y=y)))
  }
  z.values <- outer(X=x.seq,Y=y.seq,FUN=z)
  z.matrix <- matrix(z.values,nrow=n.x)
  contour(x=x.seq,y=y.seq,z=z.matrix,...)
  invisible(list(x=x.seq,y=y.seq,z=z.matrix))
}
```

could also include the function as part of the returned list



`surface.2(x^4-y^4,from.x=-1,from.y=-1)`

- In R, functions are objects, and can be arguments to other functions
 - Use `lapply` to do the same thing to many different functions
 - Separates writing the high-level operations and the first-order functions
 - Use `sapply` (etc.), wrappers, anonymous functions as adapters
- Functions can also be returned by other functions
 - Variables other than the arguments to the function are fixed by the environment of creation
 - Manipulating expressions lets us flexibly create functions