# Week 2: Homework Assignment

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Saturday, January 24, 2015

# Code Your Own Optimizer in R

This project helps understanding how non-linear optimization of log-likelihood function by the Newton-Raphson method works

Use the formulas for the steps of Newton-Raphson method from the lecture notes to code a simple version of optimizer.

## **Assignment description**

1. Create a test function that needs to be optimized.

For this project we use one-dimensional optimization, i.e. optimization with respect to only one variable.

Let the declaration of the function be my. Function<-function(my.X), where my.X is a scalar parameter with respect to which the optimization is done.

The function should cross the x-axis at least in one point.

For example, you can use

```
my.Function<-function(my.X) {</pre>
  my.X^2*3-my.X*5-5
```

#### 2. Write your optimizer.

Let the declaration of the optimizer function be <a href="my.Optimizer-function">my.Optimizer-function</a>(Start.Value, Function.To.Optimize, Epsilon)
where <a href="Start.Value">Start.Value</a> is the initial guess for the optimizer, <a href="Function.To.Optimize">Function.To.Optimize</a> is the name of your test function that needs to be optimized, <a href="Epsilon">Epsilon</a> is the stopping criterion, a small number, for example, 0.001.

The function [my.Optimizer] should repeat iterations of the Newton-Raphson algorithm while |x(i+1) - x(i)| > = Epsilon, where x(i+1) is the approximation obtained during the recent iteration and x(i) is the approximation obtained during the previous operation. You can use any of the loops in R, for example,  $[mathbb{while}(cond)]$  where  $[mathbb{cond}]$  is the condition of moving to the next iteration (|x(i+1) - x(i)| > = ??) and  $[mathbb{expr}]$  is the sequence of the commands that need to be performed at each iteration.

```
my.Optimizer<-function(Start.Value, Function.To.Optimize, Epsilon){
  #initialize check.done so that it does not satisfy while condition below
  check.done <- 1;</pre>
  #initialize 'old.value' parameter
  old.value <- Start.Value;</pre>
  #initialize the counter that determines the number of times the while loop executes
  iter <- 0
  #set an 'h' value for the derivative estimation
  h <- .00000001
  while(check.done > Epsilon){
    new.value <- old.value - (Function.To.Optimize(old.value)/</pre>
```

```
((Function.To.Optimize(old.value+h)-Function.To.Optimize(old.value))/(h)))
  check.done <- abs(new.value - old.value)</pre>
  old.value <- new.value
  iter <- iter+1
return(old.value)
```

### 3. Test the optimizer

Use your optimizer with the test function. For example, use my.Optimizer(-5, my.Function, .001). Make sure you calculate the answer manually to check the answer.

You can also test the optimizer by running uniroot(). For example, uniroot(my.Function, lower=-5, upper=+1)

The root returned by your optimizer should be the same as the output **\$root** of the object returned by **uniroot()**.

```
#first obtain the output from the optimizer that I created
my.Optimizer(-5, my.Function, .0001)
```

```
## [1] -0.7032574
```

```
#now check this against the output out uniroot()
uniroot(my.Function,lower=-5,upper=1)
```

```
## $root
## [1] -0.703257
##
## $f.root
## [1] -3.368345e-06
##
## $iter
## [1] 9
##
## $estim.prec
## [1] 6.103516e-05
```

As you can observe from the output comparisons above, my optimization function uniroot() produce the same results!

Try also to run <code>optim()</code>. Explain the difference between the two functions: <code>uniroot()</code> and <code>optim()</code>.

```
optim(-5,
    fn = my.Function,
    method="L-BFGS-B",
    hessian=TRUE,
    lower=c(-Inf,0))
```

```
## $par
## [1] 0.8333333
```

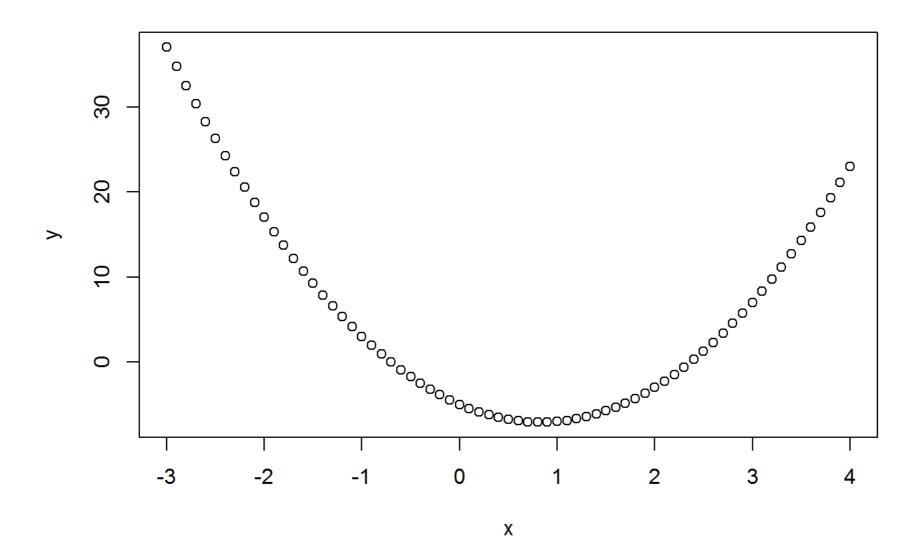
```
##
## $value
## [1] -7.083333
##
## $counts
## function gradient
         10
##
                  10
##
## $convergence
## [1] 0
##
## $message
## [1] "CONVERGENCE: REL_REDUCTION_OF_F <= FACTR*EPSMCH"
##
## $hessian
##
        [,1]
## [1,]
        6
```

### How can we reconsile their outputs?

As we see above, the <code>optim()</code> output is different than the output produced by <code>my.Optimizer()</code> and <code>uniroot()</code>.

If you observe the plot of the original function (below) you notice that, as expected, my my.Optimizer() and uniroot() identify a root of the function, while optim() identifies the minimum of the function (0.8333).

```
#show a plot of the function
x <- seq(from=-3, to=4, by=.1)
y <- my.Function(x)
plot(x,y)</pre>
```



However, our different prodcedures above are related in that if you are optimizing (finding the minimum or maximum) a function then you can either use optim() to accomplish this OR take the deravative of the function and use either

my.Optimizer or uniroot() to find the root of the derivation which is equal to the max/min of the original function.

Similarly, if you wanted to find the roots of a function you can either use <a href="my.0ptimizer">my.0ptimizer</a> or <a href="uniroot()">uniroot()</a> to accomplish this or take intergral of the function and use <a>optim()</a> to find min / max points of the intergral which is equal to the roots of the orginal function.

These findings align with our lecture and learnins in class.