CompStatLab2

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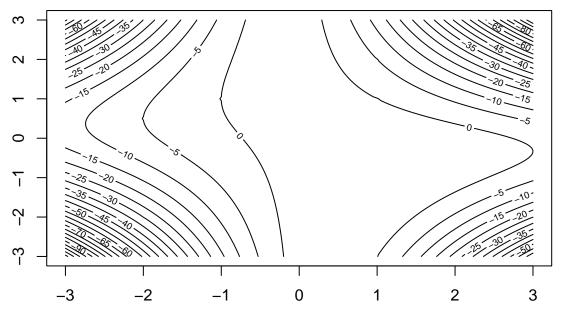
Computational Statistics 732A90 | Computer Lab 1

Question 1: Optimisation of a two-dimensional function

Consider the function $g(x,y) = -x^2 - x^2y^2 - 2xy + 2x + 2$ It is desired to determine the point (x, y), (x, y

a. Derive the gradient and the Hessian matrix in dependence of x, y. Produce a contour plot of the function g.

Answer: Please reference Appendix two_dimensional.R, and below are plot of function g.



b. Write an own algorithm based on the Newton method in order to find a local maximum of g

Answer:

Please reference Appendix two_dimensional.R

c. Use different starting values: use the three points (x, y) = (2, 0), (-1, -2), (0, 1) and a fourth point of your choice. Describe what happens when you run your algorithm for each of those starting values. If your algorithm converges to points (x, y), compute the gradient and the Hessian matrix at these points and decide about local maximum, minimum, saddle point, or neither of it. Did you find a global maximum for x, y = [-3, 3]?

Answer:

Please see below R results:

```
## 1) using starting points: (x,y) = (2,0), the newton method coverges to point:
##
## [1,] 1.0000256
## [2,] -0.9999341
## the corresponding gradient is:
## [1] 2.926312e-05 -8.049134e-05
## the corresponding hessian matrix is:
          [,1]
                   [,2]
## [1,] -3.999737 1.999839
## [2,] 1.999839 -2.000102
## This Hessian Matrix is negative definite.
## 2) using starting points: (x,y) = (-1,-2)
## [1,] 1.166791e-11
## [2,] 1.000000e+00
## the corresponding gradient is:
## [1] -2.939771e-08 -2.333581e-11
## the corresponding hessian matrix is:
      [,1]
## [1,] -4 -2.0000e+00
## [2,] -2 -2.7228e-22
## This Hessian Matrix is neither positive nor negative definite.
## 3) using starting points: (x,y) = (0,1)
     [,1]
## [1,]
## [2,]
         1
## the corresponding gradient is:
## [1] 0 0
## the corresponding hessian matrix is:
      [,1] [,2]
##
## [1,]
      -4 -2
## [2,]
      -2
## This Hessian Matrix is neither positive nor negative definite.
## 4) using starting points: (x,y) = (1,-1)
     [,1]
## [1,]
## [2,] -1
```

```
## the corresponding gradient is:
## [1] 0 0
## the corresponding hessian matrix is:
## [,1] [,2]
## [1,] -4 2
## [2,] 2 -2
## This Hessian Matrix is possible definite
```

This Hessian Matrix is negative definite.

According to above R results for 4 different starting points, it's easily to get the conclusion:

- 1) when starting point is (2,0), the coveraged point is (1.0000256, -0.9999341) which is local maximum point since the Hessian matrix at that point is negative definite.
- 2) when starting point is (-1,-2), the coveraged point is () which is saddle point since the Hessian matrix at that point is neither positive or negative definite.
- 3) when starting point is (0,1), the coveraged point is (0,1) which is saddle point since the Hessian matrix at that point is neither positive or negative definite.
- 4) when starting point is (1,-1), the coveraged point is (1,-1) which is local maximum since the Hessian matrix at that point is negative definite.

d. What would be the advantages and disadvantages when you would run a steepest ascent algorithm instead of the Newton algorithm?

Answer:

Question 2

a. Write a function for an ML-estimator for (0, 1) using the steepest ascent method with a step-size reducing line search (back-tracking). For this, you can use and modify the code for the steepest ascent example from the lecture. The function should count the number of function and gradient evaluations.

Answer:

b. Compute the ML-estimator with the function from a. for the data (xi, yi) above. Use a stopping criterion such that you can trust five digits of both parameter estimates for 0 and 1. Use the starting value (0, 1) = (-0.2, 1). The exact way to use backtracking can be varied. Try two variants and compare number of function and gradient evaluation done until convergence.

Answer:

c. Use now the function optim with both the BFGS and the Nelder-Mead algorithm. Do you obtain the same results compared with b.? Is there any difference in the precision of the result? Compare the number of function and gradient evaluations which are given in the standard output of optim.

Answer:

d. Use the function glm in R to obtain an ML-solution and compare it with your results before.

Answer:

Appendix:

 $two_dimensional.R$

```
# function g to be maximized
g <- function(x,y) {</pre>
  return(-x^2 - x^2 * y^2 - 2 * x * y + 2 * x + 2)
}
# partial derivative x for function g
deriv_x <- function(x, y) {</pre>
  return(-2 * x - 2 * x * y^2 - 2 * y + 2)
}
# partial derivative y for function g
deriv_y <- function(x, y) {</pre>
  return(-2 * x^2 * y - 2 * x)
# gradient for function g
gradient <- function (x, y) {</pre>
  return(c(deriv_x(x,y), deriv_y(x,y)))
}
# second partial derivative x for function q
deriv_x_11 <- function(x, y) {</pre>
  return(-2 - 2 * y^2)
# second partial derivative x/y for function g
deriv_x_12 <- function(x, y) {</pre>
  return(-4 * x * y - 2)
# second partial derivative y for function g
deriv_y_22 <- function(x, y) {</pre>
  return(-2 * x^2)
# hessian matrix for function g
hessian_g <- function(x, y) {
  return(matrix(c(deriv_x_11(x,y), deriv_x_12(x,y), deriv_x_12(x,y), deriv_y_22(x,y)), nrow = 2, ncol =
# produce a contour plot
xgrid \leftarrow seq(-3,3, by=0.05)
ygrid \leftarrow seq(-3,3, by=0.05)
length_x <- length(xgrid)</pre>
length_y <- length(ygrid)</pre>
dxy <- length_x * length_y</pre>
gxy <- matrix(rep(NA,dxy), nrow = length_x)</pre>
for ( i in 1:length_x) {
 for ( j in 1:length_y) {
    gxy[i, j] <- g(xgrid[i], ygrid[j])</pre>
  }
}
```

```
mgxy <- matrix(gxy, nrow = length_x, ncol = length_y)</pre>
contour(xgrid, ygrid, mgxy, nlevels=40)
# newton function
newton_g <- function(x,y, eps=0.0001) {</pre>
 xt \leftarrow c(x,y)
 xt1 < -c(x,y) + 2
 while( t(xt1 - xt) \% \% (xt1 - xt) > eps) {
   xt1 <- xt
   xt <- xt1 - solve(hessian_g(xt[1] , xt[2])) %*% gradient(xt[1],xt[2])</pre>
 return(xt)
}
# check definiteness of a hessian matrix
check_definiteness <- function(matrix) {</pre>
 eigenvalues <- eigen(matrix)$values</pre>
 if (all(eigenvalues > 0)) {
   cat("This Hessian Matrix is positive definite.\n")
 } else if (all(eigenvalues < 0)) {</pre>
   cat("This Hessian Matrix is negative definite.\n")
 } else {
   cat("This Hessian Matrix is neither positive nor negative definite.\n")
 }
}
# using different starting points for newton method
converges_point1 <- newton_g(2,0)</pre>
cat("1) using starting points: (x,y) = (2,0), the newton method coverges to point: \n")
converges_point1
grad_point1 <- gradient(converges_point1[1],converges_point1[2])</pre>
cat("the corresponding gradient is: \n")
grad_point1
hessian_point1 <- hessian_g(converges_point1[1],converges_point1[2])
cat("the corresponding hessian matrix is: \n")
hessian_point1
check_definiteness(hessian_point1)
cat("-----")
converges_point2 <- newton_g(-1,-2)</pre>
cat("2) using starting points: (x,y) = (-1,-2) \n"
converges_point2
grad_point2 <- gradient(converges_point2[1],converges_point2[2])</pre>
cat("the corresponding gradient is: \n")
grad_point2
hessian_point2 <- hessian_g(converges_point2[1],converges_point2[2])
```

```
cat("the corresponding hessian matrix is: \n")
hessian_point2
check_definiteness(hessian_point2)
converges_point3 <- newton_g(0,1)</pre>
cat("3) using starting points: (x,y) = (0,1) \ n")
converges_point3
grad_point3 <- gradient(converges_point3[1],converges_point3[2])</pre>
cat("the corresponding gradient is: \n")
grad_point3
hessian_point3 <- hessian_g(converges_point3[1],converges_point3[2])</pre>
cat("the corresponding hessian matrix is: \n")
hessian_point3
check_definiteness(hessian_point3)
converges_point4 <- newton_g(1,-1)</pre>
cat("4) using starting points: (x,y) = (1,-1) \ n")
converges_point4
grad_point4 <- gradient(converges_point4[1],converges_point4[2])</pre>
cat("the corresponding gradient is: \n")
grad_point4
hessian_point4 <- hessian_g(converges_point4[1],converges_point4[2])
cat("the corresponding hessian matrix is: \n")
hessian_point4
check_definiteness(hessian_point4)
g <- function(x,y) {
```