

# STAT 579 Homework 7

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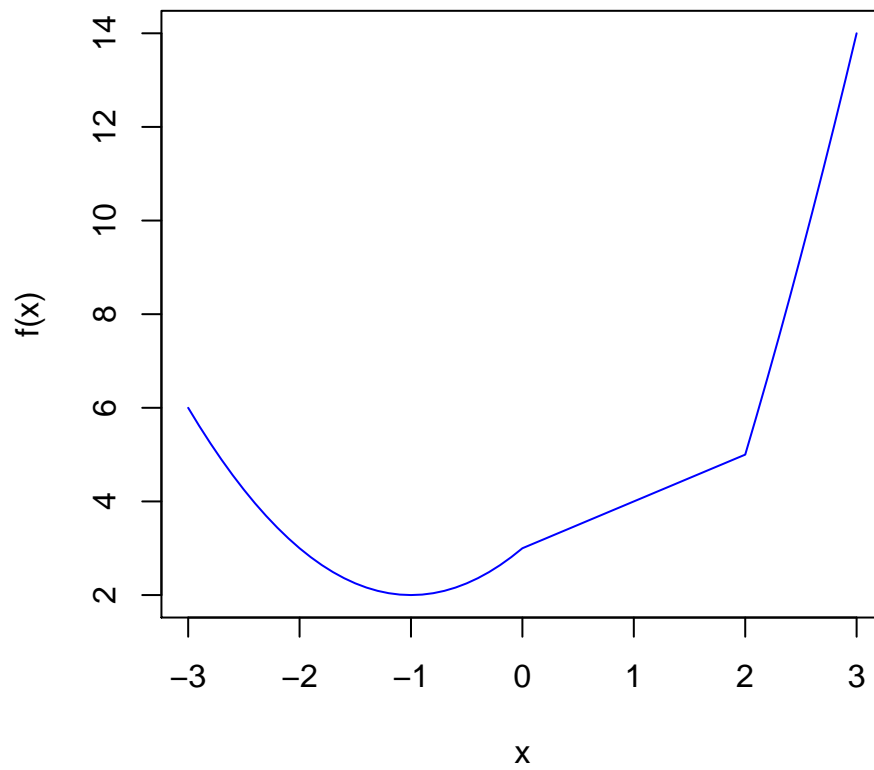
## Problem 1

```
tridiag <- function(k, n) {  
  A <- matrix(rep(0, n * n), ncol = n)  
  diag(A) <- rep(k, n)  
  diag(A[-1, -n]) <- rep(1, n - 1)  
  diag(A[-n, -1]) <- rep(1, n - 1)  
  return(A)  
}  
  
tridiag(n = 6, k = 5)
```

```
##      [,1] [,2] [,3] [,4] [,5] [,6]  
## [1,]    5    1    0    0    0    0  
## [2,]    1    5    1    0    0    0  
## [3,]    0    1    5    1    0    0  
## [4,]    0    0    1    5    1    0  
## [5,]    0    0    0    1    5    1  
## [6,]    0    0    0    0    1    5
```

## Problem 2

```
tmpFn <- function(xVec) {  
  x <- xVec  
  ifelse(x < 0, x^2 + 2 * x + 3, ifelse(x < 2, x + 3, x^2 + 4 * x - 7))  
}  
x <- seq(-3, 3, 0.1)  
fx <- tmpFn(x)  
  
plot(x = x, y = fx, "l", col = "blue", xlab = "x", ylab = "f(x)")
```



### Problem 3

```
gcd <- function(m, n) {  
  re <- m%%n  
  while (re != 0) {  
    m <- n  
    n <- re  
    re <- m%%n  
  }  
  return(n)  
}  
  
gcd(60, 130)
```

```
## [1] 10
```

#### Problem 4

```
order.matrix <- function(x) {  
  dimx <- dim(x)  
  ord <- order(x)  
  n <- length(ord)  
  dim(ord) <- dimx  
  indices <- NULL  
  for (i in 1:n) {  
    indice <- which(ord == i, arr.ind = T)  
    indices <- rbind(indices, indice)  
  }  
  indices <- cbind(1:n, indices)  
  colnames(indices) <- c("order", "row", "column")  
  return(indices)  
}  
  
x <- matrix(rchisq(n = 4 * 3, df = 1), ncol = 3)  
  
order.matrix(x)
```

```
##      order row column  
## [1,]      1  2      2  
## [2,]      2  3      1  
## [3,]      3  3      3  
## [4,]      4  2      1  
## [5,]      5  3      2  
## [6,]      6  1      1  
## [7,]      7  4      3  
## [8,]      8  1      2  
## [9,]      9  4      1  
## [10,]     10  1      3  
## [11,]     11  4      2  
## [12,]     12  2      3
```

## Problem 5

(a)

```
polaroid <- function(x) {
  p <- length(x)
  q <- rep(0, p - 1)
  R <- sqrt(sum(x^2))
  q[p - 1] <- ifelse(atan(x[p]/x[p - 1]) >= 0, atan(x[p]/x[p - 1]), atan(x[p]/x[p - 1]) + pi)
  for (i in (p - 2):2) {
    arctan <- atan((x[i + 1]/cos(q[i + 1]))/x[i])
    q[i] <- ifelse(arctan >= 0, arctan, arctan + pi)
  }
  q[1] <- ifelse(x[2]/cos(q[2]) >= 0, atan2(y = x[2]/cos(q[2]), x = x[1]), atan2(y = x[2]/cos(q[2]), x = x[1]) + 2 * pi)

  return(c(R, q))
}
```

(b)

```
normalize <- function(X) {
  norm <- sqrt(apply(X^2, MARGIN = 1, FUN = sum))
  sweep(X, MARGIN = 1, STAT = norm, FUN = "/")
}
```

(c)

```
y <- matrix(rnorm(n = 1000 * 5), ncol = 5)

z <- normalize(y)

kstest <- apply(z, MARGIN = 2, FUN = ks.test, "punif", min = -1, max = 1)

kstest
```

```
## [[1]]
##
## One-sample Kolmogorov-Smirnov test
##
## data:  newX[, i]
## D = 0.10213, p-value = 1.742e-09
## alternative hypothesis: two-sided
##
##
## [[2]]
##
## One-sample Kolmogorov-Smirnov test
##
## data:  newX[, i]
## D = 0.11491, p-value = 6.792e-12
```

```
## alternative hypothesis: two-sided
##
##
## [[3]]
##
## One-sample Kolmogorov-Smirnov test
##
## data:  newX[, i]
## D = 0.11309, p-value = 1.559e-11
## alternative hypothesis: two-sided
##
##
## [[4]]
##
## One-sample Kolmogorov-Smirnov test
##
## data:  newX[, i]
## D = 0.10606, p-value = 3.397e-10
## alternative hypothesis: two-sided
##
##
## [[5]]
##
## One-sample Kolmogorov-Smirnov test
##
## data:  newX[, i]
## D = 0.092739, p-value = 6.773e-08
## alternative hypothesis: two-sided
```

P-values are small, so they should not follow the uniform distribution on  $(-1, 1)$ .

(d)

```
polar <- apply(y, MARGIN = 1, FUN = polaroid)

R2 <- (polar[1, ])^2

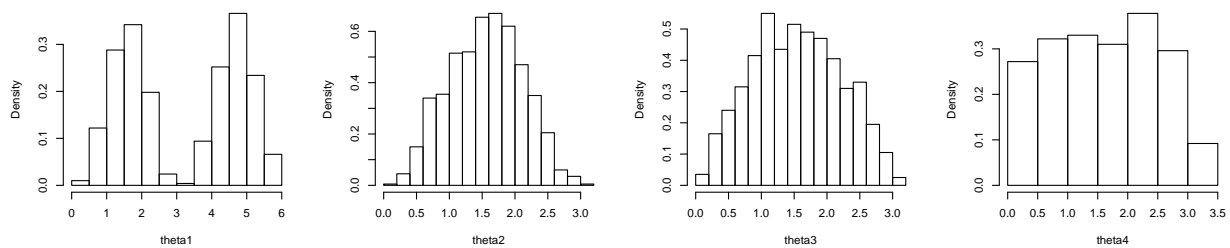
ks.test(R2, "pchisq", df = 5)
```

```
##
## One-sample Kolmogorov-Smirnov test
##
## data:  R2
## D = 0.025898, p-value = 0.5136
## alternative hypothesis: two-sided
```

```
theta <- polar[-1, ]

par(mfrow = c(1, 4))

hist(theta[1, ], xlab = "theta1", freq = F, main = "")
hist(theta[2, ], xlab = "theta2", freq = F, main = "")
hist(theta[3, ], xlab = "theta3", freq = F, main = "")
hist(theta[4, ], xlab = "theta4", freq = F, main = "")
```



```
ks.test(theta[1, ], "punif", min = 0, max = 2 * pi)
```

```
##
## One-sample Kolmogorov-Smirnov test
##
## data: theta[1, ]
## D = 0.1087, p-value = 1.089e-10
## alternative hypothesis: two-sided
```

```
ks.test(theta[2, ], "punif", min = 0, max = pi)
```

```
##
## One-sample Kolmogorov-Smirnov test
##
## data: theta[2, ]
## D = 0.18376, p-value < 2.2e-16
## alternative hypothesis: two-sided
```

```
ks.test(theta[3, ], "punif", min = 0, max = pi)
```

```
##
## One-sample Kolmogorov-Smirnov test
##
## data: theta[3, ]
## D = 0.1122, p-value = 2.329e-11
## alternative hypothesis: two-sided
```

```
ks.test(theta[4, ], "punif", min = 0, max = pi)
```

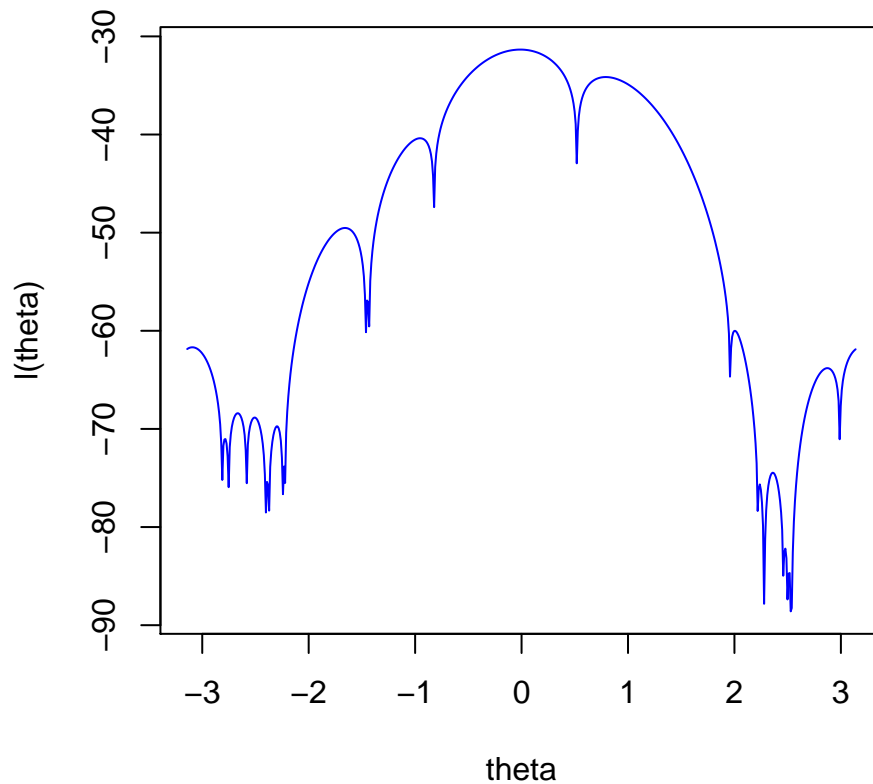
```
##
## One-sample Kolmogorov-Smirnov test
##
## data: theta[4, ]
## D = 0.034434, p-value = 0.1865
## alternative hypothesis: two-sided
```

P-value for test for  $R^2$  is large, so it should follow the  $\chi^2_5$ . P-value for  $\theta_1, \theta_2, \theta_3$  are small, so they should not follow uniform distribution. P-value for  $\theta_4$  is large, thus it should follow the uniform distribution on  $[0, \pi)$ .

## Problem 6

(a)

```
x <- c(3.91, 4.85, 2.28, 4.06, 3.7, 4.04, 5.46, 3.53, 2.28, 1.96, 2.53, 3.88,  
      2.22, 3.47, 4.82, 2.46, 2.99, 2.54, 0.52, 2.5)  
  
l <- function(theta, x) {  
  n <- length(x)  
  -n * log(2 * pi) + sum(log(1 - cos(x - theta)))  
}  
  
theta <- seq(-pi, pi, 0.01)  
  
ltheta <- apply(t(theta), MARGIN = 2, FUN = l, x = x)  
  
plot(x = theta, y = ltheta, "l", col = "blue", xlab = "theta", ylab = "l(theta)",  
     main = "")
```



(b)



```
thetamin <- optimize(f = l, x = x, interval = c(-pi, pi), maximum = T)
```

```
thetamin
```

```
## $maximum
## [1] -0.0119724
##
## $objective
## [1] -31.34291
```

(c)

```
options(digits = 20)
```

```
newton <- function(fun, derf, x0, eps) {
  iter <- 0
  repeat {
    iter <- iter + 1
    x1 <- x0 - fun(x0)/derf(x0)
    if (abs(x0 - x1) < eps || abs(fun(x1)) < 1e-10)
      break
    x0 <- x1
    cat("***** Iter. No: ", iter, " Current Iterate = ", x1, fill = T)
  }
  return(x1)
}
```

```
derl <- function(theta) {
  sum(sin(x - theta)/(1 - cos(x - theta)))
}
```

```
dderl <- function(theta) {
  -sum(1/(1 - cos(x - theta)))
}
```

```
newton(fun = derl, derf = dderl, x0 = 0, eps = 1e-04)
```

```
## ***** Iter. No: 1 Current Iterate = 0.011911932605691015
## ***** Iter. No: 2 Current Iterate = 0.035538787271073638
## ***** Iter. No: 3 Current Iterate = 0.081887140686978999
## ***** Iter. No: 4 Current Iterate = 0.17003691781937486
## ***** Iter. No: 5 Current Iterate = 0.32130252264058118
## ***** Iter. No: 6 Current Iterate = 0.49222040572661963
## ***** Iter. No: 7 Current Iterate = 0.52107506423276551
## ***** Iter. No: 8 Current Iterate = 0.52000207000084642
```

```
## [1] 0.520000000000761182
```

(d)

```
newton(fun = der1, derf = dder1, x0 = -2, eps = 1e-04)
```

```
## ***** Iter. No: 1 Current Iterate = -2.243846446394639
```

```
## ***** Iter. No: 2 Current Iterate = -2.24317446162916
```

```
## [1] -2.2431853038615284
```

```
newton(fun = der1, derf = dder1, x0 = -2.7, eps = 1e-04)
```

```
## ***** Iter. No: 1 Current Iterate = -2.7258863441687975
```

```
## ***** Iter. No: 2 Current Iterate = -2.7497175268270801
```

```
## ***** Iter. No: 3 Current Iterate = -2.75325274417618
```

```
## [1] -2.753185343300327
```

The results are different with different initial point. It is because Newton method will end with it finds a point with zero value which is nearest to the starting point. And the function has many points when its derivative is 0.