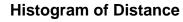
# STAT602 HW1

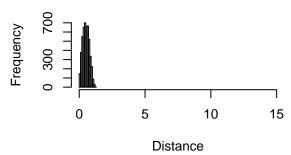
#### 1.6

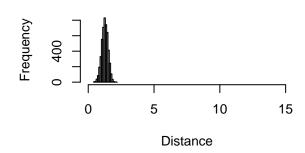
As p becomes larger, the data points become sparse. Suppose we are going to use knn, then the nearest neighbors may not be close to the point that we want to make prediction. So it may not work well when we have high dimensions.

```
n <- 100
p <- c(2, 10, 100, 1000)
old<-par(no.readonly = TRUE)</pre>
par(mfrow = c(2,2))
for(i in 1:4){
  d <- p[i]
  data <- matrix(runif(d*n), nrow = n, ncol = d)</pre>
  dis <- numeric(choose(100,2))</pre>
  cursor <- 1
  for(j in 1:(n-1))
    for(k in 2:n){
      if(k>j){
        dis[cursor] \leftarrow sum((data[j,] - data[k,])^2)^(1/2)
        cursor = cursor + 1
      }
    }
  hist(dis, xlim = c(0, 15), xlab = "Distance", main = "Histogram of Distance")
```



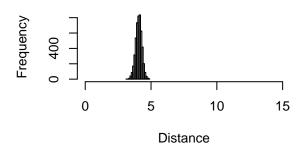
### **Histogram of Distance**

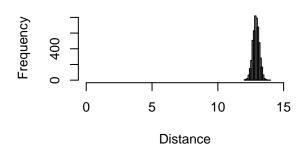




## **Histogram of Distance**

### **Histogram of Distance**





par(old)

#### 1.7

We can see that the last two combinations are too large that it is out of bound.

```
epsilon <- c(1,.1,.01)
p <- c(20, 50, 200)

i=1;j=1
v = pi^(p[j]/2)/gamma(p[j]/2+1)*epsilon[i]^p[j]
(n = 1/v)

## [1] 38.74934

i=2;j=1
v = pi^(p[j]/2)/gamma(p[j]/2+1)*epsilon[i]^p[j]
(n = 1/v)

## [1] 3.874934e+21

i=3;j=1
v = pi^(p[j]/2)/gamma(p[j]/2+1)*epsilon[i]^p[j]
(n = 1/v)</pre>
```

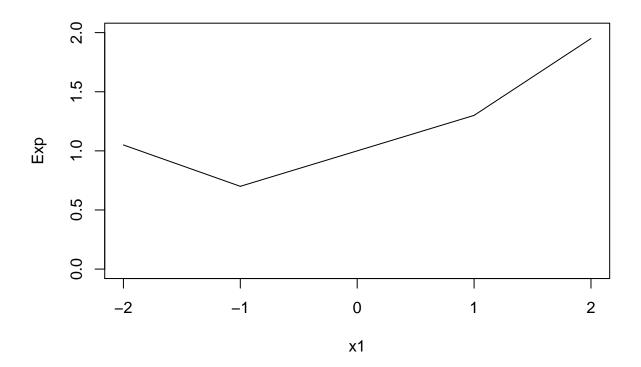
## [1] 3.874934e+41

```
i=1; j=2
v = pi^(p[j]/2)/gamma(p[j]/2+1)*epsilon[i]^p[j]
(n = 1/v)
## [1] 5.779614e+12
v = pi^(p[j]/2)/gamma(p[j]/2+1)*epsilon[i]^p[j]
(n = 1/v)
## [1] 5.779614e+62
i=3; j=2
v = pi^(p[j]/2)/gamma(p[j]/2+1)*epsilon[i]^p[j]
(n = 1/v)
## [1] 5.779614e+112
i=1; j=3
v = pi^(p[j]/2)/gamma(p[j]/2+1)*epsilon[i]^p[j]
(n = 1/v)
## [1] 1.798939e+108
i=2; j=3
v = pi^(p[j]/2)/gamma(p[j]/2+1)*epsilon[i]^p[j]
(n = 1/v)
## [1] Inf
i=3; j=3
v = pi^(p[j]/2)/gamma(p[j]/2+1)*epsilon[i]^p[j]
(n = 1/v)
## [1] Inf
```

#### 2.4

```
p \leftarrow c(0.35, 0.65)
E <- function(p, y){</pre>
  value <-p*max(0,1-y)+(1-p)*max(1+y,0)
  return(value)
}
x1 < - seq(-1,1,by = 0.01)
Exp <- numeric(length(x1))</pre>
for(i in 1:length(x1)){
  Exp[i] \leftarrow E(p[1], x1[i])
}
plot(x1, Exp, ty = "l", xlim = c(-2,2), ylim = c(0,2))
x1 < - seq(1,2,by = 0.01)
Exp <- numeric(length(x1))</pre>
for(i in 1:length(x1)){
  Exp[i] \leftarrow E(p[1], x1[i])
lines(x1, Exp)
```

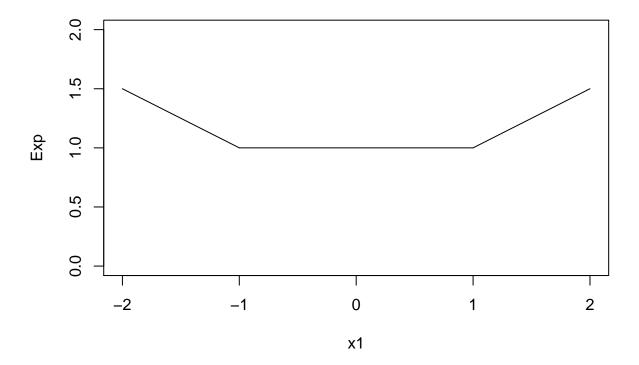
```
x1 <- seq(-2,-1,by = 0.01)
Exp <- numeric(length(x1))
for(i in 1:length(x1)){
   Exp[i] <- E(p[1], x1[i])
}
lines(x1, Exp)</pre>
```



```
p <- c(0.5,0.35)
E <- function(p, y){
   value <- p*max(0,1-y)+(1-p)*max(1+y,0)
   return(value)
}
x1 <- seq(-1,1,by = 0.01)
Exp <- numeric(length(x1))
for(i in 1:length(x1)){
   Exp[i] <- E(p[i], x1[i])
}
plot(x1, Exp, ty = "l", xlim = c(-2,2), ylim = c(0,2))

x1 <- seq(1,2,by = 0.01)
Exp <- numeric(length(x1))
for(i in 1:length(x1)){
   Exp[i] <- E(p[i], x1[i])
}
lines(x1, Exp)</pre>
```

```
x1 <- seq(-2,-1,by = 0.01)
Exp <- numeric(length(x1))
for(i in 1:length(x1)){
   Exp[i] <- E(p[1], x1[i])
}
lines(x1, Exp)</pre>
```



#### 3.4

 $\{1, sin(x), cos(x)\}\$  is the best. If the data is generated by factors that is not included in the last model, then none of these model is without model bias. If the data is generated with one of these model, then including that one, the models that include the true model are without the model bias.

```
data <- read_excel("~/Downloads/Problem3.4Data.xlsx")

K = 10

SSE = numeric(9)
for(i in 1:K)
{
   test = data[(10*(i-1)+1):(10*i),]
   train = data[-((10*(i-1)+1):(10*i)),]

# Polynomial
## 0
fp0 <- lm(data = train, y~1)</pre>
```

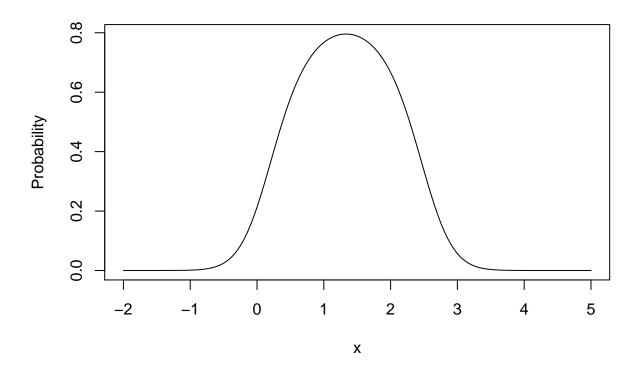
```
yp0 <- predict(fp0, test)</pre>
  SSE[1] = SSE[1] + sum((yp0-test\$y)^2)
  ## 1
  fp1 <- lm(data = train, y~x)</pre>
  yp1 <- fp1$coefficients %*% t(as.matrix(data.frame(inter = 1, x = test$x)))</pre>
  SSE[2] = SSE[2] + sum((yp1-test\$y)^2)
  ## 2
  fp2 \leftarrow lm(data = train, y~x+I(x^2))
  yp2 <- fp2$coefficients %*% t(data.frame(inter = 1, x = test$x, x2 = test$x^2))
  SSE[3] = SSE[3] + sum((yp2-test\$y)^2)
  ## 3
  fp3 <- lm(data = train, y~x+I(x^2)+I(x^3))
  yp3 <- fp3$coefficients %*% t(data.frame(inter = 1,x = test$x, x2 = test$x^2, x3 = test$x^3))
  SSE[4] = SSE[4] + sum((yp3-test\$y)^2)
  ## 4
  fp4 \leftarrow lm(data = train, y~x+I(x^2)+I(x^3)+I(x^4))
  yp4 <- fp4$coefficients %*% t(data.frame(inter = 1, x = test$x, x2 = test$x^2, x3 = test$x^3, x4 = te
  SSE[5] = SSE[5] + sum((yp4-test\$y)^2)
  ## 5
  fp5 <- lm(data = train, y~x+I(x^2)+I(x^3)+I(x^4)+I(x^5))
  yp5 <- fp5$coefficients %*% t(data.frame(inter =1,x = test$x, x2 = test$x^2, x3 = test$x^3, x4 = test
  SSE[6] = SSE[6] + sum((yp5-test\$y)^2)
  ## 6
  fp6 <- lm(data = train, y~1+I(sin(x))+I(cos(x)))
  yp6 <- fp6$coefficients %*% t(as.matrix(data.frame(inter = 1,sinx = sin(test$x), cosx = cos(test$x)))</pre>
  SSE[7] = SSE[7] + sum((yp6-test\$y)^2)
  ## 7
  fp7 <- lm(data = train, y^1+l(sin(x))+l(cos(x))+l(sin(2*x))+l(cos(2*x)))
  yp7 <- fp7$coefficients %*% t(as.matrix(data.frame(inter = 1,sinx = sin(test$x), cosx = cos(test$x),</pre>
  SSE[8] = SSE[8] + sum((yp7-test\$y)^2)
 ## 8
  fp8 <- lm(data = train, y~1+x+I(x^2)+I(x^3)+I(x^4)+I(x^5)+
              I(\sin(x))+I(\cos(x))+I(\sin(2*x))+I(\cos(2*x)))
 yp8 <- fp8$coefficients %*% t(data.frame(inter = 1 ,x = test$x, x2 = test$x^2,
                                  x3 = test$x^3, x4 = test$x^4,
                                  x5 = test$x^5, sinx = sin(test$x),
                                  cosx = cos(test\$x), sin2x = sin(2*test\$x),
                                  cos2x = cos(2*test$x)))
 SSE[9] = SSE[9] + sum((yp8-test\$y)^2)
}
SSE
## [1] 207.8989 212.4108 206.4061 190.6987 187.3633 193.0590 178.7845 184.4818
## [9] 202.8212
```

#### 3.11

(a)

```
py1x <- function(x)
{
   val = dnorm(x,1,0.5)/(dnorm(x,1,0.5)+dnorm(x,0,1))
   return(val)
}

x <- seq(-2,5,by = 0.01)
plot(x,py1x(x),type = "l", ylab = "Probability")</pre>
```



(b)

```
x1 = sqrt(2/3*(2/3+log(2)))+4/3
x2 = -sqrt(2/3*(2/3+log(2)))+4/3

(EER <- 1/2*(pnorm(x1, 0, 1)-pnorm(x2, 0, 1))+
    1/2*(pnorm(x1, 1, 0.5,lower.tail = 0)+pnorm(x2, 1, 0.5)))</pre>
```

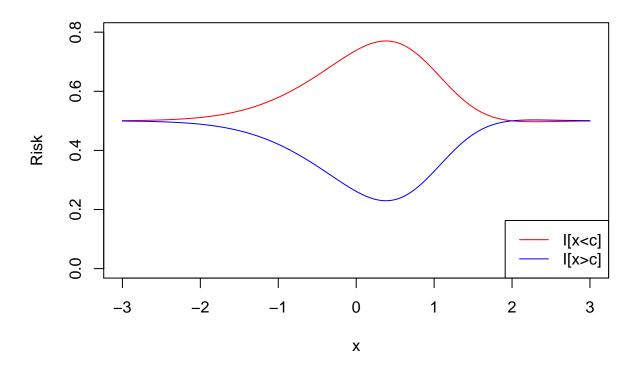
## [1] 0.2266941

(c)

```
linear1 <- function(c){
   value = (pnorm(c, 0, 1) + pnorm(c, 1, 0.5, lower.tail = 0))/2
   return(value)
}

linear2 <- function(c){
   value = (pnorm(c, 1, 0.5) + pnorm(c, 0, 1, lower.tail = 0))/2
   return(value)
}

x = seq(from = -3, to = 3, by = 0.02)
plot(x, linear1(x), ylim = c(0,0.8), type = "l", col = "red", ylab = "Risk")
legend("bottomright", c("I[x<c]", "I[x>c]"), col = c("red", "blue"), lty = c(1,1))
lines(x, linear2(x), col = "blue")
```



By minimizing the risk.

```
op <- optim(0.1, linear2, method = "Brent", lower = 0, upper = 1)
op$par</pre>
```

```
## [1] 0.381208
```

So  $g^* = I[x > 0.381]$ 

The error rate for  $g^*$  is:

```
op$value
## [1] 0.2297298
So the modeling penalty is:
op$value - EER
## [1] 0.003035748
(d)
M <- 10000
N <- 100
mini <- function(c)</pre>
 value = \min(sum(Y[X<c]==0)+sum(Y[X>c]==1), sum(Y[X<c]==1)+sum(Y[X>c]==0))
 return(value)
}
Error = 0
for(i in 1:M){
 Y <- sample(0:1, N, replace = TRUE)
 X \leftarrow rnorm(N, Y, 1-0.5*Y)
 c = optim(0, mini, method = "Brent", lower = -3, upper = 3)$par
  if(sum(Y[X<c]==0)+sum(Y[X>c]==1) <= sum(Y[X<c]==1)+sum(Y[X>c]==0)) \{ \\
   Error = Error + linear1(c)
 }
 else{
   Error = Error + linear2(c)
}
Error = Error/M
When N = 100, the error rate is:
Error
## [1] 0.2411257
The "fitting penalty" is:
Error - op$value
## [1] 0.01139588
M < -10000
N <- 50
mini <- function(c)</pre>
 return(value)
```

```
Error = 0
for(i in 1:M){
    Y <- sample(0:1, N, replace = TRUE)
    X <- rnorm(N, Y, 1-0.5*Y)

    c = optim(0, mini, method = "Brent", lower = -3, upper = 3)$par

    if(sum(Y[X<c]==0)+sum(Y[X>c]==1) <= sum(Y[X<c]==1)+sum(Y[X>c]==0)){
        Error = Error + linear1(c)
    }
    else{
        Error = Error + linear2(c)
    }
}
Error = Error/M
```

The error rate and the "fitting penalty" are:

Error; Error-op\$value

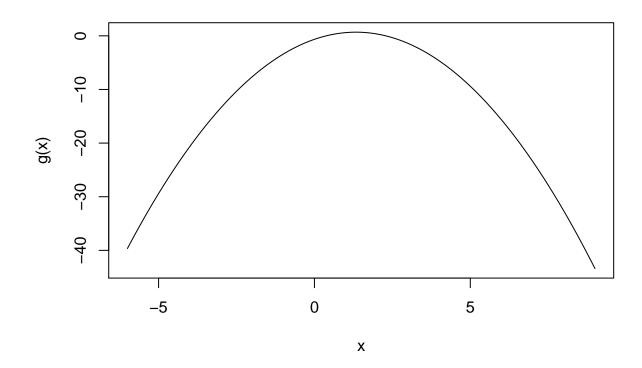
```
## [1] 0.2499291
## [1] 0.02019927
```

And, indeed, when we have a smaller sample size, the fitting penalty is greater.

#### 3.12

(a)

```
g <- function(x){
  value = 1/2*log(dnorm(x, 1, 0.5)/dnorm(x, 0, 1))
  return(value)
}
x <- seq(-6,9,by=0.02)
plot(x,g(x),ty = "l")</pre>
```



# (b)

```
data2 <- read_excel("~/Downloads/Problem3.4Data.xlsx")
xx <- data2$x; y <- data2$y

func2 <- function(para)
{
    beta0 = para[1]
    beta1 = para[2]
    beta2 = para[3]

    value = mean(exp(-y*(beta0+beta1*(xx-mean(xx))+beta2*(xx-mean(xx))^2)))
    return(value)
}

op1 <- optim(c(0,0,0),func2)
op1$par</pre>
```

**##** [1] 0.03694171 0.06596512 -0.03919403

(c)

```
func_2 <- function(x, op){</pre>
      value = oppar[1]+opppar[2]*(x-mean(xx))+opppar[3]*(x-mean(xx))^2
     return(value)
}
x < - seq(-6, 12, by=0.02)
plot(x, func_2(x, op1), ty = "l", ylab = "Value", lty = 3, col = 1, ylim = c(-7,1))
lines(x, g(x), lty = 1, col = 2)
data2 <- read_excel("~/Downloads/Problem3.4Data.xlsx")</pre>
xx <- data2$x; y <- data2$y
func3 <- function(para){</pre>
      beta0 = para[1]
      beta1 = para[2]
      beta2 = para[3]
     value = mean(exp(-y*(beta0+beta1*(xx-mean(xx)))+beta2*(xx-mean(xx))^2))) + lambda*beta2^2
     return(value)
}
lambda = 2
op5 <- optim(c(0,0,0), func3)
lines(x, func_2(x, op5), lty = 2, col = 3)
lambda = 1
op4 <- optim(c(0,0,0), func3)
lines(x, func_2(x, op4), lty = 4, col = 4)
lambda = 0.5
op3 <- optim(c(0,0,0), func3)
lines(x, func_2(x, op3), lty = 5, col = 5)
lambda = 0.25
op2 <- optim(c(0,0,0), func3)
lines(x, func_2(x, op2), lty = 6, col = 6)
legend("bottomright", lty = c(3,1,2,4,5,6), col = 1:6, legend = c("No Penalty", "Optimal", "Lambda = 2", legend("bottomright", lty = c(3,1,2,4,5,6), col = 1:6, legend("bottom
                                                                                                                                                   "Lambda = 1", "Lambda = 0.5", "Lambda = 0.25"))
```

