## STAT 37810 final project

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1

The algorithm starts from create an empty vector that stores  $\phi$ . For each draw, generates proposal phi by proposal function. Then, computes acceptance probability

$$r = \frac{p(\phi_{prop})/J_t(\phi_{prop}|\phi_{old})}{p(\phi_{old})/J_t(\phi_{old}|\phi_{phi})}$$

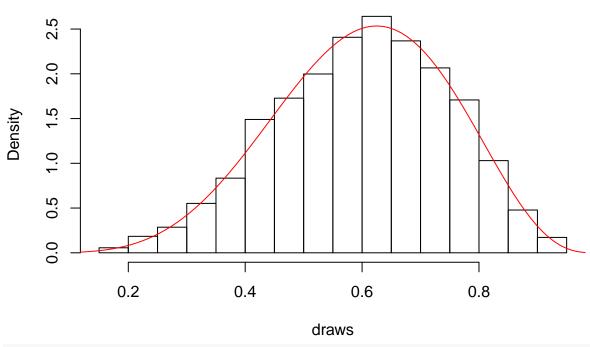
where p follows beta(6,4) and J follows proposal distribution. Then new phi is proposed phi if r is smaller than random uniform or else it is current phi.

#### 1.2

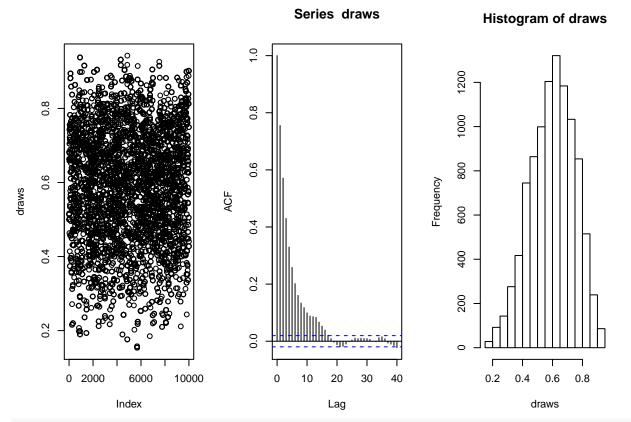
```
proposal <- function(prop,old,c){</pre>
  #This function computes probability from proposal functino
  dbeta(prop,c*old,c(1-old))
}
MHbeta <- function(nsample,initial,alpha,beta,c){</pre>
  # This function computes Metropolis-Hastings algorithm for beta
  phi=c()
  phi[1]=initial
  for (i in 1:nsample){
    current_phi=phi[i]
    proposal_phi=rbeta(1,c*current_phi,c(1-current_phi))
    A=(dbeta(proposal_phi,alpha,beta)/proposal(proposal_phi,current_phi,c))/(dbeta(current_phi,alpha,beta)
    if(runif(1)<A){</pre>
      phi[i+1]=proposal_phi
    }
    else{
      phi[i+1]=current_phi
```

```
}
  phi[2:(nsample+1)]
}
draws=MHbeta(10000,0.5,6,4,1)
hist(draws,freq =F)
xx = seq(0,1,length=100)
lines(xx,dbeta(xx,6,4),col="red")
```

# **Histogram of draws**



```
par(mfrow=c(1,3)) #1 row, 3 columns
plot(draws); acf(draws); hist(draws) #plot commands
```



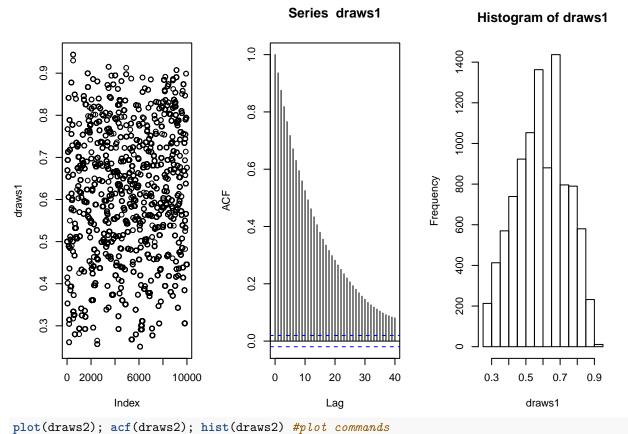
```
ks.test(draws, "pbeta", 6, 4)
```

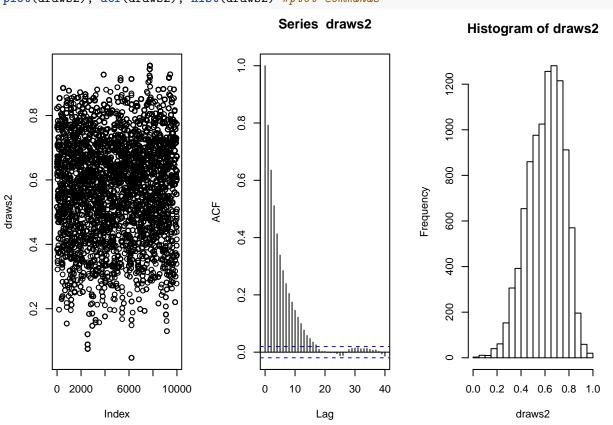
```
## Warning in ks.test(draws, "pbeta", 6, 4): ties should not be present for the
## Kolmogorov-Smirnov test
##
## One-sample Kolmogorov-Smirnov test
##
## data: draws
## D = 0.014715, p-value = 0.02632
## alternative hypothesis: two-sided
```

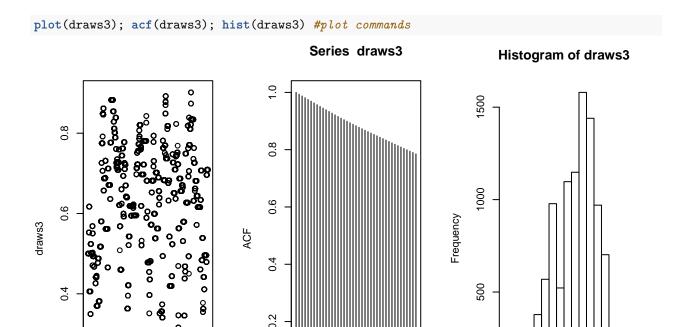
Based on the histogram, MH algorithms has simulated Beta(6,4) well. Based on Kolmogorov-Smirnov statistics, it is Beta(6,4).

#### 1.3

```
par(mfrow=c(1,3)) #1 row, 3 columns
draws1=MHbeta(10000,0.5,6,4,0.1) #c=0.1
draws2=MHbeta(10000,0.5,6,4,2.5) #c=2.5
draws3=MHbeta(10000,0.5,6,4,10) #c=10
plot(draws1); acf(draws1); hist(draws1) #plot commands
```







Based on the auto correlation plot, c=10 has highest auto correlation, c=0.1 comes next and c=2.5 has the least. This is because large or samll c will lead to slow change in phi. Therefore, it takes longer time to reach stationary distribution and thus requires higher burn-in rate.

20

Lag

30

40

0.2

0.4

0.6

draws3

0.8

0

10

0

10000

8

6000

Index

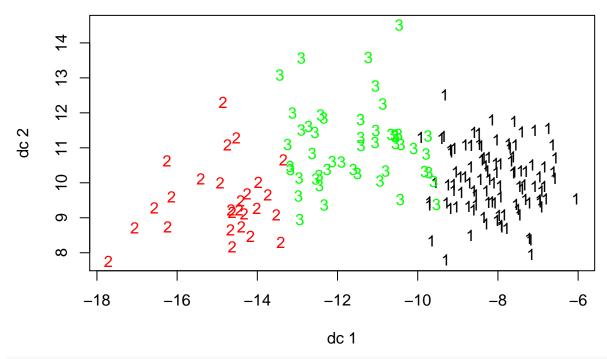
0 2000

#### 1.4

```
#not pre-allocate
MHbeta <- function(nsample,initial,alpha,beta,c){</pre>
  # This function computes Metropolis-Hastings algorithm for beta
  phi=c()
  phi[1]=initial
  for (i in 1:nsample){
    current_phi=phi[i]
    proposal_phi=rbeta(1,c*current_phi,c(1-current_phi))
    A=(dbeta(proposal_phi,alpha,beta)/proposal(proposal_phi,current_phi,c))/(dbeta(current_phi,alpha,beta)
    if(runif(1)<A){</pre>
      phi[i+1]=proposal_phi
    else{
      phi[i+1]=current_phi
  }
  phi[2:(nsample+1)]
t0=Sys.time()
draws=MHbeta(10000,0.5,6,4,1)
```

```
t1=Sys.time()
t1-t0 #not pre allocate
## Time difference of 0.1917138 secs
#pre-allocate
MHbeta <- function(nsample,initial,alpha,beta,c){</pre>
  # This function computes Metropolis-Hastings algorithm for beta
  phi=rep(0,1+nsample)
  phi[1]=initial
  for (i in 1:nsample){
    current_phi=phi[i]
    proposal_phi=rbeta(1,c*current_phi,c(1-current_phi))
    A=(dbeta(proposal_phi,alpha,beta)/proposal(proposal_phi,current_phi,c))/(dbeta(current_phi,alpha,beta)
    if(runif(1)<A){</pre>
      phi[i+1]=proposal_phi
    else{
      phi[i+1]=current_phi
 phi[2:(nsample+1)]
t0=Sys.time()
draws=MHbeta(10000,0.5,6,4,1)
t1=Sys.time()
t1-t0 #not pre allocate
## Time difference of 0.1443989 secs
Pre-allocation is faster
data(wine, package="rattle")
df=wine[,-1]
k=3
kmean <- function(df,k,tol=1){</pre>
  #This function computes k mean of dataframe df
  len=ncol(df) #number of features in data frame
  mat=matrix(0,nrow = len,ncol=k) #create empty matrix to storage k means
  for (i in 1: len){
    #computes initial value of means
    mat[i,]=runif(3,min(df[,i]),max(df[,i]))
  computelabel <- function(df,mat){</pre>
    #This functions label data point by computing and choosing minimum euclidean
    \#distance between data points and k means
    df nl <- df[,1:len] # remove label from df
    computeenclideandistance <- function(df,mat,k){</pre>
      #This sub function computes enclidean distance between mean and data points
      dist(rbind(df, t(mat)[k,]))
```

```
meandf <- apply(df_nl,1,computeenclideandistance,mat=mat,k=1) %>% as.data.frame
    for (i in 2:k){
      meandf <- meandf %>% cbind(apply(df_nl,1,computeenclideandistance,mat=mat,k=i))
    }
    label <- meandf%>%apply(1,which.min) # find minimum distance
    dfwithlabel <- cbind(df_nl,label)</pre>
    dfwithlabel
  computekmean <- function(df,mat){</pre>
    \#This\ functions\ updates\ mean\ for\ each\ iteration\ in\ k\ mean
    for (i in 1:k){
      newmean <- df%>%as.data.frame()%>%filter(label==i)%>%apply(2,mean)
      mat[,i] <- newmean[1:len]</pre>
    }
    \mathtt{mat}
  }
  while(TRUE){ #iteration
    df <- computelabel(df,mat) #compute label</pre>
    updated_mat=computekmean(df,mat) #update mean
    if (norm(updated_mat-mat)<tol){ #if smaller than tolerance</pre>
      return(as.data.frame(df))
    }
    mat=updated_mat
  }
}
test \leftarrow kmean(df,3,1) # run kmean with k=3, tol=1
#test
library('fpc')
plotcluster(df, test$label)
```



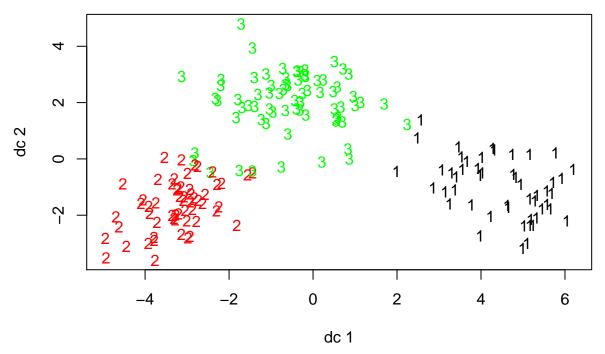
#df=data%>%mutate\_all(scale)

The data is well seperated.

### report

The algorithm starts with initialize matrix for recording k means. There are two sub functions in the algorithms: computelabel and computekmean. computelabel returns label based on the minimum euclidean distance between data points and k means. computekmean returns new mean matrix based on the current label. For each iteration, these two subfunctions are executed. New label and mean matrix is updated. The algorithm stops running when the norm of new mean matrix and previous mean matrix is below tolerance.

```
data.train <- scale(wine[-1])
test2 <- kmean(data.train,3,1)
plotcluster(data.train, test2$label)</pre>
```

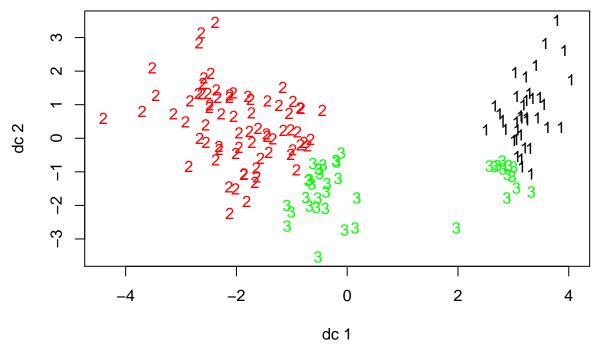


function scale normalize the data frame by first subtract the mean and then divided it by its standard deviation.

The

The results are different. This is because for unnormalized version, k mean weights more for features with larger value, which would lead to higher euclidean distance.

```
data.train <- scale(iris[,-5])
test2 <- kmean(data.train,3,1)
plotcluster(data.train, test2$label)</pre>
```



The

visualization looks well in both unnormalize and normalize version.