Lecture 17 Rscript, Hamiltonian Monte Carlo

Example dataset

Consider a logistic regression,

$$y_i \sim \operatorname{Bin}(n_i, p_i)$$

 $\log(p_i/(1-p_i)) = \mathbf{x}_i'\boldsymbol{\beta}.$

Clinical researchers are trying to determine the optimal dose level of a new medication for influenza. A group of influenza patients were recruited, and administered various doses of the experimental medication. One week after administration of medication, the researcher examined the patients to see if symptoms had improved. You were asked to investigate if there is a linear relationship between the log-odds of symptom improvement and medication dose.

Assume a flat prior for β so that your results are more comparable to a standard logistic GLM.

	Dosage: x						
y_i	0	1	2	3	4	5	6
No improvement: 0	8	5	6	6	3	5	1
Improvement: 1	2	5	4	4	7	5	9

Methods to consider

In this document, we will compare the performance of the Metropolis alogrithm to the Hamiltonian Monte Carlo algorithm when fitting a logistic regression.

Function one: Metropolis algorithm for fitting logistic regression

```
#Part one: function for performing Metropolis sampling for logistic regression normal random walk.
#Inputs:
#y: vector of responses
#n: vector (or scalar) of trial sizes.
#X: predictor matrix including intercept.
#c: rescaling for variance-covariance matrix, scalar J(theta*|theta(t-1)) = N(theta(t-1), c^2*Sigma)
\#Sigma is variance covariance matrix for parameters in J()
#iter: number of iterations
#burnin: number of initial iterations to throw out.
Metropolis.fn<-function(y,n,X,c,Sigma,iter,burnin){</pre>
p < -dim(X)[2]
                #number of parameters
library(mvtnorm)
theta0<-rnorm(p) #initial values.
theta.sim<-matrix(0,iter,p+1) #matrix to store iterations
theta.sim[1,1:p]<-theta0
for(i in 1:(iter-1)){
theta.cand <-rmvnorm(1,mean=theta.sim[i,1:p],sigma=c^2*Sigma) #draw candidate (jointly)
```

```
theta.cand <-as.numeric(theta.cand)</pre>
xbc
           <-X%*%theta.cand
           <-(1+exp(-xbc))^{(-1)}
                                   #Calculating probability of success for candidates.
p.c
хb
           <-X%*%theta.sim[i,1:p]
           <-(1+exp(-xb))^{-}(-1)
                                     #Calculating probability of success for theta(t-1).
p.b
#difference of log joint distributions.
r<-sum(dbinom(y,size=n,prob=p.c,log=TRUE)-dbinom(y,size=n,prob=p.b,log=TRUE))
#Draw an indicator whether to accept/reject candidate
ind \leftarrow rbinom(1, 1, exp(min(c(r, 0))))
theta.sim[i+1,1:p]<- ind*theta.cand + (1-ind)*theta.sim[i,1:p]</pre>
theta.sim[i+1,p+1]<- ind #counter for acceptances.
}
#Removing the iterations in burnin phase
results<-theta.sim[-c(1:burnin),]
names(results)<-c('beta0','beta1','accept') #column names</pre>
return(results)
}
```

Function two: Hamiltonian Monte Carlo algorithm for fitting logistic regression

```
#Part one: function for performing Hamiltonian Monte Carlo for logistic regression.
#Inputs:
#y: vector of responses
#n: vector (or scalar) of trial sizes.
#X: predictor matrix including intercept.
#L: number of leapfrog steps.
#M is variance covariance matrix for normal prior of momentum variable \phi. Ideally diagonal.
#iter: number of iterations
#burnin: number of initial iterations to throw out.
HMC.fn<-function(y,n,X,L,M,iter,burnin){</pre>
p \leftarrow dim(X)[2]
                #number of parameters
library(mvtnorm)
theta0<-rnorm(p) #initial values of beta
theta.sim<-matrix(0,iter,p+1) #matrix to store iterations plus acceptance.
theta.sim[1,1:p]<-theta0
                              #initial values in matrix.
epsilon<-1/L
                            \#epsilon \ assuming \ epsilon*L =1.
Minv <-solve(M)
for(i in 1:(iter-1)){
phi
          <-rmvnorm(1,mean=rep(0,p),sigma=M) #draw momentum variable.</pre>
phi
          <-as.numeric(phi)
phi0
          <-phi
                                             #saving starting phi for calculation of r.
          <-theta.sim[i,1:p]</pre>
                                              #current state of theta.
theta
            <-(1+exp(-X\%*\theta))^(-1) #calculate probabilities of success at current state.
gradtheta <- crossprod(X,y-n*p.b) #Gradient of posterior = joint distribution with respect to theta.
#leapfrog steps.
for(j in 1:L){
  phi <- phi + 0.5*epsilon*gradtheta #first half step for phi
  theta <- theta + epsilon*(Minv\%*\%phi) #full step for theta
```

```
<-(1+\exp(-X_*^*)^*(-1)  #calculate probabilities of success at candidate (sub) state.
gradtheta <- crossprod(X,y-n*p.c)</pre>
                                      #Gradient of posterior = joint distribution with respect to theta.
      <- phi + 0.5*epsilon*gradtheta #second half step for phi.</pre>
      <- as.numeric(phi)
phi
#difference of log joint distributions at final iteration of leap.frog vs current state.
r<-sum(dbinom(y,size=n,prob=p.c,log=TRUE))+dmvnorm(phi,mean=rep(0,p),sigma=M,log=TRUE)-sum(dbinom(y,si
#Draw an indicator whether to accept/reject candidate
ind \leftarrow rbinom(1,1,exp(min(c(r,0))))
theta.sim[i+1,1:p] \leftarrow ind*theta + (1-ind)*theta.<math>sim[i,1:p]
theta.sim[i+1,p+1] \leftarrow ind
#Removing the iterations in burnin phase
results<-theta.sim[-c(1:burnin),]
names(results)<-c('beta0','beta1','accept') #column names</pre>
return(results)
}
```

Formatting the data and running the code

```
x < -0:6
                        #doses
y < -c(2,5,4,4,5,7,9)
                        #responses
n<-10
                        #number in each group.
#formatting data into the correct format.
#Build predictor matrix.
pred.mat <-cbind(rep(1,length(y)),x) #First column intercept, second column dose levels
#Estimating a good choice of Sigma for the proposal distribution.
pest
        <-y/n
                                #Estimates of probability of success
logitest<-log(pest/(1-pest))</pre>
                                #Estimates of logits.
                                #fit lm with estimated logits as response.
modest <-lm(logitest~x)
#variance-covariance matrix extracted from lm object.
sigma <-vcov(modest)</pre>
#Running Metropolis algorithm
metro1<-Metropolis.fn(y=y,n=n,X=pred.mat,c=2.4/sqrt(2),Sigma=sigma,iter=10000,burnin=1000)
metro2<-Metropolis.fn(y=y,n=n,X=pred.mat,c=2.4/sqrt(2),Sigma=sigma,iter=10000,burnin=1000)
metro3<-Metropolis.fn(y=y,n=n,X=pred.mat,c=2.4/sqrt(2),Sigma=sigma,iter=10000,burnin=1000)
#Choosing candidate M
#Md <-1/diag(sigma)
M <- 1*solve(sigma)</pre>
HMC1<-HMC.fn(y=y,n=n,X=pred.mat,L=1,M=M,iter=10000,burnin=1000)
HMC2<-HMC.fn(y=y,n=n,X=pred.mat,L=1,M=M,iter=10000,burnin=1000)
HMC3<-HMC.fn(y=y,n=n,X=pred.mat,L=1,M=M,iter=10000,burnin=1000)
```

```
#Posterior means of beta0, beta1, Acceptance rate comparison
#Metropolis
metro.all<-rbind(metro1,metro2,metro3)</pre>
colMeans(metro.all)
## [1] -1.1935096 0.4206185 0.4169630
#Hamiltonian Monte Carlo
HMC.all <- rbind(HMC1,HMC2,HMC3)</pre>
colMeans(HMC.all)
## [1] -1.1936481 0.4210687 0.9263333
#Posterior standard deviations
apply(metro.all,2,FUN=sd)
## [1] 0.4751995 0.1365864 0.4930658
apply(HMC.all,2,FUN=sd)
## [1] 0.4904284 0.1412250 0.2612325
#95 % credible intervals
apply(metro.all,2,FUN=function(x) quantile(x,c(0.025,0.975)))
##
                         [,2] [,3]
               [,1]
## 2.5% -2.1558708 0.1656004
## 97.5% -0.2986174 0.6950493
apply(HMC.all,2,FUN=function(x) quantile(x,c(0.025,0.975)))
##
               [,1]
                        [,2] [,3]
## 2.5% -2.1970486 0.156878
## 97.5% -0.2642192 0.713073
#Convergence diagnostics.
library(coda)
#splitting metropolis chains for beta0, beta1 and converting into mcmc objects.
ml1<-as.mcmc.list(as.mcmc((metro1[1:4500,1:2])))
ml2<-as.mcmc.list(as.mcmc((metro2[1:4500,1:2])))
ml3<-as.mcmc.list(as.mcmc((metro3[1:4500,1:2])))
ml4<-as.mcmc.list(as.mcmc((metro1[4500+1:4500,1:2])))
ml5<-as.mcmc.list(as.mcmc((metro2[4500+1:4500,1:2])))
ml6<-as.mcmc.list(as.mcmc((metro3[4500+1:4500,1:2])))
m1<-c(m11,m12,m13,m14,m15,m16)
#Gelman-Rubin diagnostic.
gelman.diag(ml)[[1]]
       Point est. Upper C.I.
## [1,]
         1.000839
                    1.002069
## [2,]
         1.000871
                     1.002343
#effective sample size.
effectiveSize(ml)
##
       var1
                var2
## 3608.932 3438.832
```

```
#splitting Hamiltonian Monte Carlo chains for beta0, beta1 and converting into mcmc objects.
hl1<-as.mcmc.list(as.mcmc((HMC1[1:4500,1:2])))
hl2<-as.mcmc.list(as.mcmc((HMC2[1:4500,1:2])))
hl3<-as.mcmc.list(as.mcmc((HMC3[1:4500,1:2])))
hl4<-as.mcmc.list(as.mcmc((HMC1[4500+1:4500,1:2])))
h15<-as.mcmc.list(as.mcmc((HMC2[4500+1:4500,1:2])))
h16<-as.mcmc.list(as.mcmc((HMC3[4500+1:4500,1:2])))
hl<-c(hl1,hl2,hl3,hl4,hl5,hl6)
#Gelman-Rubin diagnostic.
gelman.diag(hl)[[1]]
        Point est. Upper C.I.
## [1,]
          1.001738
                    1.004692
## [2,]
          1.000690
                     1.002006
#effective sample size.
effectiveSize(hl)
##
       var1
                var2
## 5792.335 5174.447
#Checking convergence of combined Metropolis, Hamiltonian MC.
#Gelman-Rubin diagnostic.
gelman.diag(c(ml,hl))[[1]]
##
        Point est. Upper C.I.
## [1,]
          1.001041
                     1.002191
## [2,]
          1.000805
                     1.001694
#plotting HMC vs Metropolis iterations.
ylim1<-apply(rbind(metro.all,HMC.all),2,min)</pre>
ylim2<-apply(rbind(metro.all,HMC.all),2,max)</pre>
par(mfrow=c(2,2),mar=c(9.2,4.1,4.1,2.5))
for(i in 1:2){
plot(metro1[,i],type='l',ylim=c(ylim1[i],ylim2[i]),main='Metropolis output',xlab='Iteration',ylab=bquot
lines(metro2[,i],col=2)
lines(metro3[,i],col=3)
plot(HMC1[,i],type='l',ylim=c(ylim1[i],ylim2[i]),main='Hamiltonian MC output',xlab='Iteration',ylab=bqu
lines(HMC2[,i],col=2)
lines(HMC3[,i],col=3)
}
```







