

# Lab 7 solutions MAST90125: Bayesian Statistical Learning

Thursday 12 September 2019

## Writing Gibbs samplers in linear models.

In this weeks lab, we will discuss how to write code for Gibbs sampling of linear models.

## Instructions for assignment

Download USJudgeRatings.csv from LMS. Comment the codes below that purports to perform Gibbs sampling for a variety of linear models. See if you can determine what the code is doing. You may find referring back to lectures 12 and 13 useful. Run the code to see if it compares to the results obtained from using `stan` to fit the same models last week (Lab 6 version 2).

## Examples of Gibbs samplers

- Linear regression (flat prior for  $\beta$ ,  $p(\tau) \propto \tau^{-1}$ , where  $\tau = (\sigma^2)^{-1}$ ).

```
#This is a Gibbs sampler where beta is updated as a block. The arguments are
#X: matrix of predictors dimension n times p. Includes the intercept.
#y: response vector, length p.
#tau0: initial value for the residual precision.
#iter: number of iterations
#burnin: number of initial iterations to remove
Gibbs.lm1<-function(X,y,tau0,iter,burnin){
  p <- dim(X)[2] #number of predictors
  XTX <- crossprod(X)
  XTXinv <-solve(XTX)
  XTY <- crossprod(X,y)
  betahat<-solve(XTX,XTY) #betahat = (t(X)%*%X)^{-1}t(X)%*%y = mean of conditional posterior for beta
  tau <-tau0
  library(mvtnorm)

  par<-matrix(0,iter,p+1) #storing iterations, beta (length p) + tau (length 1)
  for( i in 1:iter){
    beta <- rmvnorm(1,mean=betahat,sigma=XTXinv/tau) #sample beta
    beta <- as.numeric(beta)
    err <- y-X%*%beta
    tau <- rgamma(1,0.5*n,0.5*sum(err^2)) #sample tau.
    par[i,] <-c(beta,tau) #store current round of beta, tau in par.
  }

  par <-par[-c(1:burnin),] #removing the first iterations
  return(par)
}
```

```

#Example of unblocked Gibbs sampling. We will update beta element by element. The arguments are

#X: matrix of predictors dimension n times p. Includes the intercept.
#y: response vector, length p.
#tau0: initial value for the residual precision.
#iter: number of iterations
#burnin: number of initial iterations to remove
Gibbs.lm2<-function(X,y,tau0,iter,burnin){
p <- dim(X)[2]
diagXTX <-colSums(X^2)      #calculates  $t(X)\%*\%X_{ii}$  for all i.
XTY <- crossprod(X,y)
betahat <- XTY/diagXTX      #component of conditional posteriors of beta_i's that is function of y.
tau    <-tau0

beta<-rnorm(p)
par<-matrix(0,iter,p+1)
for( i in 1:iter){
  for(j in 1:p){ #This samples beta element by element
    beta[j]<-0    #If we zero beta_j, then  $X_j\text{beta}_j = X\text{beta}$ .
    Xb    <-X%\%beta
    diff  <-t(X[,j])%\%Xb/diagXTX[j]
    beta[j] <- rnorm(1,mean=betahat[j]-diff,sd=1/sqrt(tau*diagXTX[j])) )
  }
  err  <- y-X%\%beta
  tau  <- rgamma(1,0.5*n,0.5*sum(err^2)) #samples tau from conditional posterior.
  par[i,] <-c(beta,tau)
}

par <-par[-c(1:burnin),]
return(par)
}

```

*#Example of Gibbs sampling where matrix decompositions are used to diagonalise conditional #posterior variances.This means beta is still updated as a block, but in a more efficient #way than in Gibbs.lm1. The arguments are*

```

#X: matrix of predictors dimension n times p. Includes the intercept.
#y: response vector, length p.
#tau0: initial value for the residual precision.
#iter: number of iterations
#burnin: number of initial iterations to remove
Gibbs.lm3<-function(X,y,tau0,iter,burnin){
p <- dim(X)[2] #dimension of p
svdX <-svd(X)  #matrix decomposition to speed up computation.
U    <-svdX$u   #extracting components of decompositions.
Lambda<-svdX$d
V    <-svdX$v
Vbhat <- crossprod(U,y)/Lambda #mean of conditional posterior for transformed parameters.
tau <-tau0

par<-matrix(0,iter,p+1)
for( i in 1:iter){
  sqrttau<-sqrt(tau)
  #posterior variances are diagonal for transformed parameters, so sequence of univariate normal draws

```

```

vbeta <- rnorm(p,mean=Vbhat,sd=1/(sqrttau*Lambda) )
beta <- V%*%vbeta #back transform to original parameter.
err <- y-X%*%beta
tau <- rgamma(1,0.5*n,0.5*sum(err^2)) #sample tau
par[i,] <-c(beta,tau)
}

par <-par[-c(1:burnin),] #remove initial iterations
return(par)
}

```

### Solution

```

#Formatting data, and running chains.
data<-read.csv('USJudgeRatings.csv')
response<-data$RTEN #response variable
n<-dim(data)[1]
intercept <-matrix(1,n,1) #Intercept (to be estimated without penalty)
Pred<-data[,2:12] #Predictor variables.
Pred<-as.matrix(scale(Pred))
X <-cbind(intercept,Pred)

system.time(chain1<-Gibbs.lm1(X=X,y=response,tau0=1,iter=10000,burnin=2000))

## user system elapsed
## 2.79 0.02 2.83

system.time(chain2<-Gibbs.lm1(X=X,y=response,tau0=5,iter=10000,burnin=2000))

## user system elapsed
## 2.77 0.00 2.77

system.time(chain3<-Gibbs.lm1(X=X,y=response,tau0=0.2,iter=10000,burnin=2000))

## user system elapsed
## 2.78 0.00 2.78

system.time(chain4<-Gibbs.lm2(X=X,y=response,tau0=1,iter=10000,burnin=2000))

## user system elapsed
## 1.17 0.01 1.20

system.time(chain5<-Gibbs.lm2(X=X,y=response,tau0=5,iter=10000,burnin=2000))

## user system elapsed
## 1.20 0.00 1.21

system.time(chain6<-Gibbs.lm2(X=X,y=response,tau0=0.2,iter=10000,burnin=2000))

## user system elapsed
## 1.2 0.0 1.2

system.time(chain7<-Gibbs.lm3(X=X,y=response,tau0=1,iter=10000,burnin=2000))

## user system elapsed
## 0.11 0.00 0.11

system.time(chain8<-Gibbs.lm3(X=X,y=response,tau0=5,iter=10000,burnin=2000))

## user system elapsed

```

```
##      0.07      0.00      0.07
system.time(chain9<-Gibbs.lm3(X=X,y=response,tau0=0.2,iter=10000,burnin=2000))

##      user      system elapsed
##      0.08      0.00      0.08

library(coda)
#Estimating Gelman -Rubin diagnostics.
#Note 8000 iterations were retained, so 50:50 split is iteration 1:4000 and iteration 4001:8000

#However first we must convert the output into mcmc lists for coda to interpret.
ml1<-as.mcmc.list(as.mcmc((chain1[1:4000,])))
ml2<-as.mcmc.list(as.mcmc((chain2[1:4000,])))
ml3<-as.mcmc.list(as.mcmc((chain3[1:4000,])))
ml4<-as.mcmc.list(as.mcmc((chain1[4000+1:4000,])))
ml5<-as.mcmc.list(as.mcmc((chain2[4000+1:4000,])))
ml6<-as.mcmc.list(as.mcmc((chain3[4000+1:4000,])))
estml<-c(ml1,ml2,ml3,ml4,ml5,ml6)

#Gelman-Rubin diagnostic.
gelman.diag(estml)[[1]]

##      Point est. Upper C.I.
## [1,] 1.0003413 1.0011865
## [2,] 1.0010677 1.0029840
## [3,] 1.0001089 1.0005646
## [4,] 1.0000408 1.0003908
## [5,] 1.0001809 1.0007105
## [6,] 1.0007508 1.0015460
## [7,] 1.0003355 1.0011845
## [8,] 0.9999888 1.0002935
## [9,] 1.0000326 1.0002266
## [10,] 0.9998869 1.0000718
## [11,] 0.9998638 0.9999721
## [12,] 0.9999424 1.0002215
## [13,] 1.0007316 1.0021488

#effective sample size.
effectiveSize(estml)

##      var1      var2      var3      var4      var5      var6      var7      var8
## 24000.00 24000.00 24000.00 24000.00 24000.00 24000.00 24000.00 24000.00
##      var9      var10      var11      var12      var13
## 24000.00 24000.00 24000.00 24000.00 14297.98

#However first we must convert the output into mcmc lists for coda to interpret.
ml1<-as.mcmc.list(as.mcmc((chain4[1:4000,])))
ml2<-as.mcmc.list(as.mcmc((chain5[1:4000,])))
ml3<-as.mcmc.list(as.mcmc((chain6[1:4000,])))
ml4<-as.mcmc.list(as.mcmc((chain4[4000+1:4000,])))
ml5<-as.mcmc.list(as.mcmc((chain5[4000+1:4000,])))
ml6<-as.mcmc.list(as.mcmc((chain6[4000+1:4000,])))
estml<-c(ml1,ml2,ml3,ml4,ml5,ml6)

#Gelman-Rubin diagnostic.
gelman.diag(estml)[[1]]
```

```
##          Point est. Upper C.I.
## [1,]    1.000128    1.000475
## [2,]    1.007395    1.017415
## [3,]    1.023773    1.055979
## [4,]    1.039445    1.092409
## [5,]    1.033504    1.076018
## [6,]    1.018484    1.038918
## [7,]    1.019510    1.048409
## [8,]    1.064335    1.146444
## [9,]    1.146979    1.340675
## [10,]   1.043523    1.090600
## [11,]   1.120122    1.277644
## [12,]   1.011997    1.027093
## [13,]   1.011637    1.029686
```

```
#effective sample size.
effectiveSize(estml)
```

```
##          var1          var2          var3          var4          var5          var6
## 23485.67566  610.57400  368.25729  325.25018  218.61119  300.50895
##          var7          var8          var9          var10          var11          var12
##  365.78797   77.65203   78.53642   51.70065   54.56167  253.53053
##          var13
## 1894.86162
```

```
#However first we must convert the output into mcmc lists for coda to interpret.
```

```
ml1<-as.mcmc.list(as.mcmc((chain7[1:4000,])))
ml2<-as.mcmc.list(as.mcmc((chain8[1:4000,])))
ml3<-as.mcmc.list(as.mcmc((chain9[1:4000,])))
ml4<-as.mcmc.list(as.mcmc((chain7[4000+1:4000,])))
ml5<-as.mcmc.list(as.mcmc((chain8[4000+1:4000,])))
ml6<-as.mcmc.list(as.mcmc((chain9[4000+1:4000,])))
estml<-c(ml1,ml2,ml3,ml4,ml5,ml6)
```

```
#Gelman-Rubin diagnostic.
```

```
gelman.diag(estml)[[1]]
```

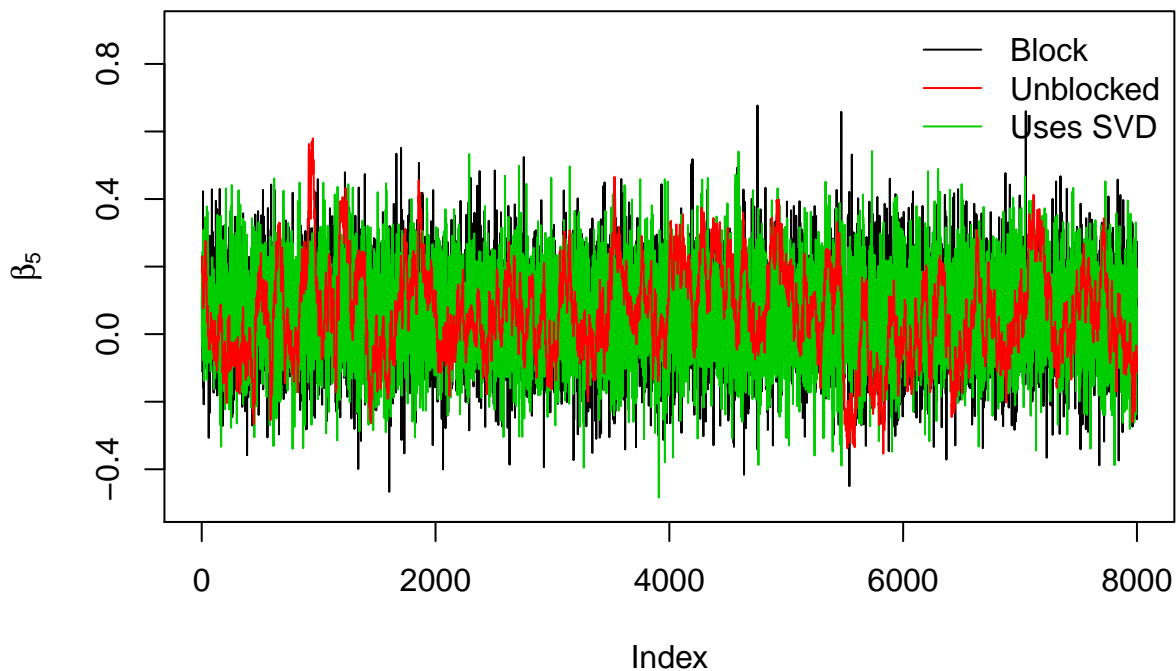
```
##          Point est. Upper C.I.
## [1,]   0.9998276   0.9998839
## [2,]   0.9999642   1.0001546
## [3,]   0.9999795   1.0001247
## [4,]   0.9999889   1.0003297
## [5,]   1.0000917   1.0002862
## [6,]   1.0001071   1.0005994
## [7,]   1.0000764   1.0003444
## [8,]   0.9998229   0.9998997
## [9,]   0.9999119   1.0000383
## [10,]  1.0000571   1.0003483
## [11,]  0.9999042   1.0001241
## [12,]  0.9998289   0.9999014
## [13,]  1.0002738   1.0009841
```

```
#effective sample size.
effectiveSize(estml)
```

```
##          var1          var2          var3          var4          var5          var6          var7          var8
## 24000.00 24724.44 24000.00 24000.00 24000.00 24000.00 25330.83 24000.00
```

```
##      var9      var10      var11      var12      var13
## 24000.00 24000.00 25121.25 24000.00 13821.75

#Comparing one co-efficient (the 5th)
plot(chain1[,5],type='l',ylim=c(-0.5,0.9),ylab=expression(beta[5]))
lines(chain7[,5],type='l',col=3,ylab=expression(beta[5]))
lines(chain4[,5],type='l',col=2,ylab=expression(beta[5]))
legend('topright',legend=c('Block','Unblocked','Uses SVD'),col=1:3,lty=1,bty='n')
```



```
#Reporting posterior means and credible intervals.
#Means
colMeans(rbind(chain1,chain2,chain3)) #Blocked

## [1] 7.602290608 0.011873407 0.280548379 0.143845680 0.059164994
## [6] -0.166363055 0.223724159 -0.002440861 -0.127163529 0.551011374
## [11] -0.064513015 0.252635156 72.335146917

colMeans(rbind(chain4,chain5,chain6)) #unblocked

## [1] 7.60239336 0.01379158 0.27763385 0.15338201 0.06821053
## [6] -0.16497641 0.22047548 -0.02897290 -0.09242135 0.53466006
## [11] -0.07101967 0.25478662 73.22236799

colMeans(rbind(chain7,chain8,chain9)) #Blocked but using svd to speed up computation.

## [1] 7.6022853183 0.0119740385 0.2809603957 0.1426368807 0.0595562056
## [6] -0.1681504228 0.2239776601 -0.0005679958 -0.1301031712 0.5567026741
## [11] -0.0669247771 0.2520019094 72.5556248923
```

```

#Credible interval
apply(rbind(chain1,chain2,chain3) ,2, FUN =function(x) quantile(x,c(0.025,0.975) )) #Blocked

##           [,1]      [,2]      [,3]      [,4]      [,5]      [,6]
## 2.5%  7.565643 -0.03743149 0.07402446 -0.06492396 -0.2027753 -0.42575267
## 97.5% 7.638450 0.06095730 0.48186976 0.35529014 0.3231502 0.09399502
##           [,7]      [,8]      [,9]      [,10]     [,11]     [,12]
## 2.5% -0.004423966 -0.4720929 -0.6451002 -0.02576225 -0.6923035 0.1329368
## 97.5% 0.451810389 0.4617845 0.3931162 1.12753074 0.5588945 0.3716479
##           [,13]
## 2.5%  41.36176
## 97.5% 111.81489

apply(rbind(chain4,chain5,chain6) ,2, FUN =function(x) quantile(x,c(0.025,0.975) )) #Unblocked

##           [,1]      [,2]      [,3]      [,4]      [,5]      [,6]
## 2.5%  7.566024 -0.03604019 0.06853033 -0.05799356 -0.1905866 -0.40733663
## 97.5% 7.638102 0.06287613 0.47676851 0.35472300 0.3433845 0.08847645
##           [,7]      [,8]      [,9]      [,10]     [,11]     [,12]
## 2.5%  0.006880288 -0.4894330 -0.5403710 -0.04866047 -0.6211989 0.1353311
## 97.5% 0.434404271 0.4290811 0.3460951 1.08043579 0.4642468 0.3740263
##           [,13]
## 2.5%  41.82221
## 97.5% 112.87112

apply(rbind(chain7,chain8,chain9) ,2, FUN =function(x) quantile(x,c(0.025,0.975) )) #Blocked with svd.

##           [,1]      [,2]      [,3]      [,4]      [,5]      [,6]
## 2.5%  7.566051 -0.03803262 0.0769504 -0.06711342 -0.2014690 -0.4260054
## 97.5% 7.638689 0.06198022 0.4846847 0.35117813 0.3229448 0.0883075
##           [,7]      [,8]      [,9]      [,10]     [,11]     [,12]
## 2.5% -0.003135034 -0.4685738 -0.6462462 -0.01599512 -0.6851662 0.1355050
## 97.5% 0.448216706 0.4671695 0.3816365 1.11751273 0.5523074 0.3693659
##           [,13]
## 2.5%  41.25019
## 97.5% 113.12326

```

- Linear mixed model/ ridge regression (flat prior for  $\beta_0$ ,  $p(\tau) = \text{Ga}(\alpha_e, \gamma_e)$ , where  $\tau = (\sigma^2)^{-1}$ ,  $\beta \sim \mathcal{N}(\mathbf{0}, \sigma_\beta^2 \mathbf{I})$ ,  $(\sigma_\beta^2)^{-1} = \tau_\beta \sim \text{Ga}(\alpha_\beta, \gamma_\beta)$ ).

```
#Arguments are
#X: matrix of predictors dimension n times p with flat prior. Includes the intercept.
#Z: matrix of predictors dimension n times q with normal prior for u.
#y: response vector, length p.
#taue_0, tauu_0: initial value for the residual and random effect precision.
#a.u, b.u. Hyper-parameter for gamma prior for tau_u
#a.e, b.e. Hyper-parameter for gamma prior for tau_e
#iter: number of iterations
#burnin: number of initial iterations to remove
normalmm.Gibbs<-function(iter,Z,X,y,burnin,taue_0,tauu_0,a.u,b.u,a.e,b.e){
  n <-length(y) #no. observations
  p <-dim(X)[2] #no of fixed effect predictors.
  q <-dim(Z)[2] #no of random effect levels
  tauu<-tauu_0
  taue<-taue_0
  #starting value for u.
  u0 <-rnorm(q,0,sd=1/sqrt(tauu))

  #Building combined predictor matrix.
  W<-cbind(X,Z) #for the joint conditional posterior for b,u
  WTW <-crossprod(W)
  library(mvtnorm)

  #storing results.
  par <-matrix(0,iter,p+q+2) #p beta coefficient, q u coefficients and 2 precision coefficient.

  #Create modified identity matrix for joint posterior.
  IO <-diag(p+q)
  diag(IO)[1:p]<-0

  for(i in 1:iter){
    #Conditional posteriors.
    tauu <-rgamma(1,a.u+0.5*q,b.u+0.5*sum(u0^2)) #sample tau_u
    #Updating component of normal posterior for beta,u
    Prec <-WTW + tauu*IO/taue
    P.mean <- solve(Prec)%*%crossprod(W,y)
    P.var <-solve(Prec)/taue
    betau <-rmvnorm(1,mean=P.mean,sigma=P.var) #sample beta, u
    betau <-as.numeric(betau)
    err <- y-W%*%betau
    taue <-rgamma(1,a.e+0.5*n,b.e+0.5*sum(err^2)) #sample tau_e
    #storing iterations for beta, u, and standard deviation of e, u.
    par[i,]<-c(betau,1/sqrt(tauu),1/sqrt(taue))
    u0 <-betau[p+1:q] #extracting u so we can update tau_u.
  }

  par <-par[-c(1:burnin),] #removing initial iterations
  colnames(par)<-c(paste('beta',1:p,sep=''),paste('u',1:q,sep=''),'sigma_b','sigma_e')
  return(par)
}
```

*Solution*



```

system.time(chain10<-normalmm.Gibbs(iter=10000,Z=Pred,X=intercept,y=response,burnin=2000,taue_0=1,tauu_0=1))

##      user  system elapsed
##      3.67    0.00    3.67

system.time(chain11<-normalmm.Gibbs(iter=10000,Z=Pred,X=intercept,y=response,burnin=2000,taue_0=0.2,tauu_0=0.2))

##      user  system elapsed
##      3.56    0.00    3.59

system.time(chain12<-normalmm.Gibbs(iter=10000,Z=Pred,X=intercept,y=response,burnin=2000,taue_0=5,tauu_0=5))

##      user  system elapsed
##      3.63    0.00    3.64

library(coda)
#Estimating Gelman -Rubin diagnostics.
#Note 8000 iterations were retained, so 50:50 split is iteration 1:4000 and iteration 4001:8000

#However first we must convert the output into mcmc lists for coda to interpret.
ml1<-as.mcmc.list(as.mcmc((chain10[1:4000,])))
ml2<-as.mcmc.list(as.mcmc((chain11[1:4000,])))
ml3<-as.mcmc.list(as.mcmc((chain12[1:4000,])))
ml4<-as.mcmc.list(as.mcmc((chain10[4000+1:4000,])))
ml5<-as.mcmc.list(as.mcmc((chain11[4000+1:4000,])))
ml6<-as.mcmc.list(as.mcmc((chain12[4000+1:4000,])))
estml<-c(ml1,ml2,ml3,ml4,ml5,ml6)

#Gelman-Rubin diagnostic.
gelman.diag(estml)[[1]]

##      Point est. Upper C.I.
## beta1      1.0005905  1.0017822
## u1         1.0002126  1.0008362
## u2         1.0002637  1.0009665
## u3         1.0003958  1.0012218
## u4         1.0000295  1.0001686
## u5         1.0001599  1.0005810
## u6         0.9998912  0.9999608
## u7         1.0000812  1.0003106
## u8         1.0001532  1.0006321
## u9         1.0004075  1.0010858
## u10        1.0000476  1.0004752
## u11        1.0000844  1.0005442
## sigma_b    1.0005441  1.0015565
## sigma_e    1.0000453  1.0004379

#effective sample size.
effectiveSize(estml)

##      beta1      u1      u2      u3      u4      u5      u6      u7
## 23250.48 24000.00 22996.89 22942.12 24000.00 19920.87 21743.77 24000.00
##      u8      u9      u10     u11  sigma_b  sigma_e
## 22309.72 19109.04 24000.00 24000.00 12994.46 15588.48

#Reporting posterior means and credible intervals.
#Means

```

```
colMeans(rbind(chain10,chain11,chain12))
```

```
##      beta1      u1      u2      u3      u4      u5
## 7.60242403 0.01390440 0.25834971 0.20378433 0.04400074 -0.06571223
##      u6      u7      u8      u9      u10      u11
## 0.15023617 0.01105592 -0.02307964 0.25012577 0.05703116 0.27183842
##      sigma_b      sigma_e
## 0.20527654 0.11821907
```

```
#95 % central Credible interval
```

```
apply(rbind(chain10,chain11,chain12) ,2, FUN =function(x) quantile(x,c(0.025,0.975) ))
```

```
##      beta1      u1      u2      u3      u4      u5
## 2.5% 7.566171 -0.02986222 0.09991092 0.0441440 -0.1425814 -0.2745050
## 97.5% 7.638656 0.05788413 0.41911658 0.3592429 0.2365529 0.1331029
##      u6      u7      u8      u9      u10      u11
## 2.5% -0.02977526 -0.2622409 -0.2999797 -0.04305112 -0.2493746 0.1784655
## 97.5% 0.33509468 0.2772295 0.2352198 0.58498483 0.3516782 0.3646120
##      sigma_b      sigma_e
## 2.5% 0.1230549 0.09333689
## 97.5% 0.3508011 0.15192766
```

- LASSO.

```

#Arguments
#X: matrix of predictors dimension n times p with flat prior. Includes the intercept.
#Z: matrix of predictors dimension n times q with normal prior for u.
#y: response vector, length p.
#taue_0: initial value for the residual precision.
#lambda: LASSO penalty
#a.e,b.e. Hyper-parameter for gamma prior for tau_e
#iter: number of iterations
#burnin: number of initial iterations to remove
normallasso.Gibbs<-function(iter,Z,X,y,burnin,taue_0,lambda,a.e,b.e){
  library(LaplacesDemon)
  n <-length(y) #no. observations
  p <-dim(X)[2] #no of fixed effect predictors.
  q <-dim(Z)[2] #no of random effect levels
  taue<-taue_0
  #Note Laplace distribution is compound of normal and exponential.
  #This results with us working with a vector tau_u = dimension of predictor with LASSO penalty.
  tauu <-rinvgaussian(q,lambda/abs(rnorm(q)),lambda^2) #Note LASSO can b

  #Building combined predictor matrix.
  W<-cbind(X,Z)
  WTW <-crossprod(W)
  library(mvtnorm)

  #storing results.
  par <-matrix(0,iter,p+q+1)

  for(i in 1:iter){
    #Conditional posteriors.

    #Updating component of normal posterior for beta,u
    Kinv <-diag(p+q)
    diag(Kinv)[1:p]<-0
    diag(Kinv)[p+1:q]<-tauu #Adding precision of predictors with LASSO penalty.

    Prec <-taue*WTW + Kinv
    P.var <-solve(Prec)
    P.mean <- taue*P.var%*%crossprod(W,y)
    betau <-rmvnorm(1,mean=P.mean,sigma=P.var) #sampling beta,u
    betau <-as.numeric(betau)
    err <- y-W%*%betau
    taue <-rgamma(1,a.e+0.5*n,b.e+0.5*sum(err^2)) #sampling tau_e, residual precision.
    #sampling tau_u, the augmented vector of precisions for predictor with LASSO penalty.
    tauu <-rinvgaussian(q,lambda/abs(betau[-c(1:p)]),lambda^2)

    #storing iterations.
    par[i,]<-c(betau,1/sqrt(taue))
  }

  par <-par[-c(1:burnin),] #removing early iterations
  colnames(par)<-c(paste('beta',1:p,sep=''),paste('u',1:q,sep=''),'sigma_e')
  return(par)
}

```

```
}
```

*Solution*

```
system.time(chain13<-normallasso.Gibbs(iter=10000,Z=Pred,X=intercept,y=response,burnin=2000,taue_0=1,la
```

```
## Warning: package 'LaplacesDemon' was built under R version 3.6.1
```

```
##
```

```
## Attaching package: 'LaplacesDemon'
```

```
## The following objects are masked from 'package:mvtnorm':
```

```
##
```

```
##      dmvt, rmvt
```

```
##      user  system elapsed
```

```
##      3.85    0.01    3.87
```

```
system.time(chain14<-normallasso.Gibbs(iter=10000,Z=Pred,X=intercept,y=response,burnin=2000,taue_0=0.2,
```

```
##      user  system elapsed
```

```
##      3.81    0.00    3.83
```

```
system.time(chain15<-normallasso.Gibbs(iter=10000,Z=Pred,X=intercept,y=response,burnin=2000,taue_0=5,la
```

```
##      user  system elapsed
```

```
##      3.83    0.00    3.85
```

```
library(coda)
```

```
#Estimating Gelman -Rubin diagnostics.
```

```
#Note 8000 iterations were retained, so 50:50 split is iteration 1:4000 and iteration 4001:8000
```

```
#However first we must convert the output into mcmc lists for coda to interpret.
```

```
ml1<-as.mcmc.list(as.mcmc((chain13[1:4000,])))
```

```
ml2<-as.mcmc.list(as.mcmc((chain14[1:4000,])))
```

```
ml3<-as.mcmc.list(as.mcmc((chain15[1:4000,])))
```

```
ml4<-as.mcmc.list(as.mcmc((chain13[4000+1:4000,])))
```

```
ml5<-as.mcmc.list(as.mcmc((chain14[4000+1:4000,])))
```

```
ml6<-as.mcmc.list(as.mcmc((chain15[4000+1:4000,])))
```

```
estml<-c(ml1,ml2,ml3,ml4,ml5,ml6)
```

```
#Gelman-Rubin diagnostic.
```

```
gelman.diag(estml)[[1]]
```

```
##      Point est. Upper C.I.
```

```
## beta1    1.0002566    1.000848
```

```
## u1       0.9999005    1.000086
```

```
## u2       0.9999033    1.000077
```

```
## u3       0.9999584    1.000242
```

```
## u4       0.9999808    1.000308
```

```
## u5       1.0000840    1.000529
```

```
## u6       0.9999940    1.000225
```

```
## u7       1.0001983    1.000805
```

```
## u8       1.0004376    1.001369
```

```
## u9       0.9999510    1.000200
```

```
## u10      1.0005808    1.001583
```

```
## u11      1.0000450    1.000421
```

```
## sigma_e  1.0001940    1.000551
```

```
#effective sample size.
```

```
effectiveSize(estml)
```

```
##      beta1      u1      u2      u3      u4      u5      u6      u7
## 24000.00 25209.60 24000.00 23195.45 24000.00 23025.30 22804.53 24000.00
##      u8      u9      u10      u11 sigma_e
## 23349.11 21831.36 22794.48 22557.86 13864.81
```

```
#Reporting posterior means and credible intervals.
```

```
#Means
```

```
colMeans(rbind(chain13,chain14,chain15))
```

```
##      beta1      u1      u2      u3      u4
## 7.602330480 0.011885803 0.280320153 0.147112797 0.057709753
##      u5      u6      u7      u8      u9
## -0.155574985 0.215273921 -0.003039792 -0.119749787 0.526449930
##      u10      u11      sigma_e
## -0.052009170 0.254331718 0.119817781
```

```
#Credible interval
```

```
apply(rbind(chain13,chain14,chain15) ,2, FUN =function(x) quantile(x,c(0.025,0.975) ))
```

```
##      beta1      u1      u2      u3      u4      u5
## 2.5% 7.566110 -0.03706777 0.07752921 -0.05359614 -0.1918535 -0.41229813
## 97.5% 7.638336 0.06088788 0.47873881 0.35473214 0.3093565 0.09508631
##      u6      u7      u8      u9      u10      u11
## 2.5% -0.0053994 -0.4412188 -0.6051434 -0.02632271 -0.6259150 0.1362780
## 97.5% 0.4359041 0.4359297 0.3602631 1.08511898 0.5237146 0.3721081
##      sigma_e
## 2.5% 0.09420474
## 97.5% 0.15491646
```