

A Mixture of Flexible Multivariate Distributions

In this document, I will compute the posterior distribution for β_j using the derivation in the document ‘next Steps’, where the prior on β_j is modeled as a mixture of multivariate normals.

```
tuto.dir=getwd()
##' @param b.gp.hat PxR matrix of standardized effect sizes across all `R` tissue types,
##' @param se.gp.hat RxR estimated covariance matrix of standardized standard errors, corresponding to
##' @param t.stat PxR matrix of t statistics for each gene-snp Pair across all R tissues
##' @param U.0kl RxR prior covariance matrix for posterior.covariance matrix K and weight l
##' @param pi LxK matrix of prior weights estimated from the EM algorithm which correspond to optimal w

b.gp.hat=na.omit(read.table("16008genesnppairs_43tissues_beta.hat.std.txt",header=F,skip=1)[-c(1,2)])
se.gp.hat=na.omit(read.table("16008genesnppairs_43tissues_beta.hat.std.txt",header=F,skip=1)[-c(1,2)])
t.stat=na.omit(read.table("16008genesnppairs_43tissues_t.stat.txt",header=F,skip=1)[-c(1,2)])
L =  #(number of grid weights to use)
R=ncol(b.gp.hat) #number of tissues
X.t=as.matrix(t.stat)
X.c=apply(X.t,2,function(x) x-mean(x)) ##Column centered matrix of t statistics
 #colMeans(X.c)
```

Now, we need to load in the prior matrices which we will try to find the optimal combination. First we load in the ‘K’ RxR prior covariance matrices specifying the relative importance of a particular tissue direction in each component.

```
##' @param R tissues from which to estimate covariance matrices
##' @param X.c a matrix of column center tstatistics
##' @param K number of PCs to keep in approximation
##' @return return K component list of covariance matrices (SFA or SVD approximations)
get.prior.covar.U.0=function(R,X.c,P,omega){

  U.0.=list()
  U.0.[[1]]=omega*diag(1,R) # the first covariance matrix will be the 'sopped up' identity
  U.0.[[2]]=omega*(t(X.c)%*%X.c)/(nrow(X.c))
  svd.X=svd(X.t) ##perform SVD on uncentered matrix
  v=svd.X$v;u=svd.X$u;d=svd.X$d

  cov.pc=1/P*v[,1:P]%*%diag(d[1:P])%*%t(v[,1:P]) ##Use the rank P summary representation

  U.0.[[3]]=omega*cov.pc
  return(U.0.)}

U.0=get.prior.covar.U.0(R,X.c,13,omega)
```

Now, for every prior covariance matrix U_k , we compute L ‘stretches’ specifying the width of this distribution.

```
##' @param function to return the K component list of prior covariance matrices
##' @param X.c a matrix of column center tstatistics
##' @param K number of PCs to keep in approximation
##' @return return L dimensional list of K dimensional lists where each L,K contains the Lth grid compon

get.prior.covar.Ukl=function(P){
```

```

    U.0kl=lapply(seq(1:L),function(l){ ##For each element of omega, computes the J covariance matrices
      omega=read.table("~/omega2.txt")[l,]
      get.prior.covar.U.0(R,X.c,P,omega)
    })
  return(U.0kl)}

```

Now that we have the prior covariance matrices, we can maximize the likelihood $L(\pi; \hat{\beta}_j, se_j)$ by computing $Pr(\hat{\beta}_j | Componentk, l)$ for each gene component at each gene SNP pair and then finding the optimal combination among all gene SNP pairs.

```

##' @return Compute a likelihood using the prior covariance matrix for each gene SNP pair
##' @param U.0kl K element list of L element list with prior covairnace matrix fo reach

U.0kl=get.prior.covar.Ukl(13)
#U.0kl[[L+1]]=diag(1,R) ##Add the identity matrix to the list
K=3
L=2
library("mvtnorm")
likelihood=function(b.gp.hat,se.gp.hat)
{
  lik.i=list()
  lapply(1:nrow(b.gp.hat),function(i){
    b.mle=b.gp.hat[i,]
    V.gp.hat=diag(se.gp.hat[i,])^2

    lik=matrix(NA,nrow=K,ncol=L)
    for(l in 1:L){
      for (k in 1:K){
        lik[k,l]=dmvnorm(x=b.mle, sigma=U.0kl[[l]][[k]] + V.gp.hat)
      }
      lik.i[[i]] = lik
    }
  })
}

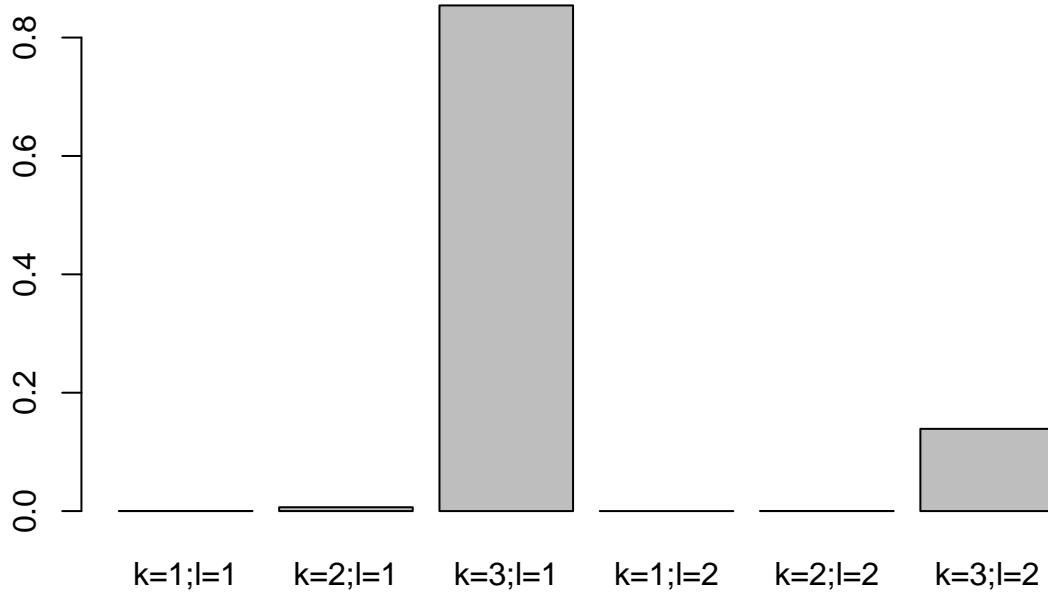
##` @return global.lik Computes the likelihood matrices for each componenet for each of j gene snp pairs
##' @return global.sum Sums the likelihood for each component across j gene snp pairs
##' @return global.norm Sums the likelihood for all components across all j pairs
updated.weights=function(b.gp.hat,se.gp.hat){
  global.lik=likelihood(b.gp.hat,se.gp.hat)
  global.sums=Reduce('+', likelihood(b.gp.hat,se.gp.hat)) ##This sums all matrices, so we get one KxL mat.
  global.norm=Reduce('+',lapply(global.lik,sum))
  return(updated.weights=global.sums/global.norm)}

names.vec=matrix(NA,nrow=K,ncol=L)
for(l in 1:L){
  for(k in 1:K){
    names.vec[k,l]=paste0("k=",k,";l=",l)}
}

x=updated.weights(b.gp.hat,se.gp.hat)
barplot(as.vector(x),names=as.vector(names.vec),main="Normalized Likelihood at Each Component")

```

Normalized Likelihood at Each Component



```
##Test with prior weights as 1/31
pi=rep(1/length(K*L),K*L)
```

Here, I've plotted the relative importance of each component after one iteration with a uniform prior weight on each π .

Recall:

$$\begin{aligned} L(\beta_j; k, l) &= Pr(\hat{b}_j | k, l) \\ &= Pr(\hat{b}_j; 0, \omega_l^2 U_k + \hat{V}) \end{aligned}$$

$$w_{jkl} = \frac{\pi_{k,l} L(\beta_j; k, l)}{\sum_{k,l} \pi_{k,l} L(\beta_j; k, l)}$$

We can then update our estimate of $\pi_{k,l}$ as

$$\pi_{kl}^{i+1} = \frac{\sum_j w_{jkl}}{\sum_{j,k,l} w_{jkl}}$$