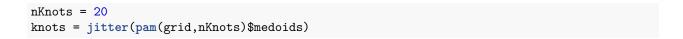
Recovering_rf_mixture

The objective of these simulations are to examine whether mixture parameters are recoverable with spatial data. We'll start by simulating some data on a grid from two gaussian random fields. We will initially assume that the fields have the same shape/scale parameters of the gaussian covariance function, but the variances are different – this allows one field to be normal, and the other fat-tailed.

```
library(cluster)
library(mvtnorm)
library(R2jags)
## Loading required package: rjags
## Loading required package: coda
## Linked to JAGS 3.4.0
## Loaded modules: basemod, bugs
##
## Attaching package: 'R2jags'
## The following object is masked from 'package:coda':
##
##
       traceplot
library(fields)
## Loading required package: spam
## Loading required package: grid
## Spam version 1.0-1 (2014-09-09) is loaded.
## Type 'help( Spam)' or 'demo( spam)' for a short introduction
## and overview of this package.
## Help for individual functions is also obtained by adding the
## suffix '.spam' to the function name, e.g. 'help( chol.spam)'.
##
## Attaching package: 'spam'
##
## The following objects are masked from 'package:base':
##
       backsolve, forwardsolve
##
##
## Loading required package: maps
grid = as.matrix(expand.grid("lon" = seq(5,15,1), "lat" = seq(5,15,1)))
nLocs = dim(grid)[1]
```

Approximating data with spatial random field

We could use RandomField, or other packages to do this. Initially we'll just simulate data using the estimation model. Use pam() to choose a number of knots on the grid. We'll specify 20 initially. We'll also jitter them slightly so knot locations don't fall exactly on stations.



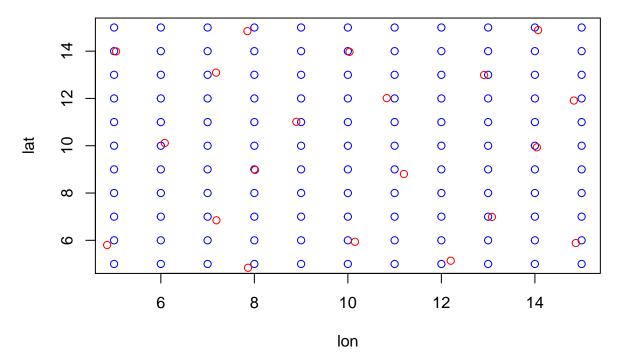


Figure 1: Simulated grid (blue) and knots for random effects (red)

Approximating data with spatial random field

Initially, we'll assume all data is observed in the same year, so there'll just be a single random effect field for normal years, and a single effect for catastrophic / abnormal years.

```
# distance matrix of knots
distKnots = as.matrix(dist(knots))
distKnotsSq = distKnots^2 # squared distances
# note: shape parameter scaled to distance matrix
gp_scale = 0.001
sigma.norm = 0.001
sigma.fat = 0.002
corKnots = exp(-gp_scale*distKnotsSq)
Sigma.normal = corKnots * sigma.norm * sigma.norm
Sigma.fat = corKnots * sigma.fat * sigma.fat
```

Calculate distance matrix from stations to knots, and covariance matrices,

```
# Calculate distance from knots to grid
distAll = as.matrix(dist(rbind(grid, knots)))^2
distKnots21Sq = t(distAll[1:nKnots, -c(1:nKnots)])

Sigma21.normal = exp(-gp_scale*distKnots21Sq) * sigma.norm * sigma.norm
Sigma21.fat = exp(-gp_scale*distKnots21Sq) * sigma.fat * sigma.fat
```

Generate single MVN field for each of the normal and 'fat-tailed' distributions.

```
re.norm = rmvnorm(1, mean = rep(0, nKnots), Sigma.normal)
re.fat = rmvnorm(1, mean = rep(0, nKnots), Sigma.fat)
```

Project to the station / grid locations, and combine the random fields using some proportion of 'fat-tailed' events. Initially, we'll use 0.05.

```
# take inverse
invSigmaKnots.norm = solve(Sigma.normal)
invSigmaKnots.fat = solve(Sigma.fat)

# Project
proj.norm = Sigma21.normal %*% invSigmaKnots.norm %*% matrix(re.norm,ncol=1)
proj.fat = Sigma21.fat %*% invSigmaKnots.fat %*% matrix(re.fat,ncol=1)

# Combine to single RF
pfat = 0.05
proj = (1-pfat)*proj.norm + pfat*proj.fat
```

Next, we'll simulate 3 kinds of data from the spatial mixture: gaussian, binomial, and poisson counts.

```
nPoints = 250
#indices = sample(seq(1,nLocs), size=nPoints, replace=T)
indices = seq(1,nLocs)
# assume no observation error, so observed/simulated gaussian data are same
x = grid[indices,]
y.gaussian = proj[indices,1]

y.binomial = rbinom(nPoints, size=1, prob=plogis(proj[indices,1]))

y.poisson = rpois(nPoints, lambda = exp(proj[indices,1]))

diagKnots = diag(nKnots)
muZeros = rep(0, nKnots)
```

Estimating normal mixture

The most important parameter we're interested in trying to recover is the contribution of the mixtures, and secondarily it's important to recover the variances scaling how extreme those events are.

As a first pass, we'll do it in JAGS, and we'll switch to TMB for the faster models linked with INLA.

```
jagsscript = cat("
model {
    # priors on parameters for gaussian process
    gp_scaleInv ~ dgamma(0.01,0.01); # shared btw normal/fat
    gp_scale <- 1/gp_scaleInv;
    gp_sigmaSqInv_fat ~ dgamma(0.01,0.01); # shared btw normal/fat
    gp_sigmaSq_fat <- 1/gp_sigmaSqInv_fat;
    gp_sigmaSqInv_norm ~ dgamma(0.01,0.01); # shared btw normal/fat
    gp_sigmaSq_norm <- 1/gp_sigmaSqInv_norm;</pre>
```

```
gp_jitterSqInv ~ dgamma(0.01,0.01); # shared btw normal/fat
      gp_jitterSq <- 1/gp_jitterSqInv;</pre>
      # This builds the 2 cov matrices needed
      # SigmaKnots is the COV matrix btween knots
      # SigmaOffDiag is the COV matrix between new locations and knots, e.g. it's (100 x 10)
      for(i in 1:nKnots) {
        for(j in 1:nKnots) {
           # this adds some jitter to the diagonal but not the off-diags
           SigmaKnotsNorm[i,j] <- (1-diagKnots[i,j]) * gp_sigmaSq_norm * exp(-gp_scale * distKnotsSq[i,j]) + gp_sigmaSq[i,j] + gp_s
          SigmaKnotsFat[i,j] <- (1-diagKnots[i,j]) * gp_sigmaSq_fat * exp(-gp_scale * distKnotsSq[i,j]) + di
        }
      }
      for(i in 1:nLocs) {
        for(j in 1:nKnots) {
           SigmaOffDiagNorm[i,j] <- gp_sigmaSq_norm * exp(-gp_scale * distKnots21Sq[i,j]);
          SigmaOffDiagFat[i,j] <- gp_sigmaSq_fat * exp(-gp_scale * distKnots21Sq[i,j]);
       invSigmaKnotsNorm <- inverse(SigmaKnotsNorm[,]); # inverse of matrix for projection and mvn distribu
       invSigmaKnotsFat <- inverse(SigmaKnotsFat[,]); # inverse of matrix for projection and mvn distributi
      # mixture contribution of fat tailed events
      pfat ~ dunif(0,0.3);
      # Spatial random effects
       spatialEffectsKnotsNorm[1:nKnots,1] ~ dmnorm(muZeros, invSigmaKnotsNorm);
       spatialEffectsKnotsFat[1:nKnots,1] ~ dmnorm(muZeros, invSigmaKnotsFat);
       spatialEffects[1:nLocs,1] <- (1-pfat)*((SigmaOffDiagNorm %*% invSigmaKnotsNorm) %*% spatialEffectsKn
       # evaluate the likelihood
      resid.tau ~ dgamma(0.001,0.001);
      for(i in 1:nPoints) {
                pred[i] <- spatialEffects[indices[i], 1];</pre>
                y.gaussian[i] ~ dnorm(pred[i], resid.tau);
} ",file="recover_rf_gaussian.txt")
```

Run the model,

```
jags.data = list("nLocs","nKnots", "y.gaussian", "distKnotsSq", "distKnots21Sq","diagKnots","muZeros","
    jags.params=c("pfat","gp_sigmaSq_fat","gp_sigmaSq_norm","spatialEffectsKnotsNorm","spatialEffectsKno
jagsmodel.gaussian = jags.parallel(jags.data, inits=NULL, parameters.to.save=jags.params, model.file="r
```

Estimating binomial mixture

We can also extend this to the binomial distribution by changing a single line in the JAGS script.

```
jagsscript = cat("
model {
      # priors on parameters for gaussian process
     gp_scaleInv ~ dgamma(0.01,0.01); # shared btw normal/fat
     gp_scale <- 1/gp_scaleInv;</pre>
     gp_sigmaSqInv_fat ~ dgamma(0.01,0.01); # not shared btw normal/fat
      gp_sigmaSq_fat <- 1/gp_sigmaSqInv_fat;</pre>
     gp_sigmaSqInv_norm ~ dgamma(0.01,0.01); # not shared btw normal/fat
     gp_sigmaSq_norm <- 1/gp_sigmaSqInv_norm;</pre>
     gp_jitterSqInv ~ dgamma(0.01,0.01); # shared btw normal/fat
     gp_jitterSq <- 1/gp_jitterSqInv;</pre>
     # This builds the 2 cov matrices needed
      # SigmaKnots is the COV matrix btween knots
      # SigmaOffDiag is the COV matrix between new locations and knots, e.g. it's (100 x 10)
     for(i in 1:nKnots) {
        for(j in 1:nKnots) {
          # this adds some jitter to the diagonal but not the off-diags
          SigmaKnotsNorm[i,j] <- (1-diagKnots[i,j]) * gp_sigmaSq_norm * exp(-gp_scale * distKnotsSq[i,j]) + gp_s
          SigmaKnotsFat[i,j] <- (1-diagKnots[i,j]) * gp_sigmaSq_fat * exp(-gp_scale * distKnotsSq[i,j]) + di
     for(i in 1:nLocs) {
        for(j in 1:nKnots) {
          SigmaOffDiagNorm[i,j] <- gp_sigmaSq_norm * exp(-gp_scale * distKnots21Sq[i,j]);</pre>
         SigmaOffDiagFat[i,j] <- gp_sigmaSq_fat * exp(-gp_scale * distKnots21Sq[i,j]);
      }
      invSigmaKnotsNorm <- inverse(SigmaKnotsNorm[,]); # inverse of matrix for projection and mvn distribu
      invSigmaKnotsFat <- inverse(SigmaKnotsFat[,]); # inverse of matrix for projection and mvn distributi
      # mixture contribution of fat tailed events
     pfat ~ dunif(0,0.3);
      # Spatial random effects
      spatialEffectsKnotsNorm[1:nKnots,1] ~ dmnorm(muZeros, invSigmaKnotsNorm);
      spatialEffectsKnotsFat[1:nKnots,1] ~ dmnorm(muZeros, invSigmaKnotsFat);
      spatialEffects[1:nLocs,1] <- (1-pfat)*((SigmaOffDiagNorm %*% invSigmaKnotsNorm) %*% spatialEffectsKn
      # evaluate the likelihood
      for(i in 1:nPoints) {
              logit(pred[i]) <- min(max(spatialEffects[indices[i], 1],-20),20);</pre>
              y.binomial[i] ~ dbern(pred[i]);
} ",file="recover_rf_binomial.txt")
```

Run the model,

```
jags.data = list("nLocs", "nKnots", "y.binomial", "distKnotsSq", "distKnots21Sq", "diagKnots", "muZeros","
    jags.params=c("pfat", "gp_sigmaSq_fat", "gp_sigmaSq_norm", "spatialEffectsKnotsNorm", "spatialEffectsKno
jagsmodel.binomial = jags.parallel(jags.data, inits=NULL, parameters.to.save=jags.params, model.file="r
```

Estimating Poisson mixture

We can also extend this to the Poisson distribution by changing a single line in the JAGS script.

```
jagsscript = cat("
model {
      # priors on parameters for gaussian process
      gp_scaleInv ~ dgamma(0.01,0.01); # shared btw normal/fat
      gp_scale <- 1/gp_scaleInv;</pre>
      gp_sigmaSqInv_fat ~ dgamma(0.01,0.01); # shared btw normal/fat
      gp_sigmaSq_fat <- 1/gp_sigmaSqInv_fat;</pre>
      gp_sigmaSqInv_norm ~ dgamma(0.01,0.01); # shared btw normal/fat
      gp_sigmaSq_norm <- 1/gp_sigmaSqInv_norm;</pre>
      gp_jitterSqInv ~ dgamma(0.01,0.01); # shared btw normal/fat
      gp_jitterSq <- 1/gp_jitterSqInv;</pre>
      # This builds the 2 cov matrices needed
      # SigmaKnots is the COV matrix btween knots
      # SigmaOffDiag is the COV matrix between new locations and knots, e.g. it's (100 x 10)
      for(i in 1:nKnots) {
        for(j in 1:nKnots) {
          # this adds some jitter to the diagonal but not the off-diags
          SigmaKnotsNorm[i,j] <- (1-diagKnots[i,j]) * gp_sigmaSq_norm * exp(-gp_scale * distKnotsSq[i,j]) + of the sigmaKnotsNorm[i,j] <- (1-diagKnots[i,j]) + of the sigmaKnotsNorm[i,j] <- (1-diagKnots[i,j]) * gp_sigmaSq_norm * exp(-gp_scale * distKnotsSq[i,j]) + of the sigmaKnotsNorm[i,j] <- (1-diagKnots[i,j]) * gp_sigmaSq_norm * exp(-gp_scale * distKnotsSq[i,j]) + of the sigmaKnotsNorm[i,j] <- (1-diagKnots[i,j]) * gp_sigmaSq_norm * exp(-gp_scale * distKnotsSq[i,j]) + of the sigmaKnotsNorm[i,j] <- (1-diagKnots[i,j]) * gp_sigmaSq_norm * exp(-gp_scale * distKnotsSq[i,j]) + of the sigmaKnotsNorm[i,j] <- (1-diagKnots[i,j]) * gp_sigmaSq_norm * exp(-gp_scale * distKnotsSq[i,j]) + of the sigmaSq_norm * exp(-gp_scale * distKnotsSq[i,j]) + of the sigmaSq[i,j] + of the sigmaSq[i,j] + of the sigmaSq_norm * exp(-gp_scale * distKnotsSq[i,
          SigmaKnotsFat[i,j] <- (1-diagKnots[i,j]) * gp_sigmaSq_fat * exp(-gp_scale * distKnotsSq[i,j]) + di
        }
      }
      for(i in 1:nLocs) {
        for(j in 1:nKnots) {
          SigmaOffDiagNorm[i,j] <- gp_sigmaSq_norm * exp(-gp_scale * distKnots21Sq[i,j]);
          SigmaOffDiagFat[i,j] <- gp_sigmaSq_fat * exp(-gp_scale * distKnots21Sq[i,j]);
      invSigmaKnotsNorm <- inverse(SigmaKnotsNorm[,]); # inverse of matrix for projection and mvn distribu
      invSigmaKnotsFat <- inverse(SigmaKnotsFat[,]); # inverse of matrix for projection and mvn distributi
      # mixture contribution of fat tailed events
      pfat ~ dunif(0,0.3);
      # Spatial random effects
      spatialEffectsKnotsNorm[1:nKnots,1] ~ dmnorm(muZeros, invSigmaKnotsNorm);
      spatialEffectsKnotsFat[1:nKnots,1] ~ dmnorm(muZeros, invSigmaKnotsFat);
      spatialEffects[1:nLocs,1] <- (1-pfat)*((SigmaOffDiagNorm %*% invSigmaKnotsNorm) %*% spatialEffectsKn
      # evaluate the likelihood
      for(i in 1:nPoints) {
               y.poisson[i] ~ dpois(exp(spatialEffects[indices[i], 1]));
} ",file="recover_rf_poisson.txt")
```

Run the model,

```
jags.data = list("nLocs", "nKnots", "y.poisson", "distKnotsSq", "distKnots21Sq", "diagKnots", "muZeros", "interface of the control of th
           jags.params=c("pfat","gp_sigmaSq_fat","gp_sigmaSq_norm","spatialEffectsKnotsNorm","spatialEffectsKno
jagsmodel.poisson = jags.parallel(jags.data, inits=NULL, parameters.to.save=jags.params, model.file="re
par(mfrow=c(2,2),mgp=c(2,1,0),mai=c(0.5,0.5,0.3,0.1))
attach.jags(jagsmodel.gaussian)
## The following object is masked _by_ .GlobalEnv:
##
##
                       pfat
hist(pfat,40,col="grey",main="Gaussian",xlab="P(fat tail)")
attach.jags(jagsmodel.binomial)
## The following object is masked _by_ .GlobalEnv:
##
##
                       pfat
hist(pfat,40,col="grey",main="Binomial",xlab="P(fat tail)")
attach.jags(jagsmodel.poisson)
## The following object is masked _by_ .GlobalEnv:
##
##
                       pfat
hist(pfat,40,col="grey",main="Poisson",xlab="P(fat tail)")
```

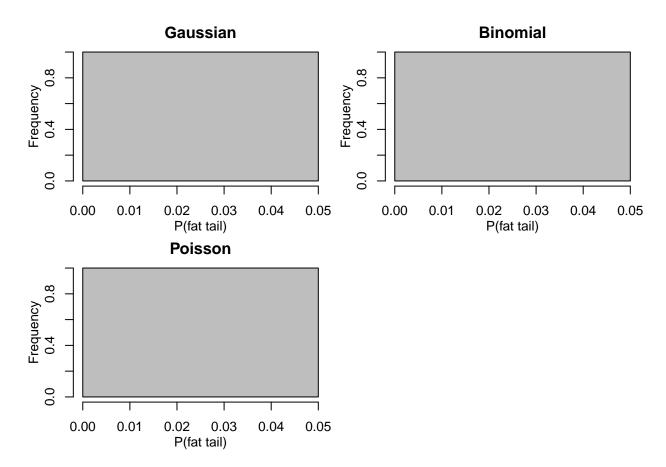


Figure 2: