

Proper/Non-intrinsic Besag model for spatial effects

Parametrization

The proper version of the Besag model for random vector $\mathbf{x} = (x_1, \dots, x_n)$ is defined as

$$x_i | x_{-i}, \tau, d \sim \mathcal{N} \left(\frac{1}{d + n_i} \sum_{i \sim j} x_j, \frac{1}{\tau(d + n_i)} \right) \quad (1)$$

where n_i is the number of neighbours of node i , $i \sim j$ indicates that the two nodes i and j are neighbours, $d > 0$ is an extra term added on the diagonal controlling the “properness” and $\tau > 0$ is a “precision-like” (or scaling) parameter.

This parameterisation corresponds to this precision matrix $Q = (Q_{ij})$, where for $j \neq i$

$$Q_{ii} = \tau(n_i + d) \quad \text{and} \quad Q_{ij} = -\tau.$$

Hyperparameters

The precision parameter τ is represented as

$$\theta_1 = \log \tau$$

and the prior is defined on θ_1 . The diagonal parameter d is represented as

$$\theta_2 = \log d$$

and the prior is defined on θ_2 .

Specification

The besag model is specified inside the `f()` function as

```
f(<whatever>, model="besagproper", graph=<graph>,  
  hyper=<hyper>)
```

The neighbourhood structure of \mathbf{x} is passed to the program through the `graph` argument. The structure of this file is described below.

Hyperparameter spesification and default values

hyper

theta1

```
name log precision  
short.name prec  
prior loggamma  
param 1 5e-04  
initial 2  
fixed FALSE  
to.theta function(x) log(x)  
from.theta function(x) exp(x)
```

theta2

```
name log diagonal
```

```

    short.name diag
    prior loggamma
    param 1 1
    initial 1
    fixed FALSE
    to.theta function(x) log(x)
    from.theta function(x) exp(x)

constr FALSE

nrow.ncol FALSE

augmented FALSE

aug.factor 1

aug.constr

n.div.by

n.required TRUE

set.default.values TRUE

status experimental

pdf besagproper

```

Example

```

## pick a graph
graph = system.file("demodata/germany.graph", package="INLA")
g = inla.read.graph(graph.file)

## we will use replicated samples in our testing
nrep = 5

## make life easy; use dense matrix algebra
d = 1.0
tau = 1.0
Q = matrix(0, g$n, g$n)
diag(Q) = tau * (d + g$nnbs)
for(i in 1:g$n) {
  if (g$nnbs[i] > 0) {
    Q[i, g$nnbs[[i]]] = -tau
    Q[g$nnbs[[i]], i] = -tau
  }
}
R = chol(Q) ## 'chol' returns the upper triangular

## simulate data with replications
y = c()
for(i in 1:nrep) {
  y = c(y, backsolve(R, rnorm(g$n)))
}

```

```

}

i = rep(1:g$n, nrep)
replicate = rep(1:nrep, each = g$n)
formula = y ~ f(i, model="besagproper", graph = graph,
               replicate=replicate,
               hyper = list(diag = list(param = c(1, 1)))) -1

## use 'exact' observations, so we fix the noise precisin to a high
## value
r = inla(formula,
          data = data.frame(y, i, replicate),
          family = "gaussian",
          control.family = list(
            hyper = list(
              prec = list(
                initial = 10,
                fixed=TRUE))))

```

Notes

If $d = 0$ and the parameter `rankdef=1` is set, then this model corresponds to the **besag** model. `constr=FALSE` is default for this model.