



LaplacesDemon Examples

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Abstract

The **LaplacesDemon** package in R enables Bayesian inference with any Bayesian model, provided the user specifies the likelihood. This vignette is a compendium of examples of how to specify different model forms.

Keywords: ~Bayesian, Bayesian Inference, Laplace's Demon, LaplacesDemon, R, STATISTICAT.

LaplacesDemon (Hall 2011), usually referred to as Laplace's Demon, is an R package that is available on CRAN (R Development Core Team 2011). A formal introduction to Laplace's Demon is provided in an accompanying vignette entitled "**LaplacesDemon** Tutorial", and an introduction to Bayesian inference is provided in the "Bayesian Inference" vignette.

The purpose of this document is to provide users of the **LaplacesDemon** package with examples of a variety of Bayesian methods. To conserve space, the examples are not worked out in detail, and only the minimum of necessary materials is provided for using the various methodologies. Necessary materials include the form expressed in notation, data (which is often simulated), initial values, and the `Model` function.

Notation in this vignette follows these standards: Greek letters represent parameters, lower case letters represent indices, lower case bold face letters represent scalars or vectors, probability distributions are represented with calligraphic font, upper case letters represent index limits, and upper case bold face letters represent matrices.

This vignette will grow over time as examples of more methods become included. Contributed examples are welcome. Please send contributed examples in a similar format in an email to laplacesdemon@statisticat.com for review and testing. All accepted contributions are, of course, credited.

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1. ANCOVA

This example is essentially the same as the two-way ANOVA (see [3](#)), except that a covariate $\mathbf{X}_{,3}$ has been added, and its parameter is δ .

1.1. Form

$$\begin{aligned}
 \mathbf{y}_i &\sim \mathcal{N}(\mu_i, \sigma_1^2) \\
 \mu_i &= \alpha + \beta[\mathbf{X}_{i,1}] + \gamma[\mathbf{X}_{i,2}] + \delta\mathbf{X}_{i,2}, \quad i = 1, \dots, N \\
 \epsilon_i &= \mathbf{y}_i - \mu_i \\
 \alpha &\sim \mathcal{N}(0, 1000) \\
 \beta_j &\sim \mathcal{N}(0, \sigma_2^2), \quad j = 1, \dots, (J-1) \\
 \beta_J &= -\sum_{j=1}^{J-1} \beta_j \\
 \gamma_k &\sim \mathcal{N}(0, \sigma_3^2), \quad k = 1, \dots, (K-1) \\
 \gamma_K &= -\sum_{k=1}^{K-1} \gamma_k \\
 \delta &\sim \mathcal{N}(0, 1000) \\
 \sigma_m &\sim \mathcal{HC}(25), \quad m = 1, \dots, 3
 \end{aligned}$$

1.2. Data

```

N <- 100
J <- 5 #Number of levels in factor (treatment) 1
K <- 3 #Number of levels in factor (treatment) 2
X <- matrix(cbind(round(runif(N,0.5,J+0.49)),round(runif(N,0.5,K+0.49))),

```

```

      runif(N,-2,2)), N, 3)
alpha <- runif(1,-1,1)
beta <- runif(J,-2,2)
beta[J] <- -sum(beta[1:(J-1)])
gamma <- runif(K,-2,2)
gamma[J] <- -sum(gamma[1:(K-1)])
delta <- runif(1,-2,2)
y <- alpha + beta[X[,1]] + gamma[X[,2]] + delta*X[,3] + rnorm(N,0,0.1)
mon.names <- c("LP","beta[5]","gamma[3]","sigma[1]","sigma[2]","sigma[3]",
  "s.beta","s.gamma","s.epsilon")
parm.names <- parm.names(list(alpha=0, beta=rep(0,J-1), gamma=rep(0,K-1),
  delta=0, log.sigma=rep(0,3)))
MyData <- list(J=J, K=K, N=N, X=X, mon.names=mon.names,
  parm.names=parm.names, y=y)

```

1.3. Initial Values

```
Initial.Values <- c(0, rep(0,(J-1)), rep(0,(K-1)), 0, rep(log(1),3))
```

1.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  alpha <- parm[1]
  beta <- rep(NA,Data$J)
  beta[1:(Data$J-1)] <- parm[2:Data$J]
  beta[J] <- -sum(beta[1:(Data$J-1)]) #Sum-to-zero constraint
  gamma <- rep(NA,Data$K)
  gamma[1:(Data$K-1)] <- parm[grepl("gamma", Data$parm.names)]
  gamma[K] <- -sum(gamma[1:(Data$K-1)]) #Sum-to-zero constraint
  delta <- parm[grepl("delta", Data$parm.names)]
  sigma <- exp(parm[grepl("log.sigma", Data$parm.names)])
  ### Log(Prior Densities)
  alpha.prior <- dnorm(alpha, 0, sqrt(1000), log=TRUE)
  beta.prior <- sum(dnorm(beta, 0, sigma[2], log=TRUE))
  gamma.prior <- sum(dnorm(gamma, 0, sigma[3], log=TRUE))
  delta.prior <- dnorm(delta, 0, sqrt(1000), log=TRUE)
  sigma.prior <- sum(dhalfcauchy(sigma, 25, log=TRUE))
  ### Log-Likelihood
  mu <- alpha + beta[Data$X[,1]] + gamma[Data$X[,2]] +
    delta*Data$X[,3]
  LL <- sum(dnorm(Data$y, mu, sigma[1], log=TRUE))
  ### Variance Components
  s.beta <- sd(beta)
  s.gamma <- sd(gamma)

```

```

s.epsilon <- sd(Data$y - mu)
### Log-Posterior
LP <- LL + alpha.prior + beta.prior + gamma.prior + delta.prior +
  sigma.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP, beta[Data$J],
  gamma[Data$K], sigma, s.beta, s.gamma, s.epsilon), yhat=mu,
  parm=parm)
return(Modelout)
}

```

2. ANOVA, One-Way

When $J = 2$, this is a Bayesian form of a t-test.

2.1. Form

$$\begin{aligned}
 \mathbf{y} &\sim \mathcal{N}(\mu, \sigma_1^2) \\
 \mu_i &= \alpha + \beta[\mathbf{x}_i], \quad i = 1, \dots, N \\
 \alpha &\sim \mathcal{N}(0, 1000) \\
 \beta_j &\sim \mathcal{N}(0, \sigma_2^2), \quad j = 1, \dots, (J-1) \\
 \beta_J &= -\sum_{j=1}^{J-1} \beta_j \\
 \sigma_{1:2} &\sim \mathcal{HC}(25)
 \end{aligned}$$

2.2. Data

```

N <- 100
J <- 5
x <- round(runif(N, 0.5, J+0.49))
alpha <- runif(1,-1,1)
beta <- runif(J,-2,2)
beta[J] <- -sum(beta[1:(J-1)])
y <- rep(NA, N)
for (i in 1:N) {y[i] <- alpha + beta[x[i]] + rnorm(1,0,0.2)}
mon.names <- c("LP","beta[1]","sigma[1]","sigma[2]")
parm.names <- parm.names(list(alpha=0, beta=rep(0,J-1),
  log.sigma=rep(0,2)))
MyData <- list(J=J, N=N, mon.names=mon.names, parm.names=parm.names, x=x,
  y=y)

```

2.3. Initial Values

```
Initial.Values <- c(0, rep(0,(J-1)), rep(log(1),2))
```

2.4. Model

```
Model <- function(parm, Data)
{
  ### Parameters
  alpha <- parm[1]
  beta <- rep(NA,Data$J)
  beta[1:(Data$J-1)] <- parm[2:Data$J]
  beta[J] <- -sum(beta[1:(Data$J-1)]) #Sum-to-zero constraint
  sigma <- exp(parm[grep("log.sigma", Data$parm.names)])
  ### Log(Prior Densities)
  alpha.prior <- dnorm(alpha, 0, sqrt(1000), log=TRUE)
  beta.prior <- sum(dnorm(beta, 0, sigma[2], log=TRUE))
  sigma.prior <- sum(dhalfcauchy(sigma, 25, log=TRUE))
  ### Log-Likelihood
  mu <- alpha + beta[Data$x]
  LL <- sum(dnorm(Data$y, mu, sigma[1], log=TRUE))
  ### Log-Posterior
  LP <- LL + alpha.prior + beta.prior + sigma.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,beta[Data$J],
    sigma), yhat=mu, parm=parm)
  return(Modelout)
}
```

3. ANOVA, Two-Way

In this representation, σ^m are the superpopulation variance components, **s.beta** and **s.gamma** are the finite-population within-variance components of the factors or treatments, and **s.error** is the finite-population between-variance component.

3.1. Form

$$\begin{aligned}
 \mathbf{y}_i &\sim \mathcal{N}(\mu_i, \sigma_1^2) \\
 \mu_i &= \alpha + \beta[\mathbf{X}_{i,1}] + \gamma[\mathbf{X}_{i,2}], \quad i = 1, \dots, N \\
 \epsilon_i &= \mathbf{y}_i - \mu_i \\
 \alpha &\sim \mathcal{N}(0, 1000) \\
 \beta_j &\sim \mathcal{N}(0, \sigma_2^2), \quad j = 1, \dots, (J-1) \\
 \beta_J &= -\sum_{j=1}^{J-1} \beta_j
 \end{aligned}$$

$$\gamma_k \sim \mathcal{N}(0, \sigma_3^2), \quad k = 1, \dots, (K-1)$$

$$\gamma_K = - \sum_{k=1}^{K-1} \gamma_k$$

$$\sigma_m \sim \mathcal{HC}(25), \quad m = 1, \dots, 3$$

3.2. Data

```

N <- 100
J <- 5 #Number of levels in factor (treatment) 1
K <- 3 #Number of levels in factor (treatment) 2
X <- matrix(cbind(round(runif(N, 0.5, J+0.49)),round(runif(N,0.5,K+0.49))),
            N, 2)
alpha <- runif(1,-1,1)
beta <- runif(J,-2,2)
beta[J] <- -sum(beta[1:(J-1)])
gamma <- runif(K,-2,2)
gamma[J] <- -sum(gamma[1:(K-1)])
y <- alpha + beta[X[,1]] + gamma[X[,2]] + rnorm(1,0,0.1)
mon.names <- c("LP","beta[5]","gamma[3]","sigma[1]","sigma[2]","sigma[3]",
              "s.beta","s.gamma","s.epsilon")
parm.names <- parm.names(list(alpha=0, beta=rep(0,J-1), gamma=rep(0,K-1),
                              log.sigma=rep(0,3)))
MyData <- list(J=J, K=K, N=N, X=X, mon.names=mon.names,
              parm.names=parm.names, y=y)

```

3.3. Initial Values

```
Initial.Values <- c(0, rep(0,(J-1)), rep(0,(K-1)), rep(log(1),3))
```

3.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  alpha <- parm[1]
  beta <- rep(NA,Data$J)
  beta[1:(Data$J-1)] <- parm[2:Data$J]
  beta[J] <- -sum(beta[1:(Data$J-1)]) #Sum-to-zero constraint
  gamma <- rep(NA,Data$K)
  gamma[1:(Data$K-1)] <- parm[grep("gamma", Data$parm.names)]
  gamma[K] <- -sum(gamma[1:(Data$K-1)]) #Sum-to-zero constraint
  sigma <- exp(parm[grep("log.sigma", Data$parm.names)])
  ### Log(Prior Densities)
  alpha.prior <- dnorm(alpha, 0, sqrt(1000), log=TRUE)
  beta.prior <- sum(dnorm(beta, 0, sigma[2], log=TRUE))

```

```

gamma.prior <- sum(dnorm(gamma, 0, sigma[3], log=TRUE))
sigma.prior <- sum(dhalfcauchy(sigma, 25, log=TRUE))
### Log-Likelihood
mu <- alpha + beta[Data$X[,1]] + gamma[Data$X[,2]]
LL <- sum(dnorm(Data$y, mu, sigma[1], log=TRUE))
### Variance Components
s.beta <- sd(beta)
s.gamma <- sd(gamma)
s.epsilon <- sd(Data$y - mu)
### Log-Posterior
LP <- LL + alpha.prior + beta.prior + gamma.prior +
      sigma.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP, beta[Data$J],
      gamma[Data$K], sigma, s.beta, s.gamma, s.epsilon), yhat=mu,
      parm=parm)
return(Modelout)
}

```

4. Autoregression, AR(1)

4.1. Form

$$\mathbf{y}_t \sim \mathcal{N}(\mu_{t-1}, \sigma^2), \quad t = 2, \dots, (T-1)$$

$$\mathbf{y}_T^{new} \sim \mathcal{N}(\mu_T, \sigma^2)$$

$$\mu_t = \alpha + \phi \mathbf{y}_t, \quad t = 1, \dots, T$$

$$\alpha \sim \mathcal{N}(0, 1000)$$

$$\phi \sim \mathcal{N}(0, 1000)$$

$$\sigma \sim \mathcal{HC}(25)$$

4.2. Data

```

T <- 100
y <- rep(0, T)
y[1] <- 0
for (t in 2:T) {y[t] <- y[t-1] + rnorm(1, 0, 0.1)}
mon.names <- c("LP", "sigma", paste("mu[", T, "]", sep=""))
parm.names <- c("alpha", "phi", "log.sigma")
MyData <- list(T=T, mon.names=mon.names, parm.names=parm.names, y=y)

```


4.3. Initial Values

```
Initial.Values <- c(rep(0,2), log(1))
```

4.4. Model

```
Model <- function(parm, Data)
{
  ### Parameters
  alpha <- parm[1]; phi <- parm[2]; sigma <- exp(parm[3])
  ### Log(Prior Densities)
  alpha.prior <- dnorm(alpha, 0, sqrt(1000), log=TRUE)
  phi.prior <- dnorm(phi, 0, sqrt(1000), log=TRUE)
  sigma.prior <- dhalfcauchy(sigma, 25, log=TRUE)
  ### Log-Likelihood
  mu <- alpha + phi*Data$y
  LL <- sum(dnorm(Data$y[2:(Data$T-1)], mu[1:(Data$T-2)],
    sigma, log=TRUE))
  ### Log-Posterior
  LP <- LL + alpha.prior + phi.prior + sigma.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,sigma,mu[Data$T]),
    yhat=mu, parm=parm)
  return(Modelout)
}
```

5. Binary Logit

5.1. Form

$$\mathbf{y} \sim \mathcal{BERN}(\eta)$$

$$\eta = \frac{1}{1 + \exp(-\mu)}$$

$$\mu = \mathbf{X}\beta$$

$$\beta_j \sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J$$

5.2. Data

```
data(demonsnacks)
N <- NROW(demonsnacks)
J <- 3
y <- ifelse(demonsnacks$Calories <= 137, 0, 1)
X <- cbind(1, as.matrix(demonsnacks[,c(7,8)]))
for (j in 2:J) {X[,j] <- CenterScale(X[,j])}
```

```

mon.names <- "LP"
parm.names <- parm.names(list(beta=rep(0,J)))
MyData <- list(J=J, X=X, mon.names=mon.names, parm.names=parm.names, y=y)

```

5.3. Initial Values

```
Initial.Values <- rep(0,J)
```

5.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  beta <- parm[1:Data$J]
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  mu <- tcrossprod(beta, Data$X)
  eta <- invlogit(mu)
  ### Log-Likelihood
  LL <- sum(dbern(Data$y, eta, log=TRUE))
  yrep <- ifelse(eta >= (sum(Data$y)/length(Data$y)),1,0)
  ### Log-Posterior
  LP <- LL + beta.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=LP,
    yhat=yrep, parm=parm)
  return(Modelout)
}

```

6. Binary Probit

6.1. Form

$$\mathbf{y} \sim \mathcal{BERN}(\mathbf{p})$$

$$\mathbf{p} = \phi(\mu)$$

$$\mu = \mathbf{X}\beta \in [-10, 10]$$

$$\beta_j \sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J$$

6.2. Data

```

data(demonsnacks)
N <- NROW(demonsnacks)

```

```

J <- 3
y <- ifelse(demonsnacks$Calories <= 137, 0, 1)
X <- cbind(1, as.matrix(demonsnacks[,c(7,8)]))
for (j in 2:J) {X[,j] <- CenterScale(X[,j])}
mon.names <- "LP"
parm.names <- parm.names(list(beta=rep(0,J)))
MyData <- list(J=J, X=X, mon.names=mon.names, parm.names=parm.names, y=y)

```

6.3. Initial Values

```
Initial.Values <- rep(0,J)
```

6.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  beta <- parm[1:Data$J]
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  ### Log-Likelihood
  mu <- tcrossprod(beta, Data$X)
  mu <- interval(mu, -10, 10)
  p <- pnorm(mu)
  LL <- sum(dbern(Data$y, p, log=TRUE))
  yrep <- ifelse(p >= (sum(Data$y)/length(Data$y)),1,0)
  ### Log-Posterior
  LP <- LL + beta.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=LP, yhat=yrep, parm=parm)
  return(Modelout)
}

```

7. Binomial Logit

7.1. Form

$$\mathbf{y} \sim \mathcal{BLN}(\mathbf{p}, \mathbf{n})$$

$$\mathbf{p} = \frac{1}{1 + \exp(-\mu)}$$

$$\mu = \beta_1 + \beta_2 \mathbf{x}$$

$$\beta_j \sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J$$

where ϕ is the inverse CDF, and $J=2$.

7.2. Data

```
#10 Trials
exposed <- c(100,100,100,100,100,100,100,100,100,100)
deaths <- c(10,20,30,40,50,60,70,80,90,100)
dose <- c(1,2,3,4,5,6,7,8,9,10)
J <- 2 #Number of parameters
mon.names <- "LP"
parm.names <- c("beta[1]","beta[2]")
MyData <- list(J=J, n=exposed, mon.names=mon.names, parm.names=parm.names,
               x=dose, y=deaths)
```

7.3. Initial Values

```
Initial.Values <- rep(0,J)
```

7.4. Model

```
Model <- function(parm, Data)
{
  ### Parameters
  beta <- parm
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  ### Log-Likelihood
  mu <- beta[1] + beta[2]*Data$x
  p <- invlogit(mu)
  LL <- sum(dbinom(Data$y, Data$n, p, log=TRUE))
  yrep <- p * Data$n
  ### Log-Posterior
  LP <- LL + beta.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=LP, yhat=yrep, parm=parm)
  return(Modelout)
}
```

8. Binomial Probit

8.1. Form

$$\mathbf{y} \sim \mathcal{BIN}(\mathbf{p}, \mathbf{n})$$

$$\mathbf{p} = \phi(\boldsymbol{\mu})$$

$$\boldsymbol{\mu} = \beta_1 + \beta_2 \mathbf{x} \in [-10, 10]$$

$$\beta_j \sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J$$

where ϕ is the inverse CDF, and $J=2$.

8.2. Data

```
#10 Trials
exposed <- c(100,100,100,100,100,100,100,100,100,100)
deaths <- c(10,20,30,40,50,60,70,80,90,100)
dose <- c(1,2,3,4,5,6,7,8,9,10)
J <- 2 #Number of parameters
mon.names <- "LP"
parm.names <- c("beta[1]","beta[2]")
MyData <- list(J=J, n=exposed, mon.names=mon.names, parm.names=parm.names,
               x=dose, y=deaths)
```

8.3. Initial Values

```
Initial.Values <- rep(0,J)
```

8.4. Model

```
Model <- function(parm, Data)
{
  ### Parameters
  beta <- parm
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  ### Log-Likelihood
  mu <- beta[1] + beta[2]*Data$x
  mu <- interval(mu, -10, 10)
  p <- pnorm(mu)
  LL <- sum(dbinom(Data$y, Data$n, p, log=TRUE))
  yrep <- p * Data$n
  ### Log-Posterior
  LP <- LL + beta.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=LP, yhat=yrep,
                  parm=parm)
  return(Modelout)
}
```

9. Cluster Analysis

This is a parametric model-based cluster analysis, also called a finite mixture model or latent class cluster analysis.

9.1. Form

$$\mathbf{Y}_{i,j} \sim \mathcal{N}(\mu_{\theta[i],j}, \sigma_{\theta[i]}^2), \quad i = 1, \dots, N, \quad j = 1, \dots, J$$

$$\theta_i = \text{Max}(\mathbf{p}_{i,1:C})$$

$$\mathbf{p}_{i,c} = \frac{\delta_{i,c}}{\sum_{c=1}^C \delta_{i,c}}$$

$$\pi_{1:C} \sim \mathcal{D}(\alpha_{1:C})$$

$$\pi_c = \frac{\sum_{i=1}^N \delta_{i,c}}{\sum \delta}$$

$$\alpha_c = 1$$

$$\delta_{i,C} = 1$$

$$\delta_{i,c} \sim \mathcal{N}(\log(\frac{1}{C}), 1000) \in [\exp(-10), \exp(10)], \quad c = 1, \dots, (C-1)$$

$$\mu_{c,j} \sim \mathcal{N}(0, \nu_j^2)$$

$$\sigma_c \sim \mathcal{HC}(25)$$

$$\nu_j \sim \mathcal{HC}(25)$$

9.2. Data

```

C <- 3 #Number of clusters
alpha <- rep(1,C) #Prior probability of cluster proportion
# Create a Y matrix
n <- 100; N <- 15 #Full sample; model sample
J <- 5 #Number of predictor variables
cluster <- round(runif(n,0.5,C+0.49))
centers <- matrix(runif(C*J, 0, 10), C, J)
Y.Full <- matrix(0, n, J)
for (i in 1:n) {for (j in 1:J)
  {Y.Full[i,j] <- rnorm(1,centers[cluster[i],j],1)}}
mean.temp <- colMeans(Y.Full)
sigma.temp <- apply(Y.Full,2,sd)
centers.cs <- (centers - matrix(rep(mean.temp,C), C, J, byrow=TRUE)) /
  (2 * matrix(rep(sigma.temp,C), C, J, byrow=TRUE))
for (j in 1:J) {Y.Full[,j] <- scale(Y.Full[,j],2)}
#summary(Y.Full)
MySample <- sample(1:n, N)
Y <- Y.Full[MySample,]
mon.names <- c("LP", parm.names(list(nu=rep(0,J), pi=rep(0,C),
  sigma=rep(0,C), theta=rep(0,N))))
parm.names <- parm.names(list(log.delta=matrix(0,N,C-1), mu=matrix(0,C,J),
  log.nu=rep(0,J), log.sigma=rep(0,C)))
MyData <- list(C=C, J=J, N=N, Y=Y, alpha=alpha, mon.names=mon.names,
  parm.names=parm.names)

```

9.3. Initial Values

```
Initial.Values <- c(runif(N*(C-1),-1,1), rep(0,C*J), rep(0,J), rep(0,C))
```

9.4. Model

```
Model <- function(parm, Data)
{
  ### Parameters
  delta <- interval(parm[grep("log.delta", Data$parm.names)], -10, 10)
  parm[grep("log.delta", Data$parm.names)] <- delta
  delta <- matrix(c(exp(delta), rep(1, Data$N)), Data$N, Data$C)
  mu <- matrix(parm[grep("mu", Data$parm.names)], Data$C, Data$J)
  nu <- exp(parm[grep("log.nu", Data$parm.names)])
  pi <- colSums(delta) / sum(delta)
  sigma <- exp(parm[grep("log.sigma", Data$parm.names)])
  ### Log(Prior Densities)
  delta.prior <- sum(dtrunc(delta, "norm", a=exp(-10), b=exp(10),
    mean=log(1/Data$C), sd=sqrt(1000), log=TRUE))
  mu.prior <- sum(dnorm(mu, 0, matrix(rep(nu, Data$C), Data$C,
    Data$J, byrow=TRUE), log=TRUE))
  nu.prior <- sum(dhalfcauchy(nu, 25, log=TRUE))
  pi.prior <- ddirichlet(pi, Data$alpha, log=TRUE)
  sigma.prior <- sum(dhalfcauchy(sigma, 25, log=TRUE))
  ### Log-Likelihood
  p <- delta / rowSums(delta)
  theta <- apply(p, 1, which.max)
  LL <- sum(dnorm(Y, mu[theta,], sigma[theta,], log=TRUE))
  Yrep <- mu[theta,]
  ### Log-Posterior
  LP <- LL + delta.prior + mu.prior + nu.prior + pi.prior +
    sigma.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP, nu, pi, sigma, theta),
    yhat=Yrep, parm=parm)
  return(Modelout)
}
```

10. Conditional Autoregression (CAR), Poisson

This CAR example is a slightly modified form of example 7.3 (Model A) in [Congdon \(2003\)](#). The Scottish lip cancer data also appears in the WinBUGS ([Spiegelhalter, Thomas, Best, and Lunn 2003](#)) examples and is a widely analyzed example. The data \mathbf{y} consists of counts for $i = 1, \dots, 56$ counties in Scotland. A single predictor \mathbf{x} is provided. The errors, ϵ , are allowed to include spatial effects as smoothing by spatial effects from areal neighbors. Interactions \mathbf{w} between counties are in terms of dummy indicators for contiguity (areal neighbors). The list of NN areal neighbors is in the *adj* variable, and cumulative positions are in variable *C*. The

vector ϵ_μ is the mean of each area's error, and is a weighted average of errors in contiguous areas.

10.1. Form

$$\begin{aligned}\mathbf{y} &\sim \mathcal{P}(\lambda) \\ \lambda &= \exp(\log(\mathbf{E}) + \beta_1 + \beta_2 \mathbf{x} + \epsilon) \\ \epsilon &\sim \mathcal{N}(\epsilon_\mu, \sigma^2) \\ \epsilon_{\mu[i]} &= \rho \sum_{j=1}^J \mathbf{w}_{i,j} \epsilon_j, \quad i = 1, \dots, N \\ \beta_j &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J \\ \rho &\sim \mathcal{U}(-1, 1) \\ \sigma &\sim \mathcal{HC}(25)\end{aligned}$$

10.2. Data

```
N <- 56 #Number of areas
NN <- 264 #Number of adjacent areas
y <- c(9,39,11,9,15,8,26,7,6,20,13,5,3,8,17,9,2,7,9,7,16,31,11,7,19,15,7,
      10,16,11,5,3,7,8,11,9,11,8,6,4,10,8,2,6,19,3,2,3,28,6,1,1,1,1,0,0)
E <- c( 1.4,8.7,3.0,2.5,4.3,2.4,8.1,2.3,2.0,6.6,4.4,1.8,1.1,3.3,7.8,4.6,
      1.1,4.2,5.5,4.4,10.5,22.7,8.8,5.6,15.5,12.5,6.0,9.0,14.4,10.2,4.8,
      2.9,7.0,8.5,12.3,10.1,12.7,9.4,7.2,5.3,18.8,15.8,4.3,14.6,50.7,8.2,
      5.6,9.3,88.7,19.6,3.4,3.6,5.7,7.0,4.2,1.8) #Expected
x <- c(16,16,10,24,10,24,10,7,7,16,7,16,10,24,7,16,10,7,7,10,7,16,10,7,1,1,
      7,7,10,10,7,24,10,7,7,0,10,1,16,0,1,16,16,0,1,7,1,1,0,1,1,0,1,1,16,10)
adj <- c(5,9,11,19, #Area 1 is adjacent to areas 5, 9, 11, and 19
      7,10, #Area 2 is adjacent to areas 7 and 10
      6,12,
      18,20,28,
      1,11,12,13,19,
      3,8,
      2,10,13,16,17,
      6,
      1,11,17,19,23,29,
      2,7,16,22,
      1,5,9,12,
      3,5,11,
      5,7,17,19,
      31,32,35,
      25,29,50,
      7,10,17,21,22,29,
```



```

7,9,13,16,19,29,
4,20,28,33,55,56,
1,5,9,13,17,
4,18,55,
16,29,50,
10,16,
9,29,34,36,37,39,
27,30,31,44,47,48,55,56,
15,26,29,
25,29,42,43,
24,31,32,55,
4,18,33,45,
9,15,16,17,21,23,25,26,34,43,50,
24,38,42,44,45,56,
14,24,27,32,35,46,47,
14,27,31,35,
18,28,45,56,
23,29,39,40,42,43,51,52,54,
14,31,32,37,46,
23,37,39,41,
23,35,36,41,46,
30,42,44,49,51,54,
23,34,36,40,41,
34,39,41,49,52,
36,37,39,40,46,49,53,
26,30,34,38,43,51,
26,29,34,42,
24,30,38,48,49,
28,30,33,56,
31,35,37,41,47,53,
24,31,46,48,49,53,
24,44,47,49,
38,40,41,44,47,48,52,53,54,
15,21,29,
34,38,42,54,
34,40,49,54,
41,46,47,49,
34,38,49,51,52,
18,20,24,27,56,
18,24,30,33,45,55)
# C has length N+1 and refers to cumulative position (-1) in the adj
# variable. For example, area 1 begins at 0 (position 1-1), and
# area 2 begins at 4 (position 5-1), etc.
C <- c(0,4,6,8,11,16,18,23,24,30,34,38,41,45,48,51,57,63,69,74,77,80,82,
      88,96,99,103,107,111,122,128,135,139,143,152,157,161,166,172,177,182,
      189,195,199,204,208,214,220,224,233,236,240,244,248,253,258,264)
mon.names <- c("LP","sigma")

```

```

parm.names <- parm.names(list(beta=rep(0,2), epsilon=rep(0,N), rho=0,
  log.sigma=0))
MyData <- list(C=C, E=E, N=N, NN=NN, adj=adj, mon.names=mon.names,
  parm.names=parm.names, x=x, y=y)

```

10.3. Initial Values

```
Initial.Values <- c(rep(0,2), rep(0,N), 0, 0)
```

10.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  beta <- parm[1:2]
  epsilon <- parm[grepl("epsilon", Data$parm.names)]
  rho <- interval(parm[grepl("rho", Data$parm.names)], -1, 1)
  parm[grepl("rho", Data$parm.names)] <- rho
  w <- epsilon[Data$adj]
  epsilon.mu <- epsilon
  for (i in 1:N) {
    epsilon.mu[i] <- rho * sum(w[(Data$C[i]+1):(Data$C[i+1])]))
  }
  sigma <- exp(parm[grepl("log.sigma", Data$parm.names)])
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  epsilon.prior <- sum(dnorm(epsilon, epsilon.mu, sigma,
    log=TRUE))
  rho.prior <- dunif(rho, -1, 1, log=TRUE)
  sigma.prior <- dhalfcauchy(sigma, 25, log=TRUE)
  ### Log-Likelihood
  lambda <- exp(log(Data$E) + beta[1] + beta[2]*Data$x/10 + epsilon)
  LL <- sum(dpois(Data$y, lambda, log=TRUE))
  ### Log-Posterior
  LP <- LL + beta.prior + epsilon.prior + rho.prior + sigma.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,sigma), yhat=lambda,
    parm=parm)
  return(Modelout)
}

```

11. Contingency Table

The two-way contingency table, matrix \mathbf{Y} , can easily be extended to more dimensions. For this example, it is vectorized as y , and used like an ANOVA data set. Contingency table \mathbf{Y} has J rows and K columns. The cell counts are fit with Poisson regression, according to intercept α , main effects β_j for each row, main effects γ_k for each column, and interaction effects $\delta_{j,k}$ for dependence effects. An omnibus (all cells) test of independence is done by estimating two models (one with δ , and one without), and a large enough Bayes Factor indicates a

violation of independence when the model with δ fits better than the model without δ . In an ANOVA-like style, main effects contrasts can be used to distinguish rows or groups of rows from each other, as well as with columns. Likewise, interaction effects contrasts can be used to test independence in groups of $\delta_{j,k}$ elements. Finally, single-cell interactions can be used to indicate violations of independence for a given cell, such as when zero is not within its 95% probability interval. Although a little different, this example is similar to a method presented by [Albert \(1997\)](#).

11.1. Form

$$\begin{aligned}
 \mathbf{Y}_{j,k} &\sim \mathcal{P}(\lambda_{j,k}), \quad j = 1, \dots, J, \quad k = 1, \dots, K \\
 \lambda_{j,k} &= \exp(\alpha + \beta_j + \gamma_k + \delta_{j,k}), \quad j = 1, \dots, J, \quad k = 1, \dots, K \\
 \alpha &\sim \mathcal{N}(0, 1000) \\
 \beta_j &\sim \mathcal{N}(0, \beta_\sigma^2), \quad j = 1, \dots, J \\
 \beta_\sigma &\sim \mathcal{HC}(25) \\
 \gamma_k &\sim \mathcal{N}(0, \gamma_\sigma^2), \quad k = 1, \dots, K \\
 \gamma_\sigma &\sim \mathcal{HC}(25) \\
 \delta_{j,k} &\sim \mathcal{N}(0, \delta_\sigma^2) \\
 \delta_\sigma &\sim \mathcal{HC}(25)
 \end{aligned}$$

11.2. Data

```

J <- 4 #Rows
K <- 4 #Columns
Y <- matrix(c(10,20,60,20, 40,30,10,40, 10,10,40,10, 40,50,1,40), J, K,
             dimnames=list(c("Chrysler","Ford","Foreign","GM"),
                           c("I-4","I-6","V-6","V-8")))
y <- as.vector(Y)
N <- length(y) #Cells
r <- rep(1:J, N/J)
c <- rep(1,K)
for (k in 2:K) {c <- c(c, rep(k, K))}
mon.names <- c("LP","beta.sigma","gamma.sigma","delta.sigma")
parm.names <- parm.names(list(alpha=0, beta=rep(0,J), gamma=rep(0,J),
                              log.b.sigma=0, log.g.sigma=0, log.d.sigma=0,
                              delta=matrix(0,J,K)))
MyData <- list(J=J, K=K, N=N, c=c, mon.names=mon.names,
               parm.names=parm.names, r=r, y=y)

```

11.3. Initial Values

```
Initial.Values <- c(0, rep(0,J), rep(0,K), rep(0,3), rep(0,J*K))
```

11.4. Model

```

Model <- function(parm, Data)
{
  ### Hyperparameters
  beta.sigma <- exp(parm[grepl("log.b.sigma", Data$parm.names)])
  gamma.sigma <- exp(parm[grepl("log.g.sigma", Data$parm.names)])
  delta.sigma <- exp(parm[grepl("log.d.sigma", Data$parm.names)])
  ### Parameters
  alpha <- parm[grepl("alpha", Data$parm.names)]
  beta <- parm[grepl("beta", Data$parm.names)]
  gamma <- parm[grepl("gamma", Data$parm.names)]
  delta <- matrix(parm[grepl("delta", Data$parm.names)],
    Data$J, Data$K)
  ### Log(Prior Densities)
  alpha.prior <- dnorm(alpha, 0, sqrt(1000), log=TRUE)
  beta.prior <- sum(dnorm(beta, 0, beta.sigma, log=TRUE))
  beta.sigma.prior <- dhalfcauchy(beta.sigma, 25, log=TRUE)
  gamma.prior <- sum(dnorm(gamma, 0, gamma.sigma, log=TRUE))
  gamma.sigma.prior <- dhalfcauchy(gamma.sigma, 25, log=TRUE)
  delta.prior <- sum(dnorm(delta, 0, delta.sigma, log=TRUE))
  delta.sigma.prior <- dhalfcauchy(delta.sigma, 25, log=TRUE)
  ### Log-Likelihood
  lambda <- exp(alpha + beta[Data$r] + gamma[Data$c] +
    diag(delta[Data$r,Data$c]))
  LL <- sum(dpois(Data$y, lambda, log=TRUE))
  ### Log-Posterior
  LP <- LL + alpha.prior + beta.prior + beta.sigma.prior +
    gamma.prior + gamma.sigma.prior + delta.prior +
    delta.sigma.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP, beta.sigma,
    gamma.sigma, delta.sigma), yhat=lambda, parm=parm)
  return(Modelout)
}

```

12. Discrete Choice, Conditional Logit

12.1. Form

$$\mathbf{y}_i \sim \mathcal{CAT}(\mathbf{p}_{i,1:J}), \quad i = 1, \dots, N, \quad j = 1, \dots, J$$

$$\mathbf{p}_{i,j} = \frac{\phi_{i,j}}{\sum_{j=1}^J \phi_{i,j}}$$

$$\phi = \exp(\mu)$$

$$\mu_{i,j} = \beta_{j,1:K} \mathbf{X}_{i,1:K} + \gamma \mathbf{Z}_{i,1:C} \in [-700, 700], \quad j = 1, \dots, (J-1)$$

$$\begin{aligned}\mu_{i,J} &= \gamma \mathbf{Z}_{i,1:C} \\ \beta_{j,k} &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, (J-1) \\ \gamma_c &\sim \mathcal{N}(0, 1000)\end{aligned}$$

12.2. Data

```

y <- x01 <- x02 <- z01 <- z02 <- c(1:300)
y[1:100] <- 1
y[101:200] <- 2
y[201:300] <- 3
x01[1:100] <- rnorm(100, 25, 2.5)
x01[101:200] <- rnorm(100, 40, 4.0)
x01[201:300] <- rnorm(100, 35, 3.5)
x02[1:100] <- rnorm(100, 2.51, 0.25)
x02[101:200] <- rnorm(100, 2.01, 0.20)
x02[201:300] <- rnorm(100, 2.70, 0.27)
z01[1:100] <- 1
z01[101:200] <- 2
z01[201:300] <- 3
z02[1:100] <- 40
z02[101:200] <- 50
z02[201:300] <- 100
N <- length(y)
J <- 3 #Number of categories in y
K <- 3 #Number of individual attributes (including the intercept)
C <- 2 #Number of choice-based attributes (intercept is not included)
X <- matrix(c(rep(1,N),x01,x02),N,K) #Design matrix of individual attrib.
Z <- matrix(c(z01,z02),N,C) #Design matrix of choice-based attributes
mon.names <- "LP"
parm.names <- parm.names(list(beta=matrix(0,J-1,K), gamma=rep(0,C)))
MyData <- list(C=C, J=J, K=K, N=N, X=X, Z=Z, mon.names=mon.names,
  parm.names=parm.names, y=y)

```

12.3. Initial Values

```
Initial.Values <- c(rep(0,(J-1)*K), rep(0,C))
```

12.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  beta <- matrix(parm[grep("beta", Data$parm.names)], Data$J-1, Data$K)
  gamma <- parm[grep("gamma", Data$parm.names)]
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))

```

```

gamma.prior <- sum(dnorm(gamma, 0, sqrt(1000), log=TRUE))
### Log-Likelihood
mu <- matrix(rep(tcrossprod(gamma, Data$Z),J),Data$N,Data$J)
mu[,1] <- mu[,1] + tcrossprod(beta[1,], Data$X)
mu[,2] <- mu[,2] + tcrossprod(beta[2,], Data$X)
mu <- interval(mu, -700, 700)
phi <- exp(mu)
p <- phi / rowSums(phi)
LL <- sum(dcat(Data$y, p, log=TRUE))
yrep <- apply(p,1,which.max)
### Log-Posterior
LP <- LL + beta.prior + gamma.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=LP, yhat=yrep, parm=parm)
return(Modelout)
}

```

13. Discrete Choice, Mixed Logit

13.1. Form

$$\begin{aligned}
\mathbf{y}_i &\sim \mathcal{CAT}(\mathbf{p}_{i,1:J}), \quad i = 1, \dots, N, \quad j = 1, \dots, J \\
\mathbf{p}_{i,j} &= \frac{\phi_{i,j}}{\sum_{j=1}^J \phi_{i,j}} \\
\phi &= \exp(\mu) \\
\mu_{i,j} &= \beta_{j,1:K} \mathbf{X}_{i,1:K} + \gamma \mathbf{Z}_{i,1:C} \in [-700, 700], \quad j = 1, \dots, (J-1) \\
\mu_{i,J} &= \gamma \mathbf{Z}_{i,1:C} \\
\beta_{j,k} &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, (J-1) \\
\gamma_c &\sim \mathcal{N}(\zeta_{\mu[c]}, \zeta_{\sigma[c]}^2) \\
\zeta_{\mu[c]} &\sim \mathcal{N}(0, 1000) \\
\zeta_{\sigma[c]} &\sim \mathcal{HC}(25)
\end{aligned}$$

13.2. Data

```

y <- x01 <- x02 <- z01 <- z02 <- c(1:300)
y[1:100] <- 1
y[101:200] <- 2
y[201:300] <- 3
x01[1:100] <- rnorm(100, 25, 2.5)
x01[101:200] <- rnorm(100, 40, 4.0)
x01[201:300] <- rnorm(100, 35, 3.5)

```

```

x02[1:100] <- rnorm(100, 2.51, 0.25)
x02[101:200] <- rnorm(100, 2.01, 0.20)
x02[201:300] <- rnorm(100, 2.70, 0.27)
z01[1:100] <- 1
z01[101:200] <- 2
z01[201:300] <- 3
z02[1:100] <- 40
z02[101:200] <- 50
z02[201:300] <- 100
N <- length(y)
J <- 3 #Number of categories in y
K <- 3 #Number of individual attributes (including the intercept)
C <- 2 #Number of choice-based attributes (intercept is not included)
X <- matrix(c(rep(1,N),x01,x02),N,K) #Design matrix of individual attrib.
Z <- matrix(c(z01,z02),N,C) #Design matrix of choice-based attributes
mon.names <- c("LP", parm.names(list(zeta.sigma=rep(0,C))))
parm.names <- parm.names(list(beta=matrix(0,J-1,K), gamma=rep(0,C),
  zeta.mu=rep(0,C), log.zeta.sigma=rep(0,C)))
MyData <- list(C=C, J=J, K=K, N=N, X=X, Z=Z, mon.names=mon.names,
  parm.names=parm.names, y=y)

```

13.3. Initial Values

```
Initial.Values <- c(rep(0,(J-1)*K), rep(0,N*C), rep(0,C), rep(0,C))
```

13.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  beta <- matrix(parm[grepl("beta", Data$parm.names)], Data$J-1, Data$K)
  gamma <- parm[grepl("gamma", Data$parm.names)]
  zeta.mu <- parm[grepl("zeta.mu", Data$parm.names)]
  zeta.sigma <- exp(parm[grepl("log.zeta.sigma", Data$parm.names)])
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  gamma.prior <- sum(dnorm(gamma, 0, sqrt(1000), log=TRUE))
  zeta.mu.prior <- sum(dnorm(zeta.mu, 0, sqrt(1000), log=TRUE))
  zeta.sigma.prior <- sum(dhalfcauchy(zeta.sigma, 25, log=TRUE))
  ### Log-Likelihood
  mu <- matrix(rep(rowSums(gamma * Data$Z),J),Data$N,Data$J)
  mu[,1] <- mu[,1] + tcrossprod(beta[1,], Data$X)
  mu[,2] <- mu[,2] + tcrossprod(beta[2,], Data$X)
  mu <- interval(mu, -700, 700)
  phi <- exp(mu)
  p <- phi / rowSums(phi)
  LL <- sum(dcat(Data$y, p, log=TRUE))
}

```

```

yrep <- apply(p,1,which.max)
### Log-Posterior
LP <- LL + beta.prior + gamma.prior + zeta.mu.prior + zeta.sigma.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,zeta.sigma.prior),
  yhat=yrep, parm=parm)
return(Modelout)
}

```

14. Discrete Choice, Multinomial Probit

14.1. Form

$$\begin{aligned}
\mathbf{Z}_{i,1:J} &\sim \mathcal{N}_J(\mu_{i,1:J}, \Sigma), \quad i = 1, \dots, N \\
\mathbf{Z}_{i,j} &\in \begin{cases} [0,10] & \text{if } \mathbf{y}_i = j \\ [-10,0] & \end{cases} \\
\mu_{1:N,j} &= \mathbf{X}\beta_{j,1:K} + \mathbf{W}\gamma[a, 1 : C] \\
\mathbf{a} &= \begin{cases} 1 & \text{if } \mathbf{y}_i < J \\ 2 & \end{cases} \\
\Sigma &\sim \mathcal{IW}(J, \mathbf{R}), \quad \mathbf{R} = \mathbf{I}_J, \quad \Sigma[1,1] = 1 \\
\beta_{j,k} &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, (J-1), \quad k = 1, \dots, K \\
\beta_{J,k} &= -\sum_{j=1}^{J-1} \beta_{j,k} \\
\gamma_{1,1:C} &\sim \mathcal{N}(0, 1000) \\
\gamma_{2,c} &= -\gamma_{1,c}, \quad c = 1, \dots, C \\
\mathbf{Z}_{i,j} &\sim \mathcal{N}(0, 1000) \in [-10, 10]
\end{aligned}$$

14.2. Data

```

y <- x1 <- x2 <- w1 <- w2 <- c(1:30)
y[1:10] <- 1
y[11:20] <- 2
y[21:30] <- 3
x1[1:10] <- rnorm(10, 25, 2.5)
x1[11:20] <- rnorm(10, 40, 4.0)
x1[21:30] <- rnorm(10, 35, 3.5)
x2[1:10] <- rnorm(10, 2.51, 0.25)
x2[11:20] <- rnorm(10, 2.01, 0.20)
x2[21:30] <- rnorm(10, 2.70, 0.27)
w1[1:10] <- 10

```



```

w1[11:20] <- 4
w1[21:30] <- 1
w2[1:10] <- 40
w2[11:20] <- 50
w2[21:30] <- 100
N <- length(y)
J <- length(unique(y)) #Number of categories in y
K <- 3 #Number of columns to be in design matrix X
R <- diag(J)
X <- matrix(c(rep(1,N),x1,x2),N,K)
C <- 2 #Number of choice-based attributes
W <- matrix(c(w1,w2),N,C) #Design matrix of choice-based attributes
mon.names <- "LP"
sigma.temp <- parm.names(list(Sigma=diag(J)), uppertri=1)
parm.names <- c(sigma.temp[2:length(sigma.temp)],
  parm.names(list(beta=matrix(0,(J-1),K), gamma=rep(0,C),
    Z=matrix(0,N,J))))
MyData <- list(J=J, K=K, N=N, R=R, W=W, X=X, mon.names=mon.names,
  parm.names=parm.names, y=y)

```

14.3. Initial Values

```

Initial.Values <- c(rep(0,length(R[upper.tri(R, diag=TRUE)]))-1),
  rep(0,(J-1)*K), rep(0,C), rep(0,N,J))

```

14.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  beta <- matrix(parm[grepl("beta", Data$parm.names)], Data$J-1, Data$K)
  beta <- rbind(beta, colSums(beta)*-1) #Sum to zero constraint
  gamma <- parm[grepl("gamma", Data$parm.names)]
  gamma <- rbind(gamma, gamma*-1) #Sum to zero constraint
  Sigma <- matrix(NA, Data$J, Data$J)
  Sigma[upper.tri(Sigma, diag=TRUE)] <- c(0, parm[grepl("Sigma",
    Data$parm.names)])
  Sigma[lower.tri(Sigma)] <- Sigma[upper.tri(Sigma)]
  diag(Sigma) <- exp(diag(Sigma))
  Z <- matrix(parm[grepl("Z", Data$parm.names)], Data$N, Data$J)
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  gamma.prior <- sum(dnorm(gamma, 0, sqrt(1000), log=TRUE))
  Sigma.prior <- dinvwishart(Sigma, Data$J, Data$R, log=TRUE)
  Z.prior <- sum(dnorm(Z, 0, sqrt(1000), log=TRUE))
  ### Log-Likelihood

```

```

mu <- matrix(0,Data$N,Data$J)
mu <- matrix(c(rep(tcrossprod(gamma[1,], Data$W),J),
  tcrossprod(gamma[2,], Data$W)),Data$N,Data$J)
for (j in 1:Data$J) {mu[,j] <- mu[,j] + tcrossprod(beta[j,], Data$X)}
Y <- indmat(Data$y)
Z <- ifelse(Z > 10, 10, Z); Z <- ifelse({Y == 0} & {Z > 0}, 0, Z)
Z <- ifelse(Z < -10, -10, Z); Z <- ifelse({Y == 1} & {Z < 0}, 0, Z)
parm[grep("Z", Data$parm.names)] <- as.vector(Z)
LL <- sum(dmvn(Z, mu, Sigma, log=TRUE))
yrep <- apply(Z, 1, which.max)
#eta <- exp(mu)
#p <- eta / rowSums(eta)
### Log-Posterior
LP <- LL + beta.prior + gamma.prior + Sigma.prior + Z.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=LP, yhat=yrep, parm=parm)
return(Modelout)
}

```

15. Dynamic Linear Model (DLM)

The data is presented so that the time-series is subdivided into three sections: modeled ($t = 1, \dots, T_m$), one-step ahead forecast ($t = T_m + 1$), and future forecast [$t = (T_m + 2), \dots, T$].

15.1. Form

$$\begin{aligned}
\mathbf{y}_t &\sim \mathcal{N}(\mu_t, \sigma_V^2), \quad t = 1, \dots, T_m \\
\mathbf{y}_t^{new} &\sim \mathcal{N}(\mu_t, \sigma_V^2), \quad t = (T_m + 1), \dots, T \\
\mu_t &= \alpha + \mathbf{x}_t \beta_t, \quad t = 1, \dots, T \\
\alpha &\sim \mathcal{N}(0, 1000) \\
\beta_1 &\sim \mathcal{N}(0, 1000) \\
\beta_t &\sim \mathcal{N}(\beta_{t-1}, \sigma_W^2), \quad t = 2, \dots, T \\
\sigma_V &\sim \mathcal{HC}(25) \\
\sigma_W &\sim \mathcal{HC}(25)
\end{aligned}$$

15.2. Data

```

T <- 20
T.m <- 14
beta.orig <- x <- rep(0,T)
for (t in 2:T) {
beta.orig[t] <- beta.orig[t-1] + rnorm(1,0,0.1)

```

```

x[t] <- x[t-1] + rnorm(1,0,0.1)}
y <- 10 + beta.orig*x + rnorm(T,0,0.1)
y[(T.m+2):T] <- NA
mon.names <- rep(NA, (T-T.m))
for (i in 1:(T-T.m)) mon.names[i] <- paste("mu[",(T.m+i),"]", sep="")
parm.names <- parm.names(list(alpha=0, beta=rep(0,T), log.beta.w.sigma=0,
  log.v.sigma=0))
MyData <- list(T=T, T.m=T.m, mon.names=mon.names, parm.names=parm.names,
  x=x, y=y)

```

15.3. Initial Values

```
Initial.Values <- rep(0,T+3)
```

15.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  alpha <- parm[1]
  beta <- parm[2:(Data$T+1)]
  beta.w.sigma <- exp(parm[Data$T+2])
  v.sigma <- exp(parm[Data$T+3])
  ### Log(Prior Densities)
  alpha.prior <- dnorm(alpha, 0, sqrt(1000), log=TRUE)
  beta.prior <- rep(0,Data$T)
  beta.prior[1] <- dnorm(beta[1], 0, sqrt(1000), log=TRUE)
  beta.prior[2:Data$T] <- dnorm(beta[2:Data$T], beta[1:(Data$T-1)],
    beta.w.sigma, log=TRUE)
  beta.w.sigma.prior <- dhalfcauchy(beta.w.sigma, 25, log=TRUE)
  v.sigma.prior <- dhalfcauchy(v.sigma, 25, log=TRUE)
  ### Log-Likelihood
  mu <- alpha + beta*Data$x
  LL <- sum(dnorm(Data$y[1:Data$T.m], mu[1:Data$T.m], v.sigma,
    log=TRUE))
  ### Log-Posterior
  LP <- LL + alpha.prior + sum(beta.prior) + beta.w.sigma.prior +
    v.sigma.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=mu[(Data$T.m+1):Data$T],
    yhat=mu, parm=parm)
  return(Modelout)
}

```

16. Factor Analysis, Confirmatory

Factor scores are in matrix \mathbf{F} , factor loadings for each variable are in vector λ , and \mathbf{f} is a

vector that indicates which variable loads on which factor.

16.1. Form

$$\begin{aligned}
\mathbf{Y}_{i,m} &\sim \mathcal{N}(\mu_{i,m}, \sigma_m^2), \quad i = 1, \dots, N, \quad m = 1, \dots, M \\
\mu_{i,m} &= \alpha_m + \lambda_m \mathbf{F}_{i, \mathbf{f}[m]}, \quad i = 1, \dots, N, \quad m = 1, \dots, M \\
\mathbf{F}_{i, 1:P} &\sim \mathcal{N}_P(\boldsymbol{\gamma}, \boldsymbol{\Omega}^{-1}), \quad i = 1, \dots, N \\
\alpha_m &\sim \mathcal{N}(0, 1000), \quad m = 1, \dots, M \\
\lambda_m &\sim \mathcal{N}(0, 1000), \quad m = 1, \dots, M \\
\sigma_m &\sim \mathcal{HC}(25), \quad m = 1, \dots, M \\
\boldsymbol{\Omega} &\sim \mathcal{W}(N, \mathbf{S}), \quad \mathbf{S} = \mathbf{I}_P
\end{aligned}$$

16.2. Data

```

data(swiss)
Y <- cbind(swiss$Agriculture, swiss$Examination, swiss$Education,
           swiss$Catholic, swiss$Infant.Mortality)
M <- NCOL(Y) #Number of variables
N <- NROW(Y) #Number of records
P <- 3 #Number of factors
f <- c(1,3,2,2,1) #Indicator f for the factor for each variable m
gamma <- rep(0,P)
S <- diag(P)
mon.names <- c("LP", "mu[1,1]")
parm.names <- parm.names(list(F=matrix(0,N,P), lambda=rep(0,M),
                                Omega=diag(P), alpha=rep(0,M), log.sigma=rep(0,M)),
                          uppertri=c(0,0,1,0,0))
MyData <- list(M=M, N=N, P=P, S=S, Y=Y, f=f, gamma=gamma,
               mon.names=mon.names, parm.names=parm.names)

```

16.3. Initial Values

```

Initial.Values <- c(rep(0,N*P), rep(0,M), S[upper.tri(S, diag=TRUE)],
                    rep(0,M), rep(0,M))

```

16.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  alpha <- parm[grep("alpha", Data$parm.names)]
  lambda <- parm[grep("lambda", Data$parm.names)]
  sigma <- exp(parm[grep("log.sigma", Data$parm.names)])
}

```

```

F <- matrix(parm[grepl("F", Data$parm.names)], Data$N, Data$P)
Omega <- matrix(NA, Data$P, Data$P)
Omega[upper.tri(Omega, diag=TRUE)] <- parm[grepl("Omega",
  Data$parm.names)]
Omega[lower.tri(Omega)] <- Omega[upper.tri(Omega)]
Sigma <- solve(Omega)
### Log(Prior Densities)
alpha.prior <- sum(dnorm(alpha, 0, sqrt(1000), log=TRUE))
lambda.prior <- sum(dnorm(lambda, 0, sqrt(1000), log=TRUE))
sigma.prior <- sum(dhalfcauchy(sigma, 25, log=TRUE))
Omega.prior <- dwishart(Omega, Data$N, Data$S, log=TRUE)
F.prior <- sum(dmvn(F, Data$gamma, Sigma, log=TRUE))
### Log-Likelihood
mu <- Data$Y
for (m in 1:Data$M) {mu[,m] <- alpha[m] + lambda[m] * F[,Data$f[m]]}
LL <- sum(dnorm(Data$Y, mu, sigma, log=TRUE))
### Log-Posterior
LP <- LL + alpha.prior + lambda.prior + sigma.prior + F.prior +
  Omega.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,mu[1,1]),
  yhat=mu, parm=parm)
return(Modelout)
}

```

17. Factor Analysis, Exploratory

Factor scores are in matrix \mathbf{F} and factor loadings are in matrix Λ . Although the calculation for the recommended number of factors to explore P is also provided below (Fokoue 2004), this example sets $P = 3$.

17.1. Form

$$\begin{aligned}
\mathbf{Y}_{i,m} &\sim \mathcal{N}(\mu_{i,m}, \sigma_m^2), \quad i = 1, \dots, N, \quad m = 1, \dots, M \\
\mu_{i,m} &= \alpha_m + \sum_{p=1}^P \nu_{i,m,p}, \quad i = 1, \dots, N, \quad m = 1, \dots, M \\
\nu_{i,m,p} &= \mathbf{F}_{i,p} \Lambda_{p,m}, \quad i = 1, \dots, N, \quad m = 1, \dots, M, \quad p = 1, \dots, P \\
\mathbf{F}_{i,1:P} &\sim \mathcal{N}_P(\gamma, \Omega^{-1}), \quad i = 1, \dots, N \\
\alpha_m &\sim \mathcal{N}(0, 1000), \quad m = 1, \dots, M \\
\gamma_p &= 0, \quad p = 1, \dots, P \\
\Lambda_{p,m} &\sim \mathcal{N}(0, 1000), \quad p = 1, \dots, P, \quad m = 1, \dots, M \\
\Omega &\sim \mathcal{W}(N, \mathbf{S}), \quad \mathbf{S} = \mathbf{I}_P
\end{aligned}$$

$$\sigma_m \sim \mathcal{HC}(25), \quad m = 1, \dots, M$$

17.2. Data

```
data(swiss)
Y <- cbind(swiss$Agriculture, swiss$Examination, swiss$Education,
swiss$Catholic, swiss$Infant.Mortality)
M <- NCOL(Y) #Number of variables
N <- NROW(Y) #Number of records
P <- trunc(0.5*(2*M + 1 - sqrt(8*M + 1))) #Number of factors to explore
P <- 3 #Number of factors to explore (override for this example)
gamma <- rep(0,P)
S <- diag(P)
mon.names <- c("LP","mu[1,1]")
parm.names <- parm.names(list(F=matrix(0,N,P), Lambda=matrix(0,P,M),
  Omega=diag(P), alpha=rep(0,M), log.sigma=rep(0,M)),
  uppertri=c(0,0,1,0,0))
MyData <- list(M=M, N=N, P=P, S=S, Y=Y, gamma=gamma, mon.names=mon.names,
  parm.names=parm.names)
```

17.3. Initial Values

```
Initial.Values <- c(rep(0,N*P), rep(0,P*M), S[upper.tri(S, diag=TRUE)],
  rep(0,M), rep(0,M))
```

17.4. Model

```
Model <- function(parm, Data)
{
  ### Parameters
  alpha <- parm[grep("alpha", Data$parm.names)]
  sigma <- exp(parm[grep("log.sigma", Data$parm.names)])
  F <- matrix(parm[grep("F", Data$parm.names)], Data$N, Data$P)
  Lambda <- matrix(parm[grep("Lambda", Data$parm.names)],
    Data$P, Data$M)
  Omega <- matrix(NA, Data$P, Data$P)
  Omega[upper.tri(Omega, diag=TRUE)] <- parm[grep("Omega",
    Data$parm.names)]
  Omega[lower.tri(Omega)] <- Omega[upper.tri(Omega)]
  Sigma <- solve(Omega)
  ### Log(Prior Densities)
  alpha.prior <- sum(dnorm(alpha, 0, sqrt(1000), log=TRUE))
  sigma.prior <- sum(dhalfcauchy(sigma, 25, log=TRUE))
  Omega.prior <- dwishart(Omega, Data$N, Data$S, log=TRUE)
  F.prior <- sum(dmvn(F, Data$gamma, Sigma, log=TRUE))
  Lambda.prior <- sum(dnorm(Lambda, 0, sqrt(1000), log=TRUE))
  ### Log-Likelihood
```

```

mu <- Data$Y
nu <- array(NA, dim=c(Data$N, Data$M, Data$P))
for (p in 1:Data$P) {nu[, ,p] <- F[,p, drop=FALSE] %*% Lambda[p,]}
for (m in 1:Data$M) {mu[,m] <- alpha[m] + rowSums(nu[,1,])}
LL <- sum(dnorm(Data$Y, mu, sigma, log=TRUE))
### Log-Posterior
LP <- LL + alpha.prior + sigma.prior + Omega.prior + F.prior +
      Lambda.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,mu[1,1]),
      yhat=mu, parm=parm)
return(Modelout)
}

```

18. Kriging

This is an example of universal kriging of \mathbf{y} given \mathbf{X} , regression effects β , and spatial effects ζ . Euclidean distance between spatial coordinates (longitude and latitude) is used for each of $i = 1, \dots, N$ records of \mathbf{y} . An additional record is created from the same data-generating process to compare the accuracy of interpolation. For the spatial component, ϕ is the rate of spatial decay and κ is the scale. κ is often difficult to identify, so it is set to 1 (Gaussian), but may be allowed to vary up to 2 (Exponential). In practice, ϕ is also often difficult to identify. While Σ is spatial covariance, spatial correlation is $\rho = \exp(-\phi\mathbf{D})$.

18.1. Form

$$\mathbf{y} \sim \mathcal{N}(\mu, \sigma_1^2)$$

$$\mu = \mathbf{X}\beta + \zeta$$

$$\mathbf{y}^{new} = \mathbf{X}\beta + \sum_{i=1}^N \left(\frac{\rho_i}{\sum \rho} \zeta_i \right)$$

$$\rho = \exp(-\phi\mathbf{D}^{new})^\kappa$$

$$\zeta \sim \mathcal{N}_N(\zeta_\mu, \Sigma)$$

$$\Sigma = \sigma_2^2 \exp(-\phi\mathbf{D})^\kappa$$

$$\beta_j \sim \mathcal{N}(0, 1000), \quad j = 1, \dots, 2$$

$$\sigma_j \sim \mathcal{HC}(25), \quad j = 1, \dots, 2$$

$$\phi \sim \mathcal{U}(1, 5)$$

$$\zeta_\mu = 0$$

$$\kappa = 1$$

18.2. Data

```

N <- 20
longitude <- runif(N+1,0,100)
latitude <- runif(N+1,0,100)
D <- as.matrix(dist(cbind(longitude,latitude), diag=TRUE, upper=TRUE))
Sigma <- 10000 * exp(-1.5 * D)
zeta <- as.vector(apply(rmvn(1000, rep(0,N+1), Sigma), 2, mean))
beta <- c(50,2)
X <- matrix(runif((N+1)*2,-2,2),(N+1),2); X[,1] <- 1
mu <- as.vector(tcrossprod(beta, X))
y <- mu + zeta
longitude.new <- longitude[N+1]; latitude.new <- latitude[N+1]
Xnew <- X[N+1,]; ynew <- y[N+1]
longitude <- longitude[1:N]; latitude <- latitude[1:N]
X <- X[1:N,]; y <- y[1:N]
D <- as.matrix(dist(cbind(longitude,latitude), diag=TRUE, upper=TRUE))
D.new <- sqrt((longitude - longitude.new)^2 + (latitude - latitude.new)^2)
mon.names <- c("LP","sigma[1]","sigma[2]","ynew")
parm.names <- parm.names(list(zeta=rep(0,N), beta=rep(0,2),
  log.sigma=rep(0,2), phi=0))
MyData <- list(D=D, D.new=D.new, N=N, X=X, Xnew=Xnew, mon.names=mon.names,
  parm.names=parm.names, y=y)

```

18.3. Initial Values

```
Initial.Values <- c(rep(0,N), rep(0,2), rep(0,2), 1)
```

18.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  beta <- parm[grep("beta", Data$parm.names)]
  zeta <- parm[grep("zeta", Data$parm.names)]
  kappa <- 1
  sigma <- exp(parm[grep("log.sigma", Data$parm.names)])
  phi <- interval(parm[grep("phi", Data$parm.names)], 1, 5)
  parm[grep("phi", Data$parm.names)] <- phi
  Sigma <- sigma[2]*sigma[2] * exp(-phi * Data$D)^kappa
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  zeta.prior <- dmvn(zeta, rep(0, Data$N), Sigma, log=TRUE)
  sigma.prior <- sum(dhalfcauchy(sigma, 25, log=TRUE))
  phi.prior <- dunif(phi, 1, 5, log=TRUE)
  ### Interpolation
  rho <- exp(-phi * Data$D.new)^kappa
  ynew <- sum(beta * Data$Xnew) + sum(rho / sum(rho) * zeta)
}

```



```

### Log-Likelihood
mu <- tcrossprod(beta, Data$X) + zeta
LL <- sum(dnorm(Data$y, mu, sigma[1], log=TRUE))
### Log-Posterior
LP <- LL + beta.prior + zeta.prior + sigma.prior + phi.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,sigma,ynew),
  yhat=mu, parm=parm)
return(Modelout)
}

```

19. Laplace Regression

This linear regression specifies that \mathbf{y} is Laplace-distributed, where it is usually Gaussian or normally-distributed. It has been claimed that it should be surprising that the normal distribution became the standard, when the Laplace distribution usually fits better and has wider tails (Kotz, Kozubowski, and Podgorski 2001). Another popular alternative is to use the t-distribution (see Robust Regression in section 31), though it is more computationally expensive to estimate, because it has three parameters. The Laplace distribution has only two parameters, location and scale like the normal distribution, and is computationally easier to fit. This example could be taken one step further, and the parameter vector β could be Laplace-distributed. Laplace's Demon recommends that users experiment with replacing the normal distribution with the Laplace distribution.

19.1. Form

$$\mathbf{y} \sim \mathcal{L}(\mu, \sigma^2)$$

$$\mu = \mathbf{X}\beta$$

$$\beta_j \sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J$$

$$\sigma \sim \mathcal{HC}(25)$$

19.2. Data

```

N <- 10000
J <- 5
X <- matrix(1,N,J)
for (j in 2:J) {X[,j] <- rnorm(N,runif(1,-3,3),runif(1,0.1,1))}
beta <- runif(J,-3,3)
e <- rlaplace(N,0,0.1)
y <- as.vector(tcrossprod(beta, X) + e)
mon.names <- c("LP", "sigma")
parm.names <- parm.names(list(beta=rep(0,J), log.sigma=0))
MyData <- list(J=J, X=X, mon.names=mon.names, parm.names=parm.names, y=y)

```

19.3. Initial Values

```
Initial.Values <- c(rep(0,J), log(1))
```

19.4. Model

```
Model <- function(parm, Data)
{
  ### Parameters
  beta <- parm[1:Data$J]
  sigma <- exp(parm[Data$J+1])
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  sigma.prior <- dhalfcauchy(sigma, 25, log=TRUE)
  ### Log-Likelihood
  mu <- tcrossprod(beta, Data$X)
  LL <- sum(dlaplace(Data$y, mu, sigma, log=TRUE))
  ### Log-Posterior
  LP <- LL + beta.prior + sigma.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP, sigma), yhat=mu,
    parm=parm)
  return(Modelout)
}
```

20. Linear Regression

20.1. Form

$$\begin{aligned} \mathbf{y} &\sim \mathcal{N}(\boldsymbol{\mu}, \sigma^2) \\ \boldsymbol{\mu} &= \mathbf{X}\boldsymbol{\beta} \\ \beta_j &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J \\ \sigma &\sim \mathcal{HC}(25) \end{aligned}$$

20.2. Data

```
N <- 10000
J <- 5
X <- matrix(1,N,J)
for (j in 2:J) {X[,j] <- rnorm(N,runif(1,-3,3),runif(1,0.1,1))}
beta <- runif(J,-3,3)
e <- rnorm(N,0,0.1)
y <- as.vector(tcrossprod(beta, X) + e)
mon.names <- c("LP", "sigma")
parm.names <- parm.names(list(beta=rep(0,J), log.sigma=0))
```

```
MyData <- list(J=J, X=X, mon.names=mon.names, parm.names=parm.names, y=y)
```

20.3. Initial Values

```
Initial.Values <- c(rep(0,J), log(1))
```

20.4. Model

```
Model <- function(parm, Data)
{
  ### Parameters
  beta <- parm[1:Data$J]
  sigma <- exp(parm[Data$J+1])
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  sigma.prior <- dgamma(sigma, 25, log=TRUE)
  ### Log-Likelihood
  mu <- tcrossprod(beta, Data$X)
  LL <- sum(dnorm(Data$y, mu, sigma, log=TRUE))
  ### Log-Posterior
  LP <- LL + beta.prior + sigma.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP, sigma), yhat=mu,
    parm=parm)
  return(Modelout)
}
```

21. Linear Regression, Multilevel

21.1. Form

$$\mathbf{y} \sim \mathcal{N}(\mu, \sigma^2)$$

$$\mu_i = \mathbf{X} \beta_{\mathbf{m}[i], 1:J}$$

$$\beta_{g, 1:J} \sim \mathcal{N}_J(\gamma, \Sigma), \quad g = 1, \dots, G$$

$$\Sigma = \Omega^{-1}$$

$$\Omega \sim \mathcal{W}(J, \mathbf{S}), \quad \mathbf{S} = \mathbf{I}_J$$

$$\gamma_j \sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J$$

$$\sigma \sim \mathcal{HC}(25)$$

where \mathbf{m} is a vector of length N , and each element indicates the multilevel group ($g = 1, \dots, G$) for the associated record.

21.2. Data

```

N <- 30
J <- 2 ### Number of predictors (including intercept)
G <- 2 ### Number of Multilevel Groups
X <- matrix(rnorm(N,0,1),N,J); X[,1] <- 1
Sigma <- matrix(runif(J*J,-1,1),J,J)
diag(Sigma) <- runif(J,1,5)
gamma <- runif(J,-1,1)
beta <- matrix(NA,G,J)
for (g in 1:G) {beta[g,] <- rmvn(1, gamma, Sigma)}
m <- round(runif(N,0.5,(G+0.49))) ### Multilevel group indicator
y <- rowSums(beta[m,] * X) + rnorm(N,0,0.1)
S <- diag(J)
mon.names <- c("LP","sigma")
parm.names <- parm.names(list(beta=matrix(0,G,J), log.sigma=0,
  gamma=rep(0,J), Omega=S), uppertri=c(0,0,0,1))
MyData <- list(G=G, J=J, N=N, S=S, X=X, m=m, mon.names=mon.names,
  parm.names=parm.names, y=y)

```

21.3. Initial.Values

```

Initial.Values <- c(rep(0,G*J), log(1), rep(0,J),
  S[upper.tri(S, diag=TRUE)])

```

21.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  beta <- matrix(parm[1:(Data$G * Data$J)], Data$G, Data$J)
  gamma <- parm[grep("gamma", Data$parm.names)]
  sigma <- exp(parm[grep("log.sigma", Data$parm.names)])
  Omega <- matrix(NA, Data$J, Data$J)
  Omega[upper.tri(Omega, diag=TRUE)] <- parm[min(grep("Omega",
    Data$parm.names)): max(grep("Omega", Data$parm.names))]
  Omega[lower.tri(Omega)] <- Omega[upper.tri(Omega)]
  Sigma <- solve(Omega)
  ### Log(Prior Densities)
  Omega.prior <- dwishart(Omega, Data$J, Data$S, log=TRUE)
  beta.prior <- sum(dmvn(beta, gamma, Sigma, log=TRUE))
  gamma.prior <- sum(dnorm(gamma, 0, sqrt(100), log=TRUE))
  sigma.prior <- dhalfcauchy(sigma, 25, log=TRUE)
  ### Log-Likelihood
  mu <- rowSums(beta[Data$m,] * Data$X)
  LL <- sum(dnorm(Data$y, mu, sigma, log=TRUE))
  ### Log-Posterior

```

```

LP <- LL + Omega.prior + beta.prior + gamma.prior + sigma.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,sigma),
  yhat=mu, parm=parm)
return(Modelout)
}

```

22. Linear Regression with Full Missingness

With ‘full missingness’, there are missing values for both the response and at least one predictor. This is a minimal example, since there are missing values in only one of the predictors. Initial values do not need to be specified for missing values in a predictor, unless another predictor variable with missing values is used to predict the missing values of a predictor. More effort is involved in specifying a model with a missing predictor that is predicted by another missing predictor. The full likelihood approach to full missingness is excellent as long as the model is identifiable. When it is not identifiable, then imputation may be done in a previous stage. In this example, $X[,2]$ is the only predictor with missing values.

22.1. Form

$$\begin{aligned}
\mathbf{y} &\sim \mathcal{N}(\mu_2, \sigma_2^2) \\
\mu_2 &= \mathbf{X}\beta \\
\mathbf{X}_{1:N,2} &\sim \mathcal{N}(\mu_1, \sigma_1^2) \\
\mu_1 &= \mathbf{X}_{1:N,(1,3:J)}\alpha \\
\alpha_j &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, (J-1) \\
\beta_j &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J \\
\sigma_k &\sim \mathcal{HC}(25), \quad k = 1, \dots, 2
\end{aligned}$$

22.2. Data

```

N <- 1000
J <- 5
X <- matrix(runif(N*J,-2,2),N,J)
X[,1] <- 1
alpha <- runif((J-1),-2,2)
X[,2] <- tcrossprod(alpha, X[, -2]) + rnorm(N,0,0.1)
beta <- runif(J,-2,2)
y <- as.vector(tcrossprod(beta, X) + rnorm(N,0,0.1))
y[sample(1:N, round(N*0.05))] <- NA
M <- ifelse(is.na(y), 1, 0)
X[sample(1:N, round(N*0.05)),2] <- NA
mon.names <- c("LP","sigma[1]","sigma[2]")

```

```

parm.names <- parm.names(list(alpha=rep(0,J-1), beta=rep(0,J),
  log.sigma=rep(0,2)))
MyData <- list(J=J, M=M, N=N, X=X, mon.names=mon.names, parm.names=parm.names,
  y=y)

```

22.3. Initial Values

```
Initial.Values <- c(rep(0,(J-1)), rep(0,J), rep(0,2))
```

22.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  alpha <- parm[1:(Data$J-1)]
  beta <- parm[Data$J:(2*Data$J - 1)]
  sigma <- exp(parm[(2*Data$J):(2*Data$J+1)])
  ### Log(Prior Densities)
  alpha.prior <- sum(dnorm(alpha, 0, sqrt(1000), log=TRUE))
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  sigma.prior <- sum(dhalfcauchy(sigma, 25, log=TRUE))
  ### Log-Likelihood
  mu1 <- tcrossprod(alpha, Data$X[,-2])
  X.imputed <- Data$X
  X.imputed[,2] <- ifelse(is.na(Data$X[,2]), mu1, Data$X[,2])
  LL1 <- sum(dnorm(X.imputed[,2], mu1, sigma[1], log=TRUE))
  mu2 <- tcrossprod(beta, X.imputed)
  y.imputed <- ifelse(is.na(Data$y), mu2, Data$y)
  LL2 <- sum((1-Data$M) * dnorm(y.imputed, mu2, sigma[2], log=TRUE))
  ### Log-Posterior
  LP <- LL1 + LL2 + alpha.prior + beta.prior + sigma.prior
  Modelout <- list(LP=LP, Dev=-2*LL2, Monitor=c(LP,sigma),
    yhat=mu2, parm=parm)
  return(Modelout)
}

```

23. Linear Regression with Missing Response

Initial values do not need to be specified for missing values in this response, \mathbf{y} . Instead, at each iteration, missing values in \mathbf{y} are replaced with their estimate in μ .

23.1. Form

$$\mathbf{y} \sim \mathcal{N}(\mu, \sigma^2)$$

$$\mu = \mathbf{X}\beta$$

$$\beta_j \sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J$$

$$\sigma \sim \mathcal{HC}(25)$$

23.2. Data

```
data(demonsnacks)
N <- NROW(demonsnacks)
J <- NCOL(demonsnacks)
y <- log(demonsnacks$Calories)
y[sample(1:N, round(N*0.05))] <- NA
M <- ifelse(is.na(y), 1, 0)
X <- cbind(1, as.matrix(demonsnacks[,c(1,3:10)]))
for (j in 2:J) {X[,j] <- CenterScale(X[,j])}
mon.names <- c("LP", "sigma")
parm.names <- parm.names(list(beta=rep(0,J), log.sigma=0))
MyData <- list(J=J, M=M, X=X, mon.names=mon.names, parm.names=parm.names, y=y)
```

23.3. Initial Values

```
Initial.Values <- c(rep(0,J), log(1))
```

23.4. Model

```
Model <- function(parm, Data)
{
  ### Parameters
  beta <- parm[1:Data$J]
  sigma <- exp(parm[Data$J+1])
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  sigma.prior <- dgamma(sigma, 25, log=TRUE)
  ### Log-Likelihood
  mu <- tcrossprod(beta, Data$X)
  y.imputed <- ifelse(is.na(Data$y), mu, Data$y)
  LL <- sum((1-Data$M) * dnorm(y.imputed, mu, sigma, log=TRUE))
  ### Log-Posterior
  LP <- LL + beta.prior + sigma.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,sigma),
    yhat=mu, parm=parm)
  return(Modelout)
}
```

24. Multinomial Logit

24.1. Form

$$\mathbf{y}_i \sim \mathcal{CAT}(\mathbf{p}_{i,1:J}), \quad i = 1, \dots, N$$

$$\mathbf{p}_{i,j} = \frac{\phi_{i,j}}{\sum_{j=1}^J \phi_{i,j}}, \quad \sum_{j=1}^J \mathbf{p}_{i,j} = 1$$

$$\phi = \exp(\mu)$$

$$\mu_{i,J} = 0, \quad i = 1, \dots, N$$

$$\mu_{i,j} = \mathbf{X}_{i,1:K} \beta_{j,1:K} \in [-700, 700], \quad j = 1, \dots, (J-1)$$

$$\beta_{j,k} \sim \mathcal{N}(0, 1000), \quad j = 1, \dots, (J-1), \quad k = 1, \dots, K$$

24.2. Data

```

y <- x01 <- x02 <- c(1:300)
y[1:100] <- 1
y[101:200] <- 2
y[201:300] <- 3
x01[1:100] <- rnorm(100, 25, 2.5)
x01[101:200] <- rnorm(100, 40, 4.0)
x01[201:300] <- rnorm(100, 35, 3.5)
x02[1:100] <- rnorm(100, 2.51, 0.25)
x02[101:200] <- rnorm(100, 2.01, 0.20)
x02[201:300] <- rnorm(100, 2.70, 0.27)
N <- length(y)
J <- 3 #Number of categories in y
K <- 3 #Number of predictors (including the intercept)
X <- matrix(c(rep(1,N),x01,x02),N,K)
mon.names <- "LP"
parm.names <- c("beta[1,1]","beta[1,2]","beta[1,3]","beta[2,1]","
  "beta[2,2]","beta[2,3]") ### Parameter Names [J,K]
MyData <- list(J=J, K=K, N=N, X=X, mon.names=mon.names,
  parm.names=parm.names, y=y)

```

24.3. Initial Values

```
Initial.Values <- c(rep(0,(J-1)*K))
```

24.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  beta <- parm[1:(Data$J-1*Data$K)]

```



```

### Log(Prior Densities)
beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
### Log-Likelihood
mu <- matrix(0,Data$N,Data$J)
mu[,1] <- tcrossprod(beta[1:3], Data$X)
mu[,2] <- tcrossprod(beta[4:6], Data$X)
mu <- interval(mu, -700, 700)
phi <- exp(mu)
p <- phi / rowSums(phi)
LL <- sum(dcat(Data$y, p, log=TRUE))
yrep <- apply(p,1,which.max)
### Log-Posterior
LP <- LL + beta.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=LP, yhat=yrep, parm=parm)
return(Modelout)
}

```

25. Multinomial Logit, Nested

25.1. Form

$$\mathbf{y}_i \sim \mathcal{CAT}(\mathbf{P}_{i,1:J}), \quad i = 1, \dots, N$$

$$\mathbf{P}_{1:N,1} = \frac{\mathbf{R}}{\mathbf{R} + \exp(\alpha \mathbf{I})}$$

$$\mathbf{P}_{1:N,2} = \frac{(1 - \mathbf{P}_{1:N,1})\mathbf{S}_{1:N,1}}{\mathbf{V}}$$

$$\mathbf{P}_{1:N,3} = \frac{(1 - \mathbf{P}_{1:N,1})\mathbf{S}_{1:N,2}}{\mathbf{V}}$$

$$\mathbf{R}_{1:N} = \exp(\mu_{1:N,1})$$

$$\mathbf{S}_{1:N,1:2} = \exp(\mu_{1:N,2:3})$$

$$\mathbf{I} = \log(\mathbf{V})$$

$$\mathbf{V}_i = \sum_{k=1}^K \mathbf{S}_{i,k}, \quad i = 1, \dots, N$$

$$\mu_{1:N,1} = \mathbf{X}\boldsymbol{\iota} \in [-700, 700]$$

$$\mu_{1:N,2} = \mathbf{X}\boldsymbol{\beta}_{2,1:K} \in [-700, 700]$$

$$\boldsymbol{\iota} = \alpha \boldsymbol{\beta}_{1,1:K}$$

$$\alpha \sim \mathcal{E}\mathcal{X}\mathcal{P}(1) \in [0, 2]$$

$$\beta_{j,k} \sim \mathcal{N}(0, 1000), \quad j = 1, \dots, (J-1) \quad k = 1, \dots, K$$

where there are $J = 3$ categories of \mathbf{y} , $K = 3$ predictors, \mathbf{R} is the non-nested alternative, \mathbf{S} is the nested alternative, \mathbf{V} is the observed utility in the nest, α is effectively 1 - correlation and has a truncated exponential distribution, and ι is a vector of regression effects for the isolated alternative after α is taken into account. The third alternative is the reference category.

25.2. Data

```
y <- x01 <- x02 <- c(1:300)
y[1:100] <- 1
y[101:200] <- 2
y[201:300] <- 3
x01[1:100] <- rnorm(100, 25, 2.5)
x01[101:200] <- rnorm(100, 40, 4.0)
x01[201:300] <- rnorm(100, 35, 3.5)
x02[1:100] <- rnorm(100, 2.51, 0.25)
x02[101:200] <- rnorm(100, 2.01, 0.20)
x02[201:300] <- rnorm(100, 2.70, 0.27)
N <- length(y)
J <- 3 #Number of categories in y
K <- 3 #Number of predictors (including the intercept)
X <- matrix(c(rep(1,N),x01,x02),N,K)
mon.names <- c("LP",parm.names(list(iota=rep(0,K))))
parm.names <- parm.names(list(alpha=0, beta=matrix(0,J-1,K)))
MyData <- list(J=J, K=K, N=N, X=X, mon.names=mon.names,
               parm.names=parm.names, y=y)
```

25.3. Initial Values

```
Initial.Values <- c(0.5, rep(0.1,(J-1)*K))
```

25.4. Model

```
Model <- function(parm, Data)
{
  ### Hyperparameters
  alpha.rate <- 1
  ### Parameters
  alpha <- interval(parm[1],0,2); parm[1] <- alpha
  beta <- matrix(parm[grep("beta", Data$parm.names)], Data$J-1, Data$K)
  ### Log(Prior Densities)
  alpha.prior <- dtrunc(alpha, "exp", a=0, b=2, rate=alpha.rate,
                        log=TRUE)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  ### Log-Likelihood
  mu <- P <- matrix(0,Data$N,Data$J)
  iota <- alpha * beta[1,]
  mu[,1] <- tcrossprod(iota, Data$X)
```

```

mu[,2] <- tcrossprod(beta[2,], Data$X)
mu <- interval(mu, -700, 700)
R <- exp(mu[,1])
S <- exp(mu[,2:3])
V <- rowSums(S)
I <- log(V)
P[,1] <- R / (R + exp(alpha*I))
P[,2] <- (1 - P[,1]) * S[,1] / V
P[,3] <- (1 - P[,1]) * S[,2] / V
LL <- sum(dcat(Data$y, P, log=TRUE))
yrep <- apply(P,1,which.max)
### Log-Posterior
LP <- LL + alpha.prior + beta.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,iota), yhat=yrep,
  parm=parm)
return(Modelout)
}

```

26. Multinomial Probit

In this form of MNP, the β parameters are sum-to-zero constraints in the reference category, and covariance matrix Σ includes all J categories of \mathbf{y} .

Note that the parameters and initial values for the upper triangular elements of Σ are read in as Σ , though the diagonal is read in as $\log(\Sigma)$, but still denoted as Σ . Apologies for any confusion this causes, and the diagonal elements could each be renamed manually in `parm.names`. The only reason this difference exists is that I am unsure of how to program that in `parm.names` for all occasions.

26.1. Form

$$\mathbf{Z}_{i,1:J} \sim \mathcal{N}_J(\mu_{i,1:J}, \Sigma), \quad i = 1, \dots, N$$

$$\mathbf{Z}_{i,j} \in \begin{cases} [0,10] & \text{if } \mathbf{y}_i = j \\ [-10,0] & \end{cases}$$

$$\mu_{1:N,j} = \mathbf{X}\beta_{j,1:K}$$

$$\Sigma \sim \mathcal{IW}(J, \mathbf{R}), \quad \mathbf{R} = \mathbf{I}_J, \quad \Sigma[1,1] = 1$$

$$\beta_{j,k} \sim \mathcal{N}(0, 1000), \quad j = 1, \dots, (J-1), \quad k = 1, \dots, K$$

$$\beta_{J,k} = -\sum_{j=1}^{J-1} \beta_{j,k}$$

$$\mathbf{Z}_{i,j} \sim \mathcal{N}(0, 1000) \in [-10, 10]$$

26.2. Data

```

y <- x1 <- x2 <- c(1:30)
y[1:10] <- 1
y[11:20] <- 2
y[21:30] <- 3
x1[1:10] <- rnorm(10, 25, 2.5)
x1[11:20] <- rnorm(10, 40, 4.0)
x1[21:30] <- rnorm(10, 35, 3.5)
x2[1:10] <- rnorm(10, 2.51, 0.25)
x2[11:20] <- rnorm(10, 2.01, 0.20)
x2[21:30] <- rnorm(10, 2.70, 0.27)
N <- length(y)
J <- 3 #Number of categories in y
K <- 3 #Number of columns to be in design matrix X
R <- diag(J)
X <- matrix(c(rep(1,N),x1,x2),N,K)
mon.names <- "LP"
sigma.temp <- parm.names(list(Sigma=diag(J)), uppertri=1)
parm.names <- c(sigma.temp[2:length(sigma.temp)],
  parm.names(list(beta=matrix(0,(J-1),K), Z=matrix(0,N,J))))
MyData <- list(J=J, K=K, N=N, R=R, X=X, mon.names=mon.names,
  parm.names=parm.names, y=y)

```

26.3. Initial Values

```

Initial.Values <- c(rep(0,length(R[upper.tri(R, diag=TRUE)]))-1),
  rep(0,(J-1)*K), rep(0,N,J))

```

26.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  beta <- matrix(parm[grepl("beta", Data$parm.names)], Data$J-1, Data$K)
  beta <- rbind(beta, colSums(beta)*-1) #Sum to zero constraint
  Sigma <- matrix(NA, Data$J, Data$J)
  Sigma[upper.tri(Sigma, diag=TRUE)] <- c(0, parm[grepl("Sigma",
    Data$parm.names)])
  Sigma[lower.tri(Sigma)] <- Sigma[upper.tri(Sigma)]
  diag(Sigma) <- exp(diag(Sigma))
  Z <- matrix(parm[grepl("Z", Data$parm.names)], Data$N, Data$J)
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  Sigma.prior <- dinvwishart(Sigma, Data$J, Data$R, log=TRUE)
  Z.prior <- sum(dnorm(Z, 0, sqrt(1000), log=TRUE))
}

```

```

### Log-Likelihood
mu <- matrix(0,Data$N,Data$J)
for (j in 1:Data$J) {mu[,j] <- tcrossprod(beta[j,], Data$X)}
Y <- indmat(Data$y)
Z <- ifelse(Z > 10, 10, Z); Z <- ifelse({Y == 0} & {Z > 0}, 0, Z)
Z <- ifelse(Z < -10, -10, Z); Z <- ifelse({Y == 1} & {Z < 0}, 0, Z)
parm[grep("Z", Data$parm.names)] <- as.vector(Z)
LL <- sum(dmvn(Z, mu, Sigma, log=TRUE))
yrep <- apply(Z, 1, which.max)
#eta <- exp(mu)
#p <- eta / rowSums(eta)
### Log-Posterior
LP <- LL + beta.prior + Sigma.prior + Z.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=LP, yhat=yrep, parm=parm)
return(Modelout)
}

```

27. Normal, Multilevel

This is Gelman's school example (Gelman, Carlin, Stern, and Rubin 2004). Note that **LaplacesDemon** is much slower to converge compared to this example that uses the **R2WinBUGS** package (Gelman 2009), an R package on CRAN. However, also note that Laplace's Demon (eventually) provides a better answer (higher ESS, lower DIC, etc.).

27.1. Form

$$\mathbf{y}_j \sim \mathcal{N}(\theta_j, \tau_j^{-1}), \quad j = 1, \dots, J$$

$$\theta_j \sim \mathcal{N}(\theta_\mu, \theta_\tau^{-1}), \quad j = 1, \dots, J$$

$$\theta_\mu \sim \mathcal{N}(0, 1000)$$

$$\theta_\tau = \frac{1}{\theta_\sigma^2}$$

$$\sigma \sim \mathcal{U}(1.0E - 100, 100)$$

$$\tau_j = \sigma_j^{-2}, \quad j = 1, \dots, J$$

27.2. Data

```

J <- 8
y <- c(28.4, 7.9, -2.8, 6.8, -0.6, 0.6, 18.0, 12.2)
sd <- c(14.9, 10.2, 16.3, 11.0, 9.4, 11.4, 10.4, 17.6)
mon.names <- c("LP", "theta.tau")
parm.names <- parm.names(list(theta=rep(0,J), theta.mu=0, sigma=0))

```

```
MyData <- list(J=J, mon.names=mon.names, parm.names=parm.names, sd=sd, y=y)
```

27.3. Initial Values

```
Initial.Values <- c(rep(0,J), 0, 1)
```

27.4. Model

```
Model <- function(parm, Data)
{
  ### Hyperparameters
  theta.mu <- parm[Data$J+1]
  sigma <- interval(parm[grep("sigma", Data$parm.names)], 1.0E-100, 100)
  parm[grep("sigma", Data$parm.names)] <- sigma
  theta.tau <- 1 / sigma^2
  tau.alpha <- 1.0E-3
  tau.beta <- 1.0E-3
  ### Parameters
  theta <- parm[1:Data$J]; tau <- 1/(Data$sd*Data$sd)
  ### Log(Hyperprior and Prior Densities)
  theta.mu.prior <- dnorm(theta.mu, sqrt(1000), log=TRUE)
  sigma.prior <- dunif(sigma, 1.0E-100, 100, log=TRUE)
  tau.prior <- sum(dgamma(tau, tau.alpha, tau.beta, log=TRUE))
  theta.prior <- sum(dnorm(theta, theta.mu, 1/sqrt(theta.tau), log=TRUE))
  ### Log-Likelihood
  LL <- sum(dnorm(Data$y, theta, 1/sqrt(tau), log=TRUE))
  ### Log-Posterior
  LP <- LL + theta.mu.prior + sigma.prior + theta.prior + tau.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP, theta.tau),
    yhat=theta, parm=parm)
  return(Modelout)
}
```

28. Panel, Autoregressive Poisson

28.1. Form

$$\mathbf{Y} \sim \mathcal{P}(\Lambda)$$

$$\Lambda_{1:N,1} = \exp(\alpha + \beta \mathbf{x})$$

$$\Lambda_{1:N,t} = \exp(\alpha + \beta \mathbf{x} + \rho \log(\mathbf{Y}_{1:N,t-1})), \quad t = 2, \dots, T$$

$$\alpha_i \sim \mathcal{N}(\alpha_\mu, \alpha_\sigma^2), \quad i = 1, \dots, N$$

$$\alpha_\mu \sim \mathcal{N}(0, 1000)$$

$$\alpha_\sigma \sim \mathcal{HC}(25)$$

$$\beta \sim \mathcal{N}(0, 1000)$$

$$\rho \sim \mathcal{N}(0, 1000)$$

28.2. Data

```

N <- 10
T <- 10
alpha <- rnorm(N,2,0.5)
rho <- 0.5
beta <- 0.5
x <- runif(N,0,1)
Y <- matrix(NA,N,T)
Y[,1] <- exp(alpha + beta*x)
for (t in 2:T) {Y[,t] <- exp(alpha + beta*x + rho*log(Y[,t-1]))}
Y <- round(Y)
mon.names <- c("LP","alpha.sigma")
parm.names <- parm.names(list(alpha=rep(0,N), alpha.mu=0,
  log.alpha.sigma=0, beta=0, rho=0))
MyData <- list(N=N, T=T, Y=Y, mon.names=mon.names, parm.names=parm.names,
  x=x)

```

28.3. Initial Values

```
Initial.Values <- c(rep(0,N), 0, log(1), 0, 0)
```

28.4. Model

```

Model <- function(parm, Data)
{
  ### Hyperparameters
  alpha.mu <- parm[Data$N+1]
  alpha.sigma <- exp(parm[Data$N+2])
  ### Parameters
  alpha <- parm[1:Data$N]
  beta <- parm[grepl("beta", Data$parm.names)]
  rho <- parm[grepl("rho", Data$parm.names)]
  ### Log(Hyperprior and Prior Densities)
  alpha.mu.prior <- dnorm(alpha.mu, 0, sqrt(1000), log=TRUE)
  alpha.sigma.prior <- dhalfcauchy(alpha.sigma, 25, log=TRUE)
  alpha.prior <- sum(dnorm(alpha, alpha.mu, alpha.sigma, log=TRUE))
  beta.prior <- dnorm(beta, 0, sqrt(1000), log=TRUE)
  rho.prior <- dnorm(rho, 0, sqrt(1000), log=TRUE)
  ### Log-Likelihood

```

```

Lambda <- Data$Y
Lambda[,1] <- exp(alpha + beta*x)
Lambda[,2:Data$T] <- exp(alpha + beta*Data$x +
  rho*log(Data$Y[,1:(Data$T-1)]))
LL <- sum(dpois(Data$Y, Lambda, log=TRUE))
### Log-Posterior
LP <- LL + alpha.prior + alpha.mu.prior + alpha.sigma.prior +
  beta.prior + rho.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,alpha.sigma),
  yhat=Lambda, parm=parm)
return(Modelout)
}

```

29. Poisson Regression

29.1. Form

$$\begin{aligned}
 \mathbf{y} &\sim \mathcal{P}(\lambda) \\
 \lambda &= \exp(\mathbf{X}\beta) \\
 \beta_j &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J
 \end{aligned}$$

29.2. Data

```

N <- 10000
J <- 5
X <- matrix(runif(N*J,-2,2),N,J); X[,1] <- 1
beta <- runif(J,-2,2)
y <- as.vector(round(exp(tcrossprod(beta, X))))
mon.names <- "LP"
parm.names <- parm.names(list(beta=rep(0,J)))
MyData <- list(J=J, X=X, mon.names=mon.names, parm.names=parm.names, y=y)

```

29.3. Initial Values

```
Initial.Values <- rep(0,J)
```

29.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  beta <- parm[1:Data$J]

```



```

### Log(Prior Densities)
beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
### Log-Likelihood
lambda <- exp(tcrossprod(beta, Data$X))
LL <- sum(dpois(Data$y, lambda, log=TRUE))
### Log-Posterior
LP <- LL + beta.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=LP,
  yhat=lambda, parm=parm)
return(Modelout)
}

```

30. Revision, Normal

This example provides both an analytic solution and numerical approximation of the revision of a normal distribution. Given a normal prior distribution (α) and data distribution (β), the posterior (γ) is the revised normal distribution. This is an introductory example of Bayesian inference, and allows the user to experiment numerical approximation, such as with MCMC in *LaplacesDemon*. Note that, regardless of the data sample size N in this example, Laplace Approximation is inappropriate due to asymptotics since the data (β) is perceived by the algorithm as a single datum rather than a collection of data. MCMC, on the other hand, is biased only by the effective number of samples taken of the posterior.

Analytic Solution

```

prior.mu <- 0
prior.sigma <- 10
N <- 10
data.mu <- 1
data.sigma <- 2
posterior.mu <- (prior.sigma^-2 * prior.mu + N * data.sigma^-2 * data.mu) /
  (prior.sigma^-2 + N * data.sigma^-2)
posterior.sigma <- sqrt(1/(prior.sigma^-2 + data.sigma^-2))
posterior.mu
posterior.sigma

```

30.1. Form

$$\begin{aligned}
 \alpha &\sim \mathcal{N}(0, 10) \\
 \beta &\sim \mathcal{N}(1, 2) \\
 \gamma &= \frac{\alpha_{\sigma}^{-2}\alpha + N\beta_{\sigma}^{-2}\beta}{\alpha_{\sigma}^{-2} + N\beta_{\sigma}^{-2}}
 \end{aligned}$$

30.2. Data

```
N <- 10
```

```
mon.names <- c("LP","gamma")
parm.names <- c("alpha","beta")
MyData <- list(N=N, mon.names=mon.names, parm.names=parm.names)
```

30.3. Initial Values

```
Initial.Values <- c(0,0)
```

30.4. Model

```
Model <- function(parm, Data)
{
  ### Hyperparameters
  alpha.mu <- 0
  alpha.sigma <- 10
  beta.mu <- 1
  beta.sigma <- 2
  ### Parameters
  alpha <- parm[1]
  beta <- parm[2]
  ### Log(Prior Density)
  alpha.prior <- dnorm(alpha, alpha.mu, alpha.sigma, log=TRUE)
  ### Log-Likelihood Density
  LL <- dnorm(beta, beta.mu, beta.sigma, log=TRUE)
  ### Posterior
  gamma <- (alpha.sigma^-2 * alpha + N * beta.sigma^-2 * beta) /
    (alpha.sigma^-2 + N * beta.sigma^-2)
  ### Log(Posterior Density)
  LP <- LL + alpha.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,gamma), yhat=LL,
    parm=parm)
  return(Modelout)
}
```

31. Robust Regression

By replacing the normal distribution with the Student t distribution, linear regression is often called robust regression. As an alternative approach to robust regression, consider Laplace regression (see section 19).

31.1. Form

$$\mathbf{y} \sim t(\mu, \sigma^2, \nu)$$

$$\mu = \mathbf{X}\beta$$

$$\beta_j \sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J$$

$$\sigma \sim \mathcal{HC}(25)$$

$$\nu \sim \mathcal{HC}(25)$$

31.2. Data

```

N <- 10000
J <- 5
X <- matrix(1,N,J)
for (j in 2:J) {X[,j] <- rnorm(N,runif(1,-3,3),runif(1,0.1,1))}
beta <- runif(J,-3,3)
e <- rnorm(N,0,0.1)
y <- as.vector(tcrossprod(beta, X) + e)
mon.names <- c("LP", "sigma", "nu")
parm.names <- parm.names(list(beta=rep(0,J), log.sigma=0, log.nu=0))
MyData <- list(J=J, X=X, mon.names=mon.names, parm.names=parm.names, y=y)

```

31.3. Initial Values

```
Initial.Values <- c(rep(0,J), log(1), log(2))
```

31.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  beta <- parm[1:Data$J]
  sigma <- exp(parm[Data$J+1])
  nu <- exp(parm[Data$J+2])
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  sigma.prior <- dhalfcauchy(sigma, 25, log=TRUE)
  nu.prior <- dhalfcauchy(nu, 25, log=TRUE)
  ### Log-Likelihood
  mu <- tcrossprod(beta, Data$X)
  LL <- sum(dst(Data$y, mu, sigma, nu, log=TRUE))
  ### Log-Posterior
  LP <- LL + beta.prior + sigma.prior + nu.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,sigma,nu), yhat=mu,
    parm=parm)
  return(Modelout)
}

```

32. Seemingly Unrelated Regression (SUR)

The following data was used by Zellner (1962) when introducing the Seemingly Unrelated Regression methodology.

32.1. Form

$$\begin{aligned} \mathbf{Y}_{t,k} &\sim \mathcal{N}_K(\mu_{t,k}, \Sigma), \quad t = 1, \dots, T; \quad k = 1, \dots, K \\ \mu_{1,t} &= \alpha_1 + \alpha_2 \mathbf{X}_{t,1} + \alpha_3 \mathbf{X}_{t,2}, \quad t = 1, \dots, T \\ \mu_{2,t} &= \beta_1 + \beta_2 \mathbf{X}_{t,3} + \beta_3 \mathbf{X}_{t,4}, \quad t = 1, \dots, T \\ \Sigma &= \Omega^{-1} \\ \Omega &\sim \mathcal{W}(K, \mathbf{S}), \quad \mathbf{S} = \mathbf{I}_K \\ \alpha_j &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J \\ \beta_j &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J \end{aligned}$$

where J=3, K=2, and T=20.

32.2. Data

```
T <- 20
year <- c(1935, 1936, 1937, 1938, 1939, 1940, 1941, 1942, 1943, 1944, 1945, 1946,
          1947, 1948, 1949, 1950, 1951, 1952, 1953, 1954)
IG <- c(33.1, 45.0, 77.2, 44.6, 48.1, 74.4, 113.0, 91.9, 61.3, 56.8, 93.6, 159.9,
        147.2, 146.3, 98.3, 93.5, 135.2, 157.3, 179.5, 189.6)
VG <- c(1170.6, 2015.8, 2803.3, 2039.7, 2256.2, 2132.2, 1834.1, 1588.0, 1749.4,
        1687.2, 2007.7, 2208.3, 1656.7, 1604.4, 1431.8, 1610.5, 1819.4, 2079.7,
        2371.6, 2759.9)
CG <- c(97.8, 104.4, 118.0, 156.2, 172.6, 186.6, 220.9, 287.8, 319.9, 321.3, 319.6,
        346.0, 456.4, 543.4, 618.3, 647.4, 671.3, 726.1, 800.3, 888.9)
IW <- c(12.93, 25.90, 35.05, 22.89, 18.84, 28.57, 48.51, 43.34, 37.02, 37.81,
        39.27, 53.46, 55.56, 49.56, 32.04, 32.24, 54.38, 71.78, 90.08, 68.60)
VW <- c(191.5, 516.0, 729.0, 560.4, 519.9, 628.5, 537.1, 561.2, 617.2, 626.7,
        737.2, 760.5, 581.4, 662.3, 583.8, 635.2, 723.8, 864.1, 1193.5, 1188.9)
CW <- c(1.8, 0.8, 7.4, 18.1, 23.5, 26.5, 36.2, 60.8, 84.4, 91.2, 92.4, 86.0, 111.1,
        130.6, 141.8, 136.7, 129.7, 145.5, 174.8, 213.5)
Y <- matrix(c(IG, IW), T, 2)
S <- diag(NCOL(Y))
mon.names <- c("LP", "Sigma[1,1]", "Sigma[2,1]", "Sigma[1,2]", "Sigma[2,2]")
parm.names <- parm.names(list(alpha=rep(0,3), beta=rep(0,3),
                             Omega=diag(2)), uppertri=c(0,0,1))
MyData <- list(S=S, T=T, Y=Y, CG=CG, CW=CW, IG=IG, IW=IW, VG=VG, VW=VW,
               mon.names=mon.names, parm.names=parm.names)
```

32.3. Initial Values

```
Initial.Values <- c(rep(0,3), rep(0,3), S[upper.tri(S, diag=TRUE)])
```

32.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  alpha <- parm[1:3]
  beta <- parm[4:6]
  Omega <- matrix(parm[c(7,8,8,9)], NROW(Data$S), NROW(Data$S))
  Sigma <- solve(Omega)
  ### Log(Prior Densities)
  alpha.prior <- sum(dnorm(alpha, 0, sqrt(1000), log=TRUE))
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  Omega.prior <- dwishart(Omega, NROW(Data$S), Data$S, log=TRUE)
  ### Log-Likelihood
  mu <- matrix(0, Data$T, 2)
  mu[,1] <- alpha[1] + alpha[2]*Data$CG + alpha[3]*Data$VG
  mu[,2] <- beta[1] + beta[2]*Data$CW + beta[3]*Data$VW
  LL <- sum(dmvn(Data$Y, mu, Sigma, log=TRUE))
  ### Log-Posterior
  LP <- LL + alpha.prior + beta.prior + Omega.prior
  Modelout <- list(LP=LP, Dev=-2*LL,
    Monitor=c(LP, as.vector(Sigma)), yhat=mu, parm=parm)
  return(Modelout)
}

```

33. Space-Time, Nonseparable

This approach to space-time or spatiotemporal modeling applies kriging both to the stationary spatial and temporal components, where space is continuous and time is discrete. Matrix Ξ contains the space-time effects. Spatial coordinates are given in longitude and latitude for $s = 1, \dots, S$ points in space and measurements are taken across time-periods $t = 1, \dots, T$ for $\mathbf{Y}_{s,t}$. The dependent variable is also a function of design matrix \mathbf{X} and regression effects vector β . For more information on kriging, see section 18. This example uses a nonseparable, stationary covariance function in which space and time are separable only when $\psi = 0$.

33.1. Form

$$\begin{aligned}
 \mathbf{Y}_{s,t} &\sim \mathcal{N}(\mu_{s,t}, \sigma_1^2), \quad s = 1, \dots, S, \quad t = 1, \dots, T \\
 \mu &= \mathbf{X}\beta + \Xi \\
 \Xi &\sim \mathcal{N}_{ST}(\Xi_\mu, \Sigma) \\
 \Sigma &= \sigma_2^2 \exp \left(-\frac{\mathbf{D}_S^\kappa}{\phi_1} - \frac{\mathbf{D}_T^\lambda}{\phi_2} - \psi \frac{\mathbf{D}_S^\kappa \mathbf{D}_T^\lambda}{\phi_1 \phi_2} \right) \\
 \beta_j &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J
 \end{aligned}$$

$$\phi_k \sim \mathcal{U}(1, 5), \quad k = 1, \dots, 2$$

$$\sigma_k \sim \mathcal{HC}(25), \quad k = 1, \dots, 2$$

$$\psi \sim \mathcal{HC}(25)$$

$$\Xi_\mu = 0$$

$$\kappa = 1, \quad \lambda = 1$$

33.2. Data

```

S <- 10
T <- 5
longitude <- runif(S,0,100)
latitude <- runif(S,0,100)
D.S <- as.matrix(dist(cbind(rep(longitude,T),rep(latitude,T)), diag=TRUE,
  upper=TRUE))
D.T <- as.matrix(dist(cbind(rep(1:T,each=S),rep(1:T,each=S)), diag=TRUE,
  upper=TRUE))
Sigma <- 10000 * exp(-D.S/3 - D.T/2 - 0.2*(D.S/3)*(D.T/2))
Xi <- as.vector(apply(rmvn(1000, rep(0,S*T), Sigma), 2, mean))
Xi <- matrix(Xi,S,T)
beta <- c(50,2)
X <- matrix(runif(S*2,-2,2),S,2); X[,1] <- 1
mu <- as.vector(tcrossprod(beta, X))
Y <- mu + Xi
mon.names <- c("LP","psi","sigma[1]","sigma[2]")
parm.names <- parm.names(list(Xi=matrix(0,S,T), beta=rep(0,2),
  phi=rep(0,2), log.sigma=rep(0,2), log.psi=0))
MyData <- list(D.S=D.S, D.T=D.T, S=S, T=T, X=X, Y=Y, mon.names=mon.names,
  parm.names=parm.names)

```

33.3. Initial Values

```
Initial.Values <- c(rep(0,S*T), mean(Y), 0, rep(1,2), rep(0,2), 0)
```

33.4. Model

```

Model <- function(parm, Data)
{
  ### Hyperparameters
  Xi.mu <- rep(0,Data$S*Data$T)
  ### Parameters
  beta <- parm[grep("beta", Data$parm.names)]
  Xi <- parm[grep("Xi", Data$parm.names)]
  kappa <- 1; lambda <- 1
  sigma <- exp(parm[grep("log.sigma", Data$parm.names)])
}

```

```

phi <- interval(parm[grep("phi", Data$parm.names)], 1, 5)
parm[grep("phi", Data$parm.names)] <- phi
psi <- exp(parm[grep("log.psi", Data$parm.names)])
Sigma <- sigma[2]*sigma[2] * exp(-(Data$D.S / phi[1])^kappa -
  (Data$D.T / phi[2])^lambda -
  psi*(Data$D.S / phi[1])^kappa * (Data$D.T / phi[2])^lambda)
### Log(Prior Densities)
beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
Xi.prior <- dmvn(Xi, Xi.mu, Sigma, log=TRUE)
sigma.prior <- sum(dhalfcauchy(sigma, 25, log=TRUE))
phi.prior <- sum(dunif(phi, 1, 5, log=TRUE))
psi.prior <- dhalfcauchy(psi, 25, log=TRUE)
### Log-Likelihood
Xi <- matrix(Xi, Data$S, Data$T)
mu <- as.vector(tcrossprod(beta, Data$X)) + Xi
LL <- sum(dnorm(Data$Y, mu, sigma[1], log=TRUE))
### Log-Posterior
LP <- LL + beta.prior + Xi.prior + sigma.prior + phi.prior + psi.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,psi,sigma),
  yhat=mu, parm=parm)
return(Modelout)
}

```

34. Space-Time, Separable

This introductory approach to space-time or spatiotemporal modeling applies kriging both to the stationary spatial and temporal components, where space is continuous and time is discrete. Vector ζ contains the spatial effects and vector θ contains the temporal effects. Spatial coordinates are given in longitude and latitude for $s = 1, \dots, S$ points in space and measurements are taken across time-periods $t = 1, \dots, T$ for $\mathbf{Y}_{s,t}$. The dependent variable is also a function of design matrix \mathbf{X} and regression effects vector β . For more information on kriging, see section 18. This example uses separable space-time covariances, which is more convenient but usually less appropriate than a nonseparable covariance function.

34.1. Form

$$\begin{aligned}
\mathbf{Y}_{s,t} &\sim \mathcal{N}(\mu_{s,t}, \tau_1^{-1}), \quad s = 1, \dots, S, \quad t = 1, \dots, T \\
\mu_{s,t} &= \mathbf{X}_{s,1:T} \beta + \zeta_s + \Theta_{s,t} \\
\Theta_{s,1:T} &= \theta \\
\theta &\sim \mathcal{N}(\theta_\mu, \Sigma_T) \\
\Sigma_T &= \frac{1}{\tau_3} \exp(-\phi_2 \mathbf{D}_T)^\lambda \\
\zeta &\sim \mathcal{N}(\zeta_\mu, \Sigma_S)
\end{aligned}$$

$$\begin{aligned}\Sigma_S &= \frac{1}{\tau_2} \exp(-\phi_1 \mathbf{D}_S)^\kappa \\ \beta_j &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, 2 \\ \tau_k &\sim \mathcal{G}(0.001, 0.001), \quad k = 1, \dots, 3 \\ \phi_k &\sim \mathcal{U}(1, 5), \quad k = 1, \dots, 2 \\ \zeta_\mu &= 0 \\ \theta_\mu &= 0 \\ \kappa &= 1, \quad \lambda = 1\end{aligned}$$

34.2. Data

```
S <- 20
T <- 10
longitude <- runif(S,0,100)
latitude <- runif(S,0,100)
D.S <- as.matrix(dist(cbind(longitude,latitude), diag=TRUE, upper=TRUE))
Sigma.S <- 10000 * exp(-1.5 * D.S)
zeta <- as.vector(apply(rmvn(1000, rep(0,S), Sigma.S), 2, mean))
D.T <- as.matrix(dist(cbind(c(1:T),c(1:T)), diag=TRUE, upper=TRUE))
Sigma.T <- 10000 * exp(-3 * D.T)
theta <- as.vector(apply(rmvn(1000, rep(0,T), Sigma.T), 2, mean))
Theta <- matrix(theta,S,T,byrow=TRUE)
beta <- c(50,2)
X <- matrix(runif(S*2,-2,2),S,2); X[,1] <- 1
mu <- as.vector(tcrossprod(beta, X))
Y <- mu + zeta + Theta + matrix(rnorm(S*T,0,0.1),S,T)
mon.names <- c("LP","tau[1]","tau[2]","tau[3]")
parm.names <- parm.names(list(zeta=rep(0,S), theta=rep(0,T),
  beta=rep(0,2), phi=rep(0,2), log.tau=rep(0,3))
MyData <- list(D.S=D.S, D.T=D.T, S=S, T=T, X=X, Y=Y, mon.names=mon.names,
  parm.names=parm.names)
```

34.3. Initial Values

```
Initial.Values <- c(rep(0,S), rep(0,T), rep(0,2), rep(1,2), rep(0,3))
```

34.4. Model

```
Model <- function(parm, Data)
{
  ### Hyperparameters
  beta.mu <- 0
  beta.tau <- 1.0E-3
  zeta.mu <- rep(0,Data$S)
```



```

theta.mu <- rep(0,Data$T)
tau.alpha <- rep(1.0E-3,3)
tau.beta <- rep(1.0E-3,3)
### Parameters
beta <- parm[grep("beta", Data$parm.names)]
zeta <- parm[grep("zeta", Data$parm.names)]
theta <- parm[grep("theta", Data$parm.names)]
kappa <- 1; lambda <- 1
tau <- exp(parm[grep("log.tau", Data$parm.names)])
phi <- interval(parm[grep("phi", Data$parm.names)], 1, 5)
parm[grep("phi", Data$parm.names)] <- phi
Sigma.S <- 1/tau[2] * exp(-phi[1] * Data$D.S)^kappa
Sigma.T <- 1/tau[3] * exp(-phi[2] * Data$D.T)^lambda
### Log(Prior Densities)
beta.prior <- sum(dnorm(beta, beta.mu, 1/sqrt(beta.tau), log=TRUE))
zeta.prior <- dmvn(zeta, zeta.mu, Sigma.S, log=TRUE)
theta.prior <- dmvn(theta, theta.mu, Sigma.T, log=TRUE)
tau.prior <- sum(dgamma(tau, tau.alpha, tau.beta, log=TRUE))
phi.prior <- sum(dunif(phi, 1, 5, log=TRUE))
### Log-Likelihood
Theta <- matrix(theta, Data$S, Data$T, byrow=TRUE)
mu <- as.vector(tcrossprod(beta, Data$X)) + zeta + Theta
LL <- sum(dnorm(Data$Y, mu, 1/sqrt(tau[1]), log=TRUE))
### Log-Posterior
LP <- LL + beta.prior + zeta.prior + theta.prior + tau.prior +
  phi.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP,tau),
  yhat=mu, parm=parm)
return(Modelout)
}

```

35. Survival Analysis

Although the dependent variable is usually denoted as \mathbf{t} in survival analysis, it is denoted here as \mathbf{y} so Laplace's Demon recognizes it as a dependent variable for posterior predictive checks.

35.1. Form

$$\mathbf{y}_i \sim \mathcal{WETB}(\gamma, \mu_i), \quad i = 1, \dots, N$$

$$\mu = \exp(\mathbf{X}\beta)$$

$$\beta_j \sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J$$

$$\gamma \sim \mathcal{G}(1, 0.001)$$

35.2. Data

```

N <- 50
J <- 5
X <- matrix(runif(N*J,-2,2),N,J); X[,1] <- 1
beta <- runif(J,-1,1)
y <- as.vector(round(exp(tcrossprod(beta, X)))) + 1
mon.names <- c("LP","gamma")
parm.names <- parm.names(list(beta=rep(0,J), log.gamma=0))
MyData <- list(J=J, N=N, X=X, mon.names=mon.names, parm.names=parm.names,
               y=y)

```

35.3. Initial Values

```
Initial.Values <- c(rep(0,J), log(1))
```

35.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  beta <- parm[1:Data$J]
  gamma <- exp(parm[Data$J+1])
  ### Log(Prior Densities)
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  gamma.prior <- dgamma(gamma, 1, 1.0E-3, log=TRUE)
  ### Log-Likelihood
  mu <- exp(tcrossprod(beta, Data$X)) + 1
  h <- (gamma/lambda)*(Data$y/lambda)^(gamma-1)
  S <- exp(-mu * Data$y^gamma)
  LL <- sum(dweibull(Data$y, gamma, mu, log=TRUE))
  ### Log-Posterior
  LP <- LL + beta.prior + gamma.prior
  Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP, gamma),
                  yhat=mu, parm=parm)
  return(Modelout)
}

```

36. Variable Selection

This example uses a modified form of the random-effects (or global adaptation) Stochastic Search Variable Selection (SSVS) algorithm presented in [O'Hara and Sillanpaa \(2009\)](#), which selects variables according to practical significance rather than statistical significance. Here, SSVS is applied to linear regression, though this method is widely applicable. For J variables, each regression effects vector β_j is conditional on γ_j , a binary inclusion variable. Each β_j is a discrete mixture distribution with respect to $\gamma_j = 0$ or $\gamma_j = 1$, with precision 100 or $\beta_\sigma = 0.1$,

respectively. As with other representations of SSVS, these precisions may require tuning.

With other representations of SSVS, each γ_j is Bernoulli-distributed, though this would be problematic in Laplace's Demon, because γ_j would be in the list of parameters (rather than monitors), and would not be stationary due to switching behavior. To keep γ in the monitors, an uninformative normal density is placed on each prior δ_j , with mean $1/J$ for J variables and variance 1000. Each δ_j is transformed with the inverse logit and rounded to γ_j . Note that $\lfloor x + 0.5 \rfloor$ means to round x . The prior for δ can be manipulated to influence sparseness. When the goal is to select the best model, each $\mathbf{X}_{1:N,j}$ is retained for a future run when the posterior mean of $\gamma_j \geq 0.5$. When the goal is model-averaging, the results of this model may be used directly.

36.1. Form

$$\begin{aligned} \mathbf{y} &\sim \mathcal{N}(\mu, \sigma^2) \\ \mu &= \mathbf{X}\beta \\ (\beta_j | \gamma_j) &\sim (1 - \gamma_j)\mathcal{N}(0, 0.01) + \gamma_j\mathcal{N}(0, \beta_\sigma^2) \quad j = 1, \dots, J \\ \beta_\sigma &\sim \mathcal{HC}(25) \\ \gamma_j &= \lfloor \frac{1}{1 + \exp(-\delta_j)} + 0.5 \rfloor, \quad j = 1, \dots, J \\ \delta_j &\sim \mathcal{N}(0, 10) \in [-100, 100], \quad j = 1, \dots, J \\ \sigma &\sim \mathcal{HC}(25) \end{aligned}$$

36.2. Data

```
data(demonsnacks)
N <- NROW(demonsnacks)
J <- NCOL(demonsnacks)
y <- log(demonsnacks$Calories)
X <- cbind(1, as.matrix(demonsnacks[,c(1,3:10)]))
for (j in 2:J) {X[,j] <- CenterScale(X[,j])}
mon.names <- c("LP", "min.beta.sigma", "sigma",
  parm.names(list(gamma=rep(0,J)))
parm.names <- parm.names(list(beta=rep(0,J), delta=rep(0,J),
  log.beta.sigma=0, log.sigma=0))
MyData <- list(J=J, X=X, mon.names=mon.names, parm.names=parm.names, y=y)
```

36.3. Initial Values

```
Initial.Values <- c(rep(0,J), rep(0,J), log(1), log(1))
```

36.4. Model

```
Model <- function(parm, Data)
{
```

```

### Hyperparameters
beta.sigma <- exp(parm[grepl("log.beta.sigma", Data$parm.names)])
delta <- interval(parm[grepl("delta", Data$parm.names)], -100, 100)
parm[grepl("delta", Data$parm.names)] <- delta
### Parameters
beta <- parm[1:Data$J]
gamma <- round(invlogit(delta))
beta.sigma <- ifelse(gamma == 0, 0.1, beta.sigma)
sigma <- exp(parm[grepl("log.sigma", Data$parm.names)])
### Log(Hyperprior and Prior Densities)
beta.prior <- sum(dnorm(beta, 0, beta.sigma, log=TRUE))
beta.sigma.prior <- sum(dhalfcauchy(beta.sigma, 25, log=TRUE))
delta.prior <- sum(dtrunc(delta, "norm", a=-100, b=100,
  mean=logit(1/Data$J), sd=sqrt(1000), log=TRUE))
sigma.prior <- dhalfcauchy(sigma, 25, log=TRUE)
### Log-Likelihood
mu <- tcrossprod(beta, Data$X)
LL <- sum(dnorm(y, mu, sigma, log=TRUE))
### Log-Posterior
LP <- LL + beta.prior + beta.sigma.prior + delta.prior + sigma.prior
Modelout <- list(LP=LP, Dev=-2*LL, Monitor=c(LP, min(beta.sigma),
  sigma, gamma), yhat=mu, parm=parm)
return(Modelout)
}

```

37. Zero-Inflated Poisson (ZIP)

37.1. Form

$$\begin{aligned}
\mathbf{y} &\sim \mathcal{P}(\Lambda_{1:N,2}) \\
\mathbf{z} &\sim \mathcal{BERN}(\Lambda_{1:N,1}) \\
\mathbf{z}_i &= \begin{cases} 1 & \text{if } y_i = 0 \\ 0 & \text{otherwise} \end{cases} \\
\Lambda_{i,2} &= \begin{cases} 0 & \text{if } \Lambda_{i,1} \geq 0.5 \\ \Lambda_{i,2} & \text{otherwise} \end{cases} \\
\Lambda_{1:N,1} &= \frac{1}{1 + \exp(-\mathbf{X}_1 \alpha)} \\
\Lambda_{1:N,2} &= \exp(\mathbf{X}_2 \beta) \\
\alpha_j &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J_1 \\
\beta_j &\sim \mathcal{N}(0, 1000), \quad j = 1, \dots, J_2
\end{aligned}$$

37.2. Data

```

N <- 1000
J1 <- 4
J2 <- 3
X1 <- matrix(runif(N*J1,-2,2),N,J1); X1[,1] <- 1
X2 <- matrix(runif(N*J2,-2,2),N,J2); X2[,1] <- 1
alpha <- runif(J1,-1,1)
beta <- runif(J2,-1,1)
p <- as.vector(invlogit(tcrossprod(alpha, X1) + rnorm(N,0,0.1)))
mu <- as.vector(round(exp(tcrossprod(beta, X2) + rnorm(N,0,0.1))))
y <- ifelse(p > 0.5, 0, mu)
z <- ifelse(y == 0, 1, 0)
mon.names <- "LP"
parm.names <- parm.names(list(alpha=rep(0,J1), beta=rep(0,J2)))
MyData <- list(J1=J1, J2=J2, N=N, X1=X1, X2=X2, mon.names=mon.names,
               parm.names=parm.names, y=y, z=z)

```

37.3. Initial Values

```
Initial.Values <- rep(0,J1+J2)
```

37.4. Model

```

Model <- function(parm, Data)
{
  ### Parameters
  alpha <- parm[1:Data$J1]
  beta <- parm[grep("beta", Data$parm.names)]
  ### Log(Prior Densities)
  alpha.prior <- sum(dnorm(alpha, 0, sqrt(1000), log=TRUE))
  beta.prior <- sum(dnorm(beta, 0, sqrt(1000), log=TRUE))
  ### Log-Likelihood
  Lambda <- matrix(NA, Data$N, 2)
  Lambda[,1] <- invlogit(tcrossprod(alpha, Data$X1))
  Lambda[,2] <- exp(tcrossprod(beta, Data$X2))
  Lambda[,2] <- ifelse(Lambda[,1] >= 0.5, 0, Lambda[,2])
  LL1 <- sum(dbern(Data$z, Lambda[,1], log=TRUE))
  LL2 <- sum(dpois(Data$y, Lambda[,2], log=TRUE))
  ### Log-Posterior
  LP <- LL1 + LL2 + alpha.prior + beta.prior
  Modelout <- list(LP=LP, Dev=-2*LL2, Monitor=LP,
                  yhat=Lambda[,2], parm=parm)
  return(Modelout)
}

```

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