Airlines Fatalities

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Airlines Fatalities

Introduction

In this project we are analyzing data collected from 1976 to 2001 about airline fatal accidents by the International Civil Aviation Organization in Montreal , Canada (www.icao.int).

Our goal is to get a good prediction about future fatalitis through a bayesian approach.

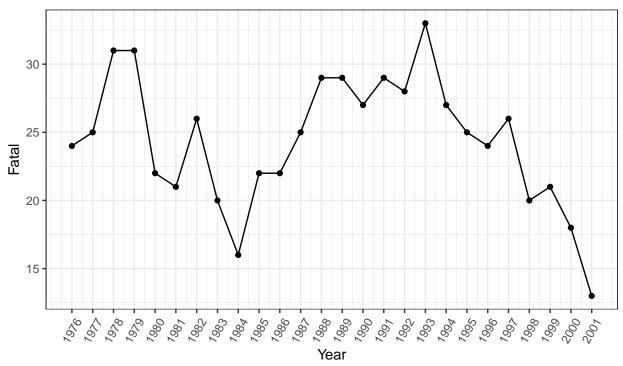
Our data is structured with four columns, year, fatal, miles, rate.

"Passenger miles" are in units of 10^{11} and the "accident rate" is the number of fatal accidents per 10^{11} passenger miles.

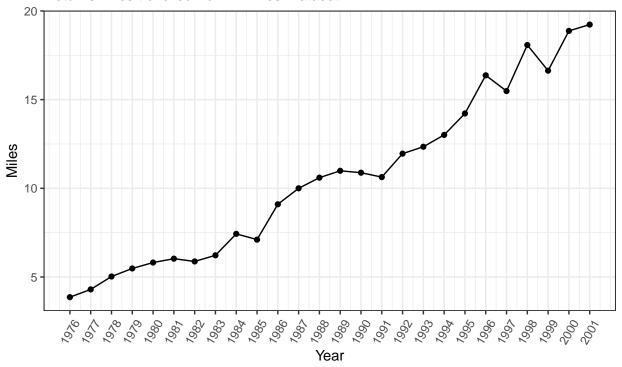
year	fatal	miles	rate
1976	24	3.863	6.213
1977	25	4.3	5.814
1978	31	5.027	6.167
1979	31	5.481	5.656
1980	22	5.814	3.784
1981	21	6.033	3.481

Following three descriptive time series plot respectively about the number of fatal accidents, the miles flown and the accident rate per each year.

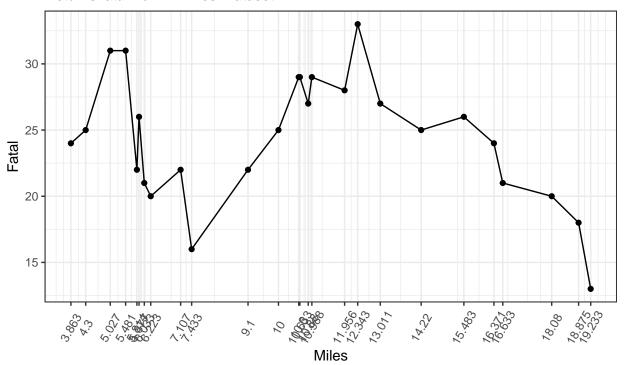
Time Series Chart Returns fatal from 'Airlines' Dataset



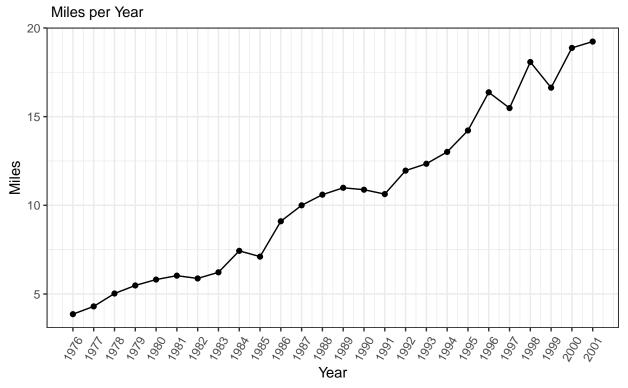
Time Series Chart Returns miles travelled from 'Airlines' Dataset



Time Series Chart Returns fatal from 'Airlines' Dataset

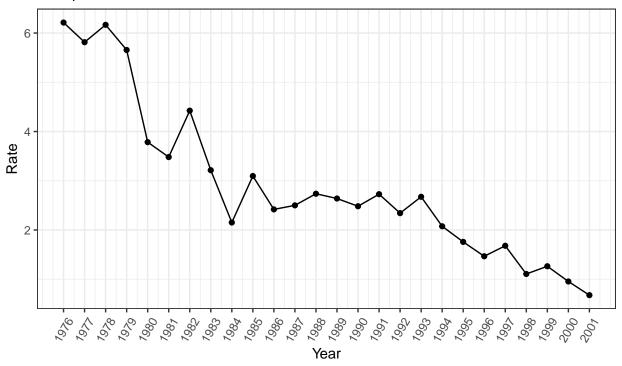


Time Series Chart



Time Series Chart

Rate per Year



Source: American Airlines

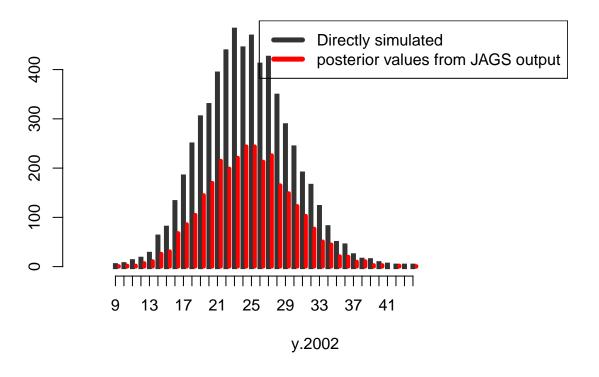
JAGS MODELS:

First model

```
cat( "model {
  for( i in 1:I ) {
    fatal[i] ~ dpois(mu)
    mu ~ dgamma(0.01,0.01)
 }", file="a1.jag" )
a1.par <- c("mu", "fatal[27]", "fatal[28]", "fatal[29]")</pre>
a1.ini <- list(list( mu=22 ),
              list( mu=23 ),
              list( mu=24 ) )
a1.dat <- list( fatal = c(airline$fatal, NA, NA, NA), I=29 )
\# Model compilation and burn-in
a1.mod <- jags.model( file = "a1.jag",
                     data = a1.dat,
                     inits = a1.ini,
                    n.chains = 3,
                    n.adapt = 1000)
```

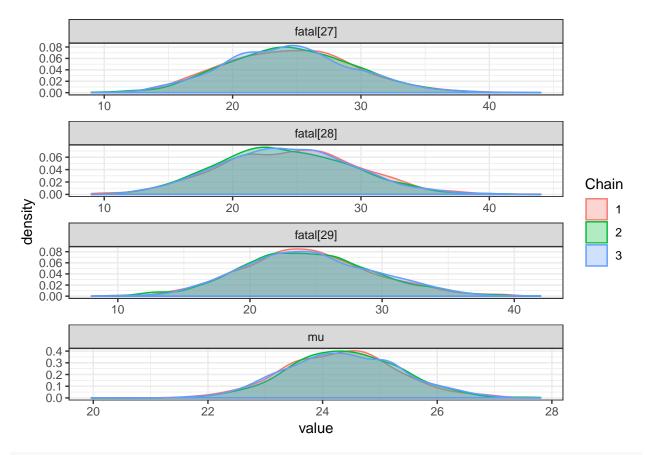
```
## Compiling model graph
##
      Resolving undeclared variables
##
      Allocating nodes
## Graph information:
##
      Observed stochastic nodes: 26
##
      Unobserved stochastic nodes: 4
##
      Total graph size: 32
##
## Initializing model
update(a1.mod, 1000)
# Sampling from the posterior
a1.res <- coda.samples( a1.mod,
                      var = a1.par,
                      n.iter = 10000,
                      thin = 10)
summary( a1.res )
##
## Iterations = 1010:11000
## Thinning interval = 10
## Number of chains = 3
## Sample size per chain = 1000
## 1. Empirical mean and standard deviation for each variable,
##
      plus standard error of the mean:
##
                       SD Naive SE Time-series SE
##
              Mean
## fatal[27] 24.46 4.9572 0.09051
                                          0.09053
## fatal[28] 24.26 5.1762 0.09450
                                          0.09448
## fatal[29] 24.41 4.9616 0.09059
                                          0.08926
## mu
             24.37 0.9799 0.01789
                                          0.01749
## 2. Quantiles for each variable:
##
              2.5%
                     25%
                           50%
##
                                 75% 97.5%
## fatal[27] 15.00 21.00 24.00 28.00 34.00
## fatal[28] 15.00 21.00 24.00 28.00 35.00
## fatal[29] 15.00 21.00 24.00 28.00 35.00
             22.48 23.68 24.36 25.04 26.34
theta <- rgamma(6000, 634, 26)
y.2002 <- rpois(6000,theta)
plot( main = 'Posterior predictive distribution of y in 2002', table(y.2002), type="h", lwd=5, lend=2, c
tpr <- table( as.matrix( a1.res[,"fatal[27]"] ) )</pre>
points( as.numeric(names(tpr))+0.4, tpr, type="h", col="red", lwd=4 )
legend("topright", c('Directly simulated', 'posterior values from JAGS output'),col=c(gray(0.2),'red'),l
```

Posterior predictive distribution of y in 2002

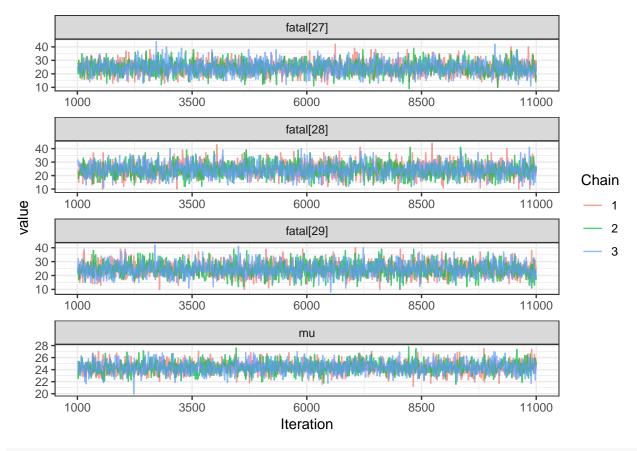


Diagnostics for model 1

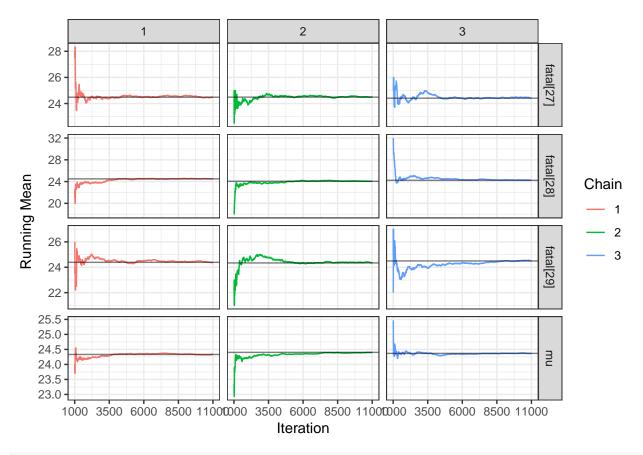
result = ggs(a1.res)
ggs_density(result)



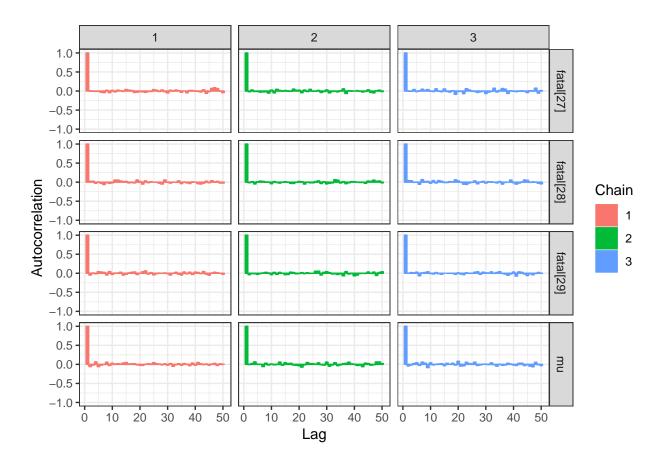
ggs_traceplot(result)



ggs_running(result)



ggs_autocorrelation(result)



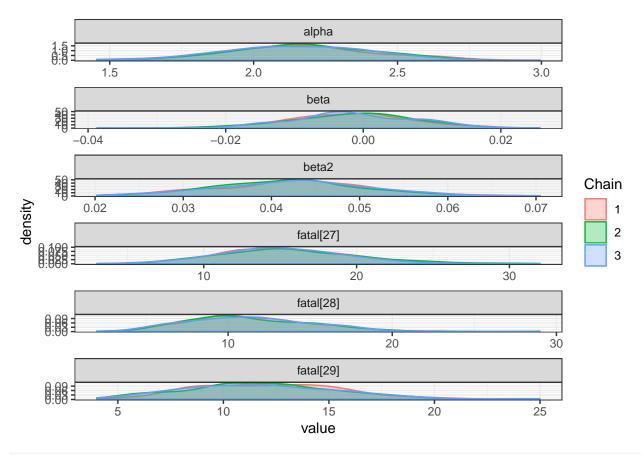
Third model

```
library(rjags)
 cat( "model
for( i in 1:I )
 fatal[i] ~ dpois( mu[i] )
 log(mu[i]) <- alpha + beta*miles[i]+ beta2 * miles[i]*rate[i]</pre>
 }
 alpha ~ dnorm(0,0.0001)
 beta ~ dnorm(0.0001,0.00001)
 beta2 ~ dnorm(0.0001,0.00001)
}",
 file="a3.jag" )
 a3.ini= list(
    list(alpha = 0.01, beta = 0.6, beta2=0.1),
    list(alpha = 0.2, beta = 0.2, beta2=0.2),
    list(alpha = 0.7, beta = 0.7, beta2=0.3))
 a3.dat <- list( rate= c(airline$rate, 0.7080, 0.3, 0.4), fatal=c(airline$fatal, NA, NA, NA), miles=c(airline$
 a3.par <- c("alpha", "beta", "beta2", "fata1[27]", "fata1[28]", "fata1[29]")
```

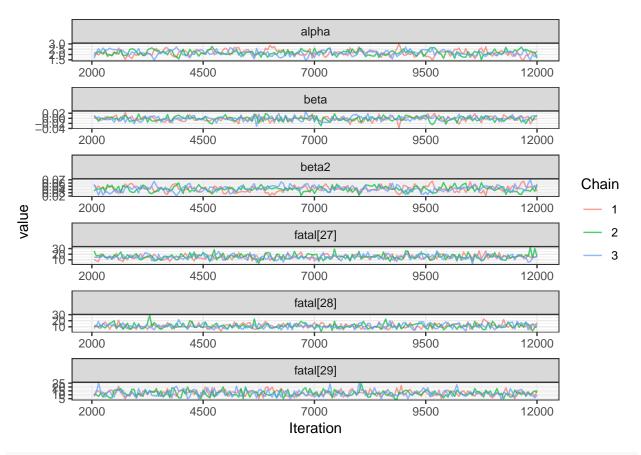
```
# Model compilation and burn-in
a3.mod <- jags.model(file = "a3.jag",data = a3.dat,inits = a3.ini,n.chains = 3,n.adapt = 1000)
## Compiling model graph
##
      Resolving undeclared variables
##
      Allocating nodes
## Graph information:
##
      Observed stochastic nodes: 26
##
      Unobserved stochastic nodes: 6
##
      Total graph size: 210
##
## Initializing model
update(a3.mod, 1000)
 # Sampling from the posterior
a3.res <- coda.samples( a3.mod,var = a3.par,n.iter = 10000,thin = 50 )
summary( a3.res )
##
## Iterations = 2050:12000
## Thinning interval = 50
## Number of chains = 3
## Sample size per chain = 200
##
## 1. Empirical mean and standard deviation for each variable,
##
      plus standard error of the mean:
##
##
                             SD Naive SE Time-series SE
                  Mean
## alpha
              2.157789 0.255229 0.0104197
                                               0.0156726
             -0.001531 0.008970 0.0003662
## beta
                                               0.0003665
## beta2
              0.042269 0.008507 0.0003473
                                               0.0005480
## fatal[27] 15.340000 4.278424 0.1746659
                                               0.1625348
## fatal[28] 11.373333 3.783448 0.1544586
                                               0.1423056
## fatal[29] 11.998333 3.683424 0.1503751
                                              0.1366078
## 2. Quantiles for each variable:
##
##
                 2.5%
                            25%
                                      50%
                                                 75%
                                                        97.5%
## alpha
             1.64099 1.990990 2.162948 2.326468 2.63506
## beta
             -0.01880 -0.006728 -0.001383 0.004598 0.01544
## beta2
              0.02564 \quad 0.036597 \quad 0.042649 \quad 0.047512 \quad 0.05911
## fatal[27] 8.00000 12.000000 15.000000 18.000000 24.02500
## fatal[28] 5.00000 9.000000 11.000000 14.000000 19.00000
## fatal[29] 5.00000 9.000000 12.000000 14.000000 19.02500
```

Model 3 diagnostics

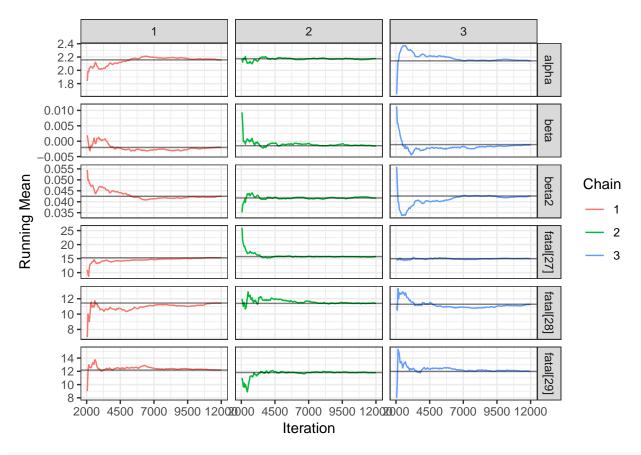
```
result3 = ggs(a3.res)
ggs_density(result3)
```



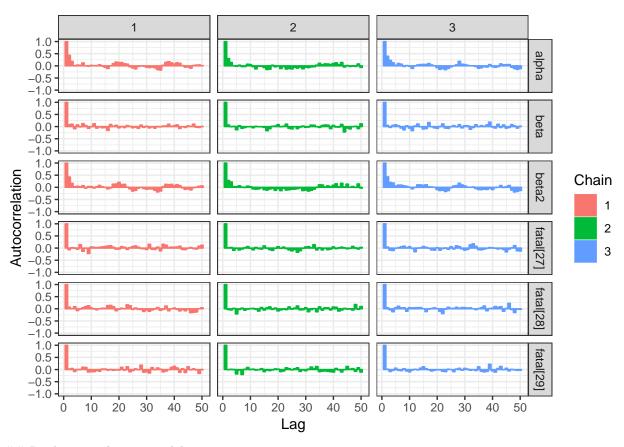
ggs_traceplot(result3)



ggs_running(result3)

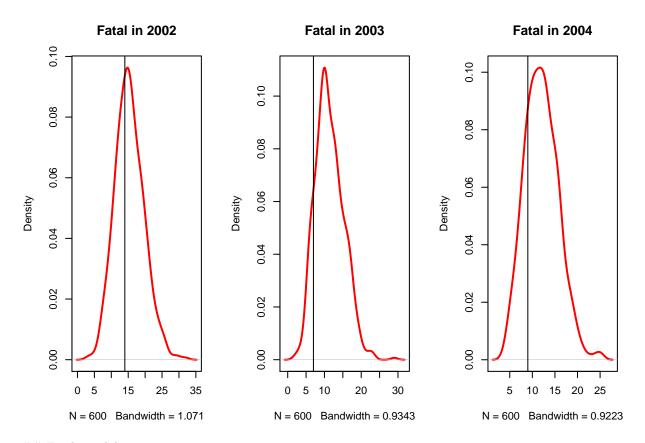


ggs_autocorrelation(result3)



```
## Predicting values in model 3
```

```
par(mfrow=c(1,3))
a3.m <- as.matrix(a3.res)
plot(density(a3.m[,"fatal[27]"]),main="Fatal in 2002", col="red", lwd=2)
abline( v=14,col="black")
a3.m <- as.matrix(a3.res)
plot(density(a3.m[,"fatal[28]"]),main="Fatal in 2003", col="red", lwd=2)
abline( v=7,col="black")
a3.m <- as.matrix(a3.res)
plot(density(a3.m[,"fatal[29]"]),main="Fatal in 2004", col="red", lwd=2)
abline( v=9,col="black")</pre>
```



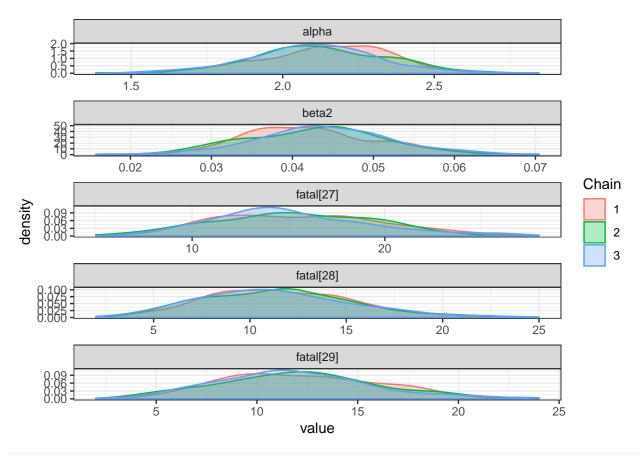
Forth model

```
library(rjags)
 cat( "model
for( i in 1:I )
 fatal[i] ~ dpois( mu[i] )
 log(mu[i]) <- alpha + beta2 * miles[i]*rate[i]</pre>
 alpha ~ dnorm(0,0.0001)
 beta2 ~ dnorm(0.0001,0.00001)
 }",
file="a4.jag" )
 a4.ini= list(
    list(alpha = 0.01, beta2=0.1),
    list(alpha = 0.2, beta2=0.2),
    list(alpha = 0.7, beta2=0.3))
 a4.dat <- list( rate= c(airline$rate,0.7080,0.3,0.4),fatal=c(airline$fatal,NA,NA,NA), miles=c(airline$
 a4.par <- c("alpha", "beta2", "fatal[27]", "fatal[28]", "fatal[29]")
 # Model compilation and burn-in
 a4.mod <- jags.model(file = "a4.jag",data = a4.dat,inits = a4.ini,n.chains = 3,n.adapt = 1000)
```

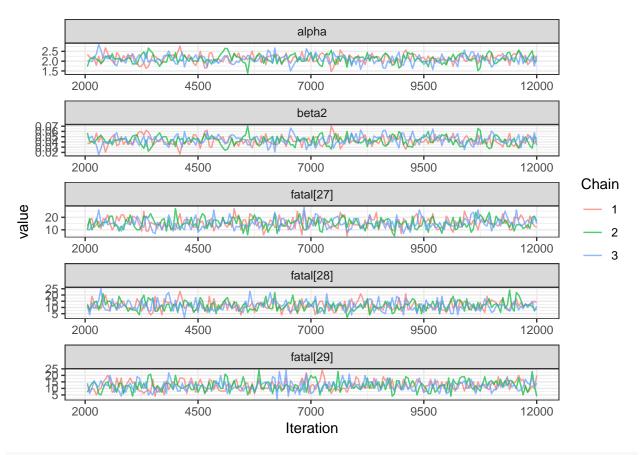
```
## Compiling model graph
##
      Resolving undeclared variables
##
      Allocating nodes
## Graph information:
##
      Observed stochastic nodes: 26
##
      Unobserved stochastic nodes: 5
##
      Total graph size: 180
##
## Initializing model
update(a4.mod, 1000)
 # Sampling from the posterior
a4.res <- coda.samples( a4.mod, var = a4.par, n.iter = 10000, thin = 50 )
summary( a4.res )
##
## Iterations = 2050:12000
## Thinning interval = 50
## Number of chains = 3
## Sample size per chain = 200
## 1. Empirical mean and standard deviation for each variable,
     plus standard error of the mean:
##
##
                            SD Naive SE Time-series SE
                 Mean
              2.12438 0.216094 0.0088220
                                              0.0106781
## alpha
## beta2
              0.04294 0.008454 0.0003451
                                              0.0004023
## fatal[27] 15.36667 4.179873 0.1706426
                                              0.1845807
## fatal[28] 11.45667 3.872525 0.1580952
                                              0.1583297
## fatal[29] 12.07667 3.890503 0.1588291
                                              0.1589362
##
## 2. Quantiles for each variable:
##
               2.5%
                         25%
                                  50%
                                           75%
                                                  97.5%
## alpha
             1.6888 1.99182 2.12813 2.27721 2.52951
## beta2
             0.0272 0.03714 0.04288 0.04846 0.06034
## fatal[27] 8.0000 12.00000 15.00000 18.00000 24.00000
## fatal[28] 5.0000 9.00000 11.00000 14.00000 20.00000
## fatal[29] 5.0000 9.00000 12.00000 15.00000 20.00000
```

Model 4 diagnostics

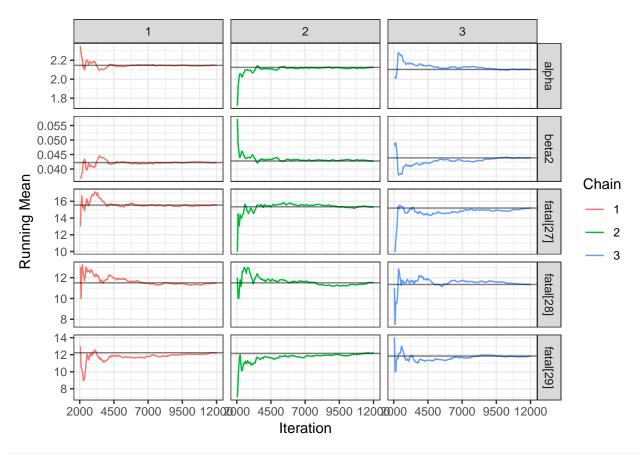
```
result4 = ggs(a4.res)
ggs_density(result4)
```



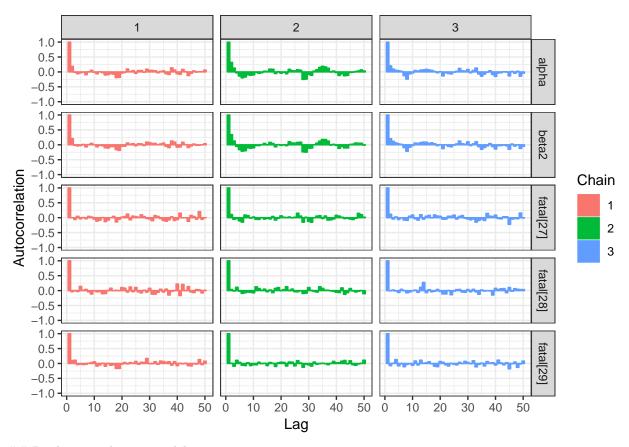
ggs_traceplot(result4)



ggs_running(result4)

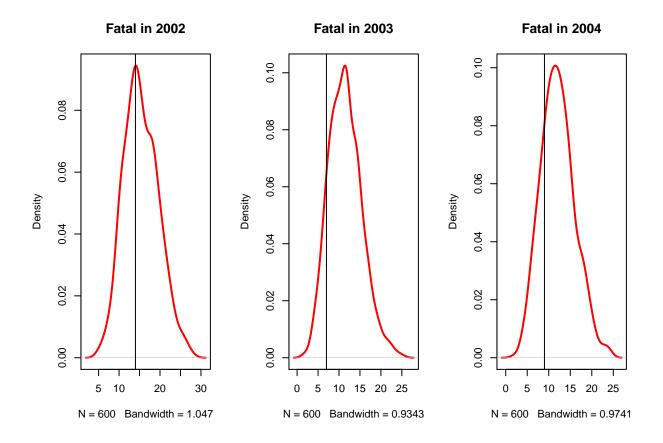


ggs_autocorrelation(result4)



Predicting values in model 4

```
par(mfrow=c(1,3))
a4.m <- as.matrix(a4.res)
plot(density(a4.m[,"fatal[27]"]),main="Fatal in 2002", col="red", lwd=2)
abline( v=14,col="black")
a4.m <- as.matrix(a4.res)
plot(density(a4.m[,"fatal[28]"]),main="Fatal in 2003", col="red", lwd=2)
abline( v=7,col="black")
a4.m <- as.matrix(a4.res)
plot(density(a4.m[,"fatal[29]"]),main="Fatal in 2004", col="red", lwd=2)
abline( v=9,col="black")</pre>
```



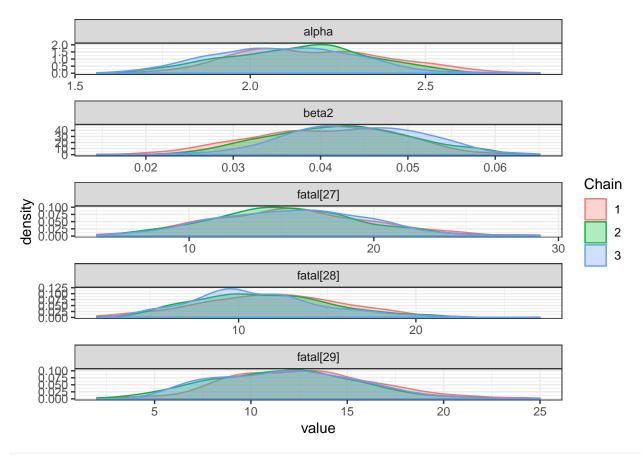
Model 5 test with normal approximation

```
library(rjags)
 cat( "model
for( i in 1:I )
 fatal[i] ~ dnorm( mu[i],1/mu[i] )
 mu[i] <- alpha + beta2 * miles[i]</pre>
 }
 alpha ~ dnorm(0,0.0001)
 beta2 ~ dnorm(0,0.0001)
 file="a5.jag" )
 a5.ini= list(
    list(alpha = 0.01, beta2=3),
    list(alpha = 0.2, beta2=1),
    list(alpha = 0.7, beta2=2))
 a5.dat <- list( rate = c(airline$rate,0.7080,0.3,0.4),fatal=c(airline$fatal,NA,NA,NA), miles=c(airline
 a5.par <- c("alpha", "beta2", "fatal[27]", "fatal[28]", "fatal[29]")
 \# Model compilation and burn-in
 a5.mod <- jags.model(file = "a5.jag",data = a5.dat,inits = a5.ini,n.chains = 3,n.adapt = 1000)
```

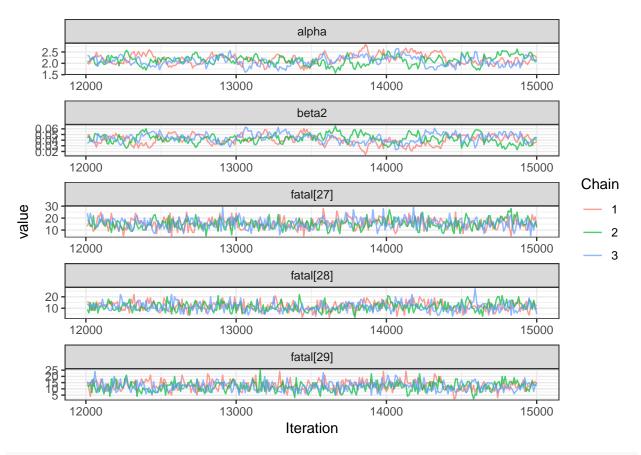
```
## Warning in jags.model(file = "a5.jag", data = a5.dat, inits = a5.ini,
## n.chains = 3, : Unused variable "rate" in data
## Compiling model graph
      Resolving undeclared variables
##
##
      Allocating nodes
## Graph information:
##
      Observed stochastic nodes: 26
##
      Unobserved stochastic nodes: 5
     Total graph size: 151
##
##
## Initializing model
 # Sampling from the posterior
update(a5.mod, 1000)
a5.res <- coda.samples( a4.mod,var = a4.par,n.iter = 3000,thin = 10 )
summary( a5.res )
##
## Iterations = 12010:15000
## Thinning interval = 10
## Number of chains = 3
## Sample size per chain = 300
## 1. Empirical mean and standard deviation for each variable,
     plus standard error of the mean:
##
##
                            SD Naive SE Time-series SE
              2.13946 0.207691 0.0069230
                                              0.0211400
## alpha
## beta2
              0.04228 0.008088 0.0002696
                                              0.0008388
## fatal[27] 15.48333 4.200581 0.1400194
                                              0.1954485
## fatal[28] 11.43000 3.832794 0.1277598
                                              0.1548190
## fatal[29] 12.10444 3.779251 0.1259750
                                              0.1680841
##
## 2. Quantiles for each variable:
##
##
                2.5%
                          25%
                                   50%
                                            75%
                                                   97.5%
             1.75429 1.99413 2.13811 2.28086 2.55063
## alpha
            0.02626 0.03673 0.04242 0.04796 0.05746
## fatal[27] 8.00000 13.00000 15.00000 18.00000 24.00000
## fatal[28] 5.00000 9.00000 11.00000 14.00000 19.52500
## fatal[29] 5.00000 9.00000 12.00000 15.00000 20.00000
```

Model5 diagnostics

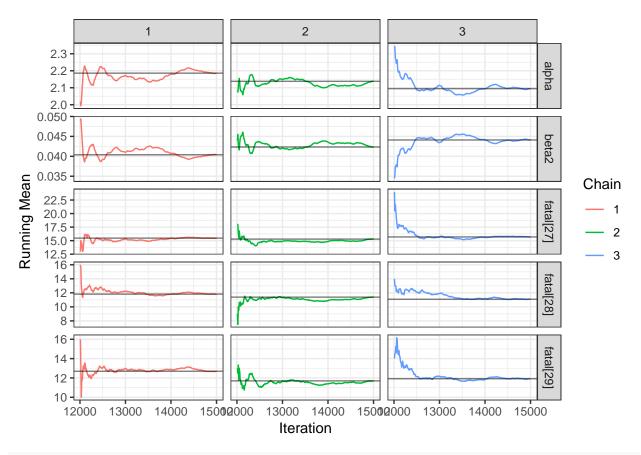
```
result5 = ggs(a5.res)
ggs_density(result5)
```



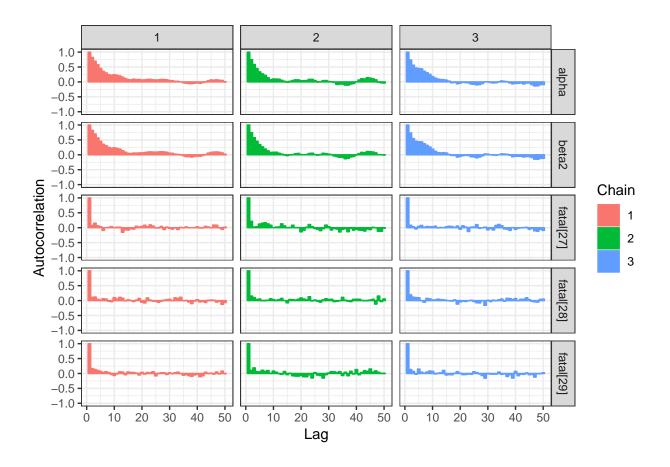
ggs_traceplot(result5)



ggs_running(result5)



ggs_autocorrelation(result5)



Predicting values in model 5

```
par(mfrow=c(1,3))
a5.m <- as.matrix(a5.res)
plot(density(a5.m[,"fatal[27]"]),main="Fatal in 2002", col="red", lwd=2)
abline( v=14,col="black")
a5.m <- as.matrix(a5.res)
plot(density(a5.m[,"fatal[28]"]),main="Fatal in 2003", col="red", lwd=2)
abline( v=7,col="black")
a5.m <- as.matrix(a5.res)
plot(density(a5.m[,"fatal[29]"]),main="Fatal in 2004", col="red", lwd=2)
abline( v=9,col="black")</pre>
```

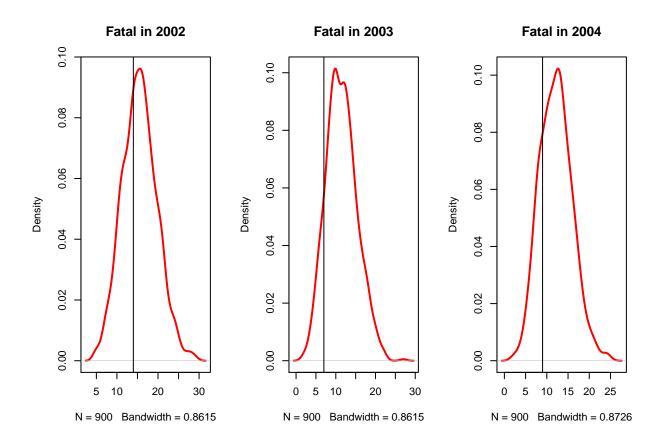


Table of true values

Year	Fatal accidents	Passenger miles	Accident rate
2002	14	19.775	0.7080
2003	7	23.300	0.3004
2004	9	20.300	0.4433

Model comparison throughout DIC index

A good way to select the model that fits best our data, is using the DIC measure. The DIC (deviance information criterion) is a Bayesian criterion for model comparison, defined as:

$$DIC = D(\bar{\theta}) + 2_{p_D}$$

Where: $D(\theta) = -2logL(\theta)$, $p_D = \bar{D} - D(\bar{\theta})$, is the effective number of parameters of the model, the larger p_D the easier the model fits on the data. $\bar{D} = E[D(\theta)]$, its the mean deviance. is a measure of how well the model fits the data, the larger the worse the fit. $D(\bar{\theta}) = D[E(\theta)]$

the basic idea is that the models with smaller DIC should be preferred. Models are penalized both by the value of \bar{D} , which favors a good fit, but also by the effective number of parameters p_D . Since \bar{D} will decrease as the number of parameters in a model increases, the p_D term compensates for this effect by favoring models with a smaller number of parameters.

```
#Model Comparison
DICmodel1 = dic.samples(a1.mod, n.iter=10000, thin=10, type="pD")
DICmodel3 = dic.samples(a3.mod, n.iter=10000, thin =10, type="pD")
DICmodel4 = dic.samples(a4.mod, n.iter=10000, thin=10, type="pD")
DICmodel5 = dic.samples(a5.mod, n.iter=10000, thin =10, type="pD")
DICmodel1
## Mean deviance: 156.2
## penalty 0.9657
## Penalized deviance: 157.2
DICmodel3
## Mean deviance: 133.7
## penalty 2.83
## Penalized deviance: 136.6
DICmodel4
## Mean deviance: 132.9
## penalty 1.918
## Penalized deviance: 134.8
DICmodel5
## Mean deviance: 154.8
## penalty 2.303
## Penalized deviance: 157.1
```