

# R2OpenBUGS: A Package for Running OpenBUGS from R

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## Abstract

The **R2OpenBUGS** package provides convenient functions to call **OpenBUGS** from R. It automatically writes the data and scripts in a format readable by **OpenBUGS** for processing in batch mode, which is possible since version 1.4. After the **OpenBUGS** process has finished, it is possible either to read the resulting data into R by the package itself—which gives a compact graphical summary of inference and convergence diagnostics—or to use the facilities of the **coda** package for further analyses of the output. Examples are given to demonstrate the usage of this package.

*Keywords:* R, OpenBUGS, interface, MCMC.

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An earlier version of this vignette has been published by the Journal of Statistical Software: Sturtz S, Ligges U, Gelman A (2005): “**R2WinBUGS**: A Package for Running WinBUGS from R.” *Journal of Statistical Software*, 12(3), 1–16. **R2OpenBUGS** was adapted from **R2WinBUGS** by Neal Thomas.

## 1. Introduction

The usage of Markov chain Monte Carlo (MCMC) methods became very popular within the last decade. **OpenBUGS** (**B**ayesian inference **U**sing **G**ibbs **S**ampling, Spiegelhalter, Thomas, Best, and Lunn 2003) is a popular software for analyzing complex statistical models using MCMC methods. This software uses Gibbs sampling (Geman and Geman 1984; Gelfand and Smith 1990; Casella and George 1992) and the Metropolis algorithm (Metropolis, Rosenbluth, Rosenbluth, Teller, and Teller 1953) to generate a Markov chain by sampling from full conditional distributions. The **OpenBUGS** software is available for free at <http://www.OpenBUGS.info>. An introduction to MCMC methods is given in Gilks, Richardson, and Spiegelhalter (1996).

Using **OpenBUGS**, the user must specify the model to run, and to load data and initial values for a specified number of Markov chains. Then it is possible to run the Markov chain(s) and to save the results for the parameters the user is interested in. Summary statistics of these data, convergence diagnostics, kernel estimates etc. are available as well. Nevertheless, some users of this software might be interested in saving the output and reading it into R (R Development Core Team 2004) for further analyses. **OpenBUGS** comes with the ability to run the software in batch mode using scripts.

The **R2OpenBUGS** package makes use of this feature and provides the tools to call **OpenBUGS** directly after data manipulation in R. Furthermore, it is possible to work with the results after importing them back into R again, for example to create posterior predictive simulations or, more generally, graphical displays of data and posterior simulations (Gelman 2004). Embedding in R can also be useful for frequently changed data or processing a bunch of data sets, because it is much more convenient to use some R functions (possibly within a loop) rather than using “copy &

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paste” to update data in OpenBUGS each time; however difficulties have been encountered in this area because both R and OpenBUGS can lock up RAM in the Windows operating system.

R is a “language for data analysis and graphics” and an open source and freely available statistical software package implementing that language, see <http://www.R-project.org/>. Historically, R is an implementation of the award-winning S language and system (Becker and Chambers 1984; Becker, Chambers, and Wilks 1988; Chambers and Hastie 1992; Chambers 1998). R and **R2OpenBUGS** are available from CRAN (Comprehensive R Archive Network), i.e., <http://CRAN.R-Project.org> or one of its mirrors. **R2OpenBUGS** could be ported to the commercial S implementation S-PLUS. Minor adaptations would be needed since S-PLUS lacks some of R’s functions and capabilities. If an internet connection is available, **R2OpenBUGS** can be installed by typing `install.packages("R2openBUGS")` at the R command prompt. Do not forget to load the package with `library("R2OpenBUGS")`.

The package **coda** by Plummer, Best, Cowles, and Vines (2004) is very useful for the analysis of OpenBUGS’ output, the reader might want to install this package as well. The CRAN package **boa** (Bayesian Output Analysis Program) by Smith (2004) has similar aims. JAGS (Just Another Gibbs Sampler) by Plummer (2003) is a program for analysis of Bayesian hierarchical models using Gibbs sampling that aims for the same functionality as classic BUGS. JAGS is developed to work closely together with R and the **coda** package.

In this paper, we give two examples, involving educational testing experiments in schools (cf. Section 2.1), and incidence of childhood leukaemia depending on benzene emissions (cf. Section 2.2). Details on the functions of **R2OpenBUGS** are given in Section 3. These functions automatically write the data and a script in a format readable by OpenBUGS for processing in batch mode, and call OpenBUGS from R. After the OpenBUGS process has finished, it is possible either to read the resulting data into R by the package itself or to use the facilities of the **coda** package for further analyses of the output. In Section 4, we demonstrate how to apply the functions provided by **R2OpenBUGS** on the examples’ data, and how to analyze the output both with package **coda** and with **R2OpenBUGS**’s methods to `plot()` and `print()` the output.

## 2. Examples

In this Section, we introduce two examples which will be continued in Section 4.

### 2.1. Schools data

The Scholastic Aptitude Test (SAT) measures the aptitude of high-schoolers in order to help colleges to make admissions decisions. It is divided into two parts, verbal (SAT-V) and mathematical (SAT-M). Our data comes from the SAT-V (Scholastic Aptitude Test-Verbal) on eight different high schools, from an experiment conducted in the late 1970s. SAT-V is a standard multiple choice test administered by the Educational Testing Service. This Service was interested in the effects of coaching programs for each of the selected schools.

The study included coached and uncoached pupils, about sixty in each of the eight different schools; see Rubin (1981). All of them had already taken the PSAT (Preliminary SAT) which results were used as covariates. For each school, the estimated treatment effect and the standard error of the effect estimate are given. These are calculated by an analysis of covariance adjustment appropriate for a completely randomized experiment (Rubin 1981). This example was analyzed using a hierarchical normal model in Rubin (1981) and Gelman, Carlin, Stern, and Rubin (2003, Section 5.5).

### 2.2. Leukaemia registration data

Spatial data usually arises on different, non-nesting spatial scales. One example is childhood leukaemia registration data analyzed by Best, Cocksings, Bennett, Wakefield, and Elliott (2001) using ecologic regression. Data are given for Greater London bounded by the M25 orbital motorway. The data are not available as an example in **R2OpenBUGS** but we use the example here to illustrate alternative calls to the `bugs()` function and output analysis using the **coda** package.

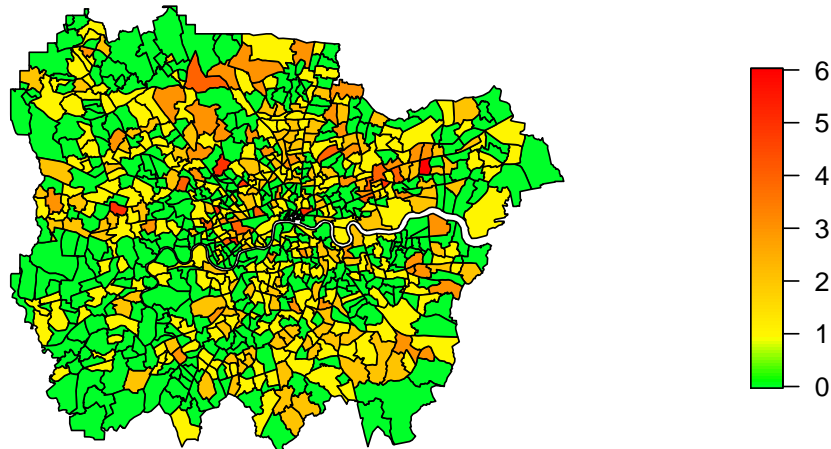


Figure 1: Observed number of cases of childhood leukaemia in 1985–1996

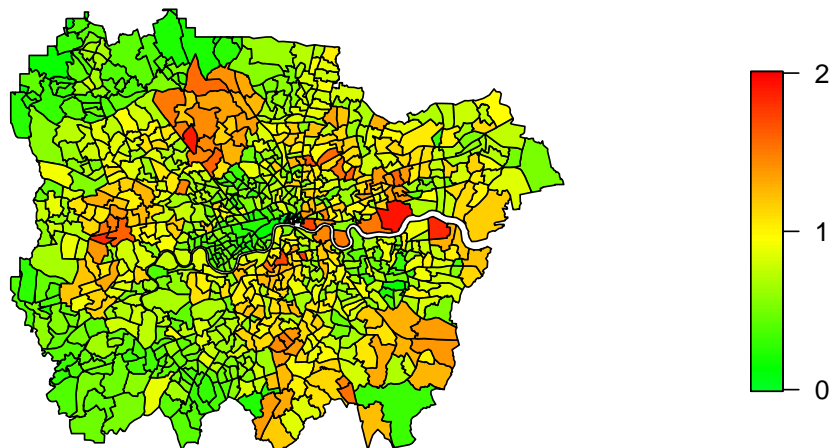


Figure 2: Expected number of cases of childhood leukaemia in 1985–1996

The observed number of leukaemia cases among children under 15 years old is given at ward level. Census wards are administrative areas containing approximately 5000 to 10000 people. Central London is divided into 873 wards. The number of incident cases of leukaemia in children is available from 1985 until 1996 from the Office of National Statistics and the Thames Cancer Registry. A plot of these numbers is given in Figure 1.

Additionally, the number of expected cases (cf. Fig. 2) is calculated on the same resolution using population numbers for different age-sex-strata and the national leukaemia rate for the corresponding strata, for details see [Best \*et al.\* \(2001\)](#).

It is assumed that benzene emissions have an effect on the incidence rate of leukaemia. Benzene emission rates are available in tonnes per year from an atmospheric emissions inventory for London ([Buckingham, Clewley, Hutchinson, Sadler, and Shah 1997](#)) produced by the London Research Centre. They are provided at  $1\text{km} \times 1\text{km}$  grid cells, giving 2132 grid cells in total. Their spatial distribution is shown in Figure 3.

For further details on the data see [Best \*et al.\* \(2001\)](#).

We model these data by Poisson-Gamma models introduced by [Best, Ickstadt, and Wolpert \(2000\)](#) using OpenBUGS. A linking matrix containing information which grid cell belongs to which ward

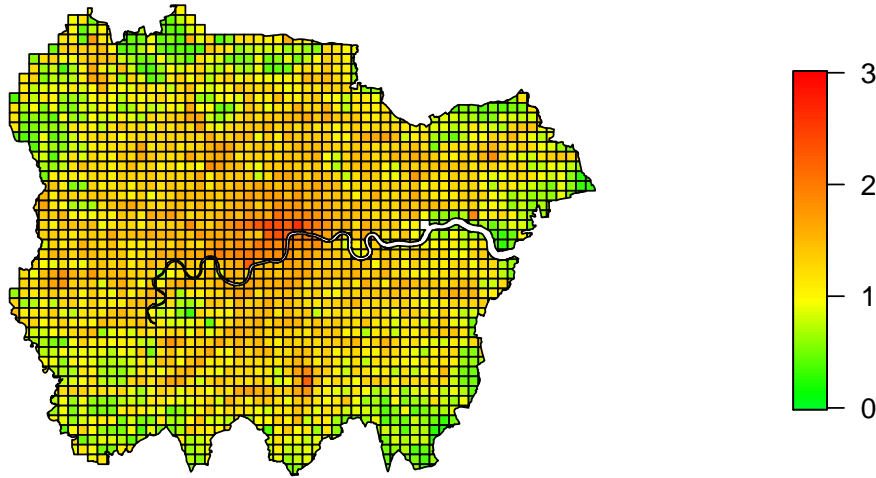


Figure 3: Benzene emissions in tonnes per year

and to which amount is required. This matrix is calculated using R. Unfortunately, **OpenBUGS** does not support a list format such as directly produced by R. Therefore, the data must be provided as a matrix with 2132 rows and 873 columns (or vice versa). Most of the entries of this matrix are zeroes, but using `dump()` to export it from R yields in a file size of 14.2 MB. Unfortunately, opening a file of such size really slows **OpenBUGS** down, and it was not even possible on some of our PCs. Importing data written by our **R2OpenBUGS** package does not make any problems using the batch mode, probably due to memory management issues in **OpenBUGS**.

### 3. Implementation

The implementation of the **R2OpenBUGS** package is straightforward. The “main” function `bugs()` is intended to be called by the user. In principle, it is a wrapper for several other functions called therein step by step as follows:

1. `bugs.data.inits()` writes the data files ‘data.txt’, and ‘inits1.txt’, ‘inits2.txt’, ... into the working directory. These files will be used by **OpenBUGS** during batch processing.

In particular, input for **OpenBUGS** must not exceed a certain number of digits. Moreover, it needs an **E** instead of an **e** in scientific notation. Scientific notation is particularly desirable because of the “number of digits” limitation. The default (`digits = 5`) is to, e.g., reformat the number 123456.789 to 1.23457E+05.

2. `bugs.script()` writes the file ‘script.txt’ that is used by **OpenBUGS** for batch processing.
3. `bugs.run()` updates the lengths of the adaptive phases in the WinBUGS registry (using a function `bugs.update.settings()`), calls WinBUGS, and runs it in batch mode with ‘script.txt’.
4. `bugs.sims()` is only called if the argument `codaPkg` has been set to **FALSE** (the default). Otherwise `bugs()` returns the filenames of stored data. These can, for example, be imported by package **coda** (see the example in Section 4.2, page 10), which provides functions for convergence diagnostics, calculation of Monte Carlo estimates, trace plots, and so forth.

The function `bugs.sims()` reads simulations from **OpenBUGS** into R (not necessarily called by `bugs()` itself), formats them, monitors convergence, performs convergence checks, and computes medians and quantiles. It also prepares the output for `bugs()` itself.

These functions are not intended to be called by the user directly. Arguments are passed from `bugs()` to the other functions, if appropriate. A shortened help file of `bugs()` listing all arguments is given in Appendix A; for the full version type `?bugs` in R after having installed and loaded the package **R2OpenBUGS** (see Section 1).

As known from **OpenBUGS**, one must specify the `data` in form of a list, with list names equal to the names of data in the corresponding **OpenBUGS** model. Alternatively, it is possible to specify a vector or list of names (of mode `character`). In that case objects of that names are looked for in the environment in which `bugs()` has been called (usually that is the user's Workspace, `.GlobalEnv`). If data have already been written in a file called `'data.txt'` to the working directory, it is possible to specify `data = "data.txt"`. One will usually want to supply initial values. This can be done either in the form of a function `inits()` that creates these values, so that different chains can be automatically initialized at different points (see Section 4.1), or by specifying them directly (see Section 4.2). If `inits()` is not specified, `bugs()` just uses the starting values created by **OpenBUGS**; but in practice **OpenBUGS** can crash when reasonable initial values are not specified, and so we recommend constructing a simple `inits()` function to simulate reasonable starting points (Gelman *et al.* 2003, Section C.2). It is also necessary to specify which parameters should be saved for monitoring by specifying `parameters.to.save`.

The user might also want to change the defaults for the length of the burn-in (`n.burnin`, which defaults to half the length of the chain) period for every MCMC run and the number of iterations (`n.iter`, default value 3) that are used to calculate Monte Carlo estimates. The specification of a thinning parameter (`n.thin`) is possible as well; this is useful when the number of parameters is large, to keep the saved output to a reasonably-sized R object.

By setting the argument `debug = TRUE`, **OpenBUGS** remains open after the run. This way it is possible to find errors in the code or the data structure, or even to work with that software as in a usual run. This feature is not available with Linux execution.

It is possible to run one or more Markov chains. The number of chains (`n.chains`) must be specified together with the chains' initial values (`inits`). If more than one Markov chain is requested and `codaPkg` is set to `FALSE`, the convergence diagnostic  $\hat{R}$  (Brooks and Gelman 1998) is calculated by `bugs.sims()` for each of the saved parameters.

Since the communication between **OpenBUGS** and R is based on files, rather huge files will be saved in the working directory by the `bugs()` call, either files to be read in by `bugs()` itself, or by the **coda** package. The user might want to delete those files after the desired contents has been imported into R, and save those objects, e.g., as compressed R data files.

The function `bugs()` returns a rather complex object of class `bugs`, if called with argument `codaPkg = FALSE`. In order to look at the structure of such an object, type `str(objectname)`. For convenience, **R2OpenBUGS** provides methods corresponding to class `bugs` for the generic functions `print()` and `plot()`.

So that user will not be overwhelmed with information; summaries of the output are provided by the `print()` method. That is, some parameters of the `bugs()` call are summarized, and mean, standard deviation, several quantiles of the parameters and convergence diagnostics based on Gelman and Rubin (1992) are printed. See the example in Section 4.1, page 7, for a typical output. As with Spiegelhalter, Best, Carlin, and van der Linde (2002), the DIC computed by `bugs.sims()` is defined as the posterior mean of the deviance plus  $p_D$ , the estimated effective number of parameters in the posterior distribution. We define  $p_D$  as half the posterior variance of the deviance and estimate it as half the average of the within-chain variances of the deviance.<sup>1</sup>

The `plot()` for objects of class `bugs` provides information condensed in some plots conveniently arranged within the same graphics device. For an example, see Figure 4 in Section 4.1. It is

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<sup>1</sup>In contrast, Spiegelhalter *et al.* (2002), and **OpenBUGS**, define  $p_D$  as the posterior mean of the deviance evaluated at the posterior mean of the parameter values. We cannot use that definition because the deviance function is not available to our program, which calls **OpenBUGS** from the "outside". Both definitions of  $p_D$ —ours and that introduced by Spiegelhalter *et al.* (2002)—can be derived from the asymptotic  $\chi^2$  distribution of the deviance relative to its minimum (Gelman *et al.* 2003, Section 6.7). We make no claim that our measure of  $p_D$  is superior to that of Spiegelhalter *et al.* (2002); we choose this measure purely because it is computationally possible given what is available to us from the **OpenBUGS** output.

intended to adapt this function to work with MCMC output in general, even if obtained from software other than OpenBUGS.

## 4. Examples continued

The Examples introduced in Section 4 are continued in this Section. We apply the functions provided by **R2OpenBUGS** to the examples' data and analyze the output.

### 4.1. Schools data

Schools example data (see Section 2.1) are available with the **R2OpenBUGS** package:

```
> data(schools)
> schools
  school estimate    sd
1      A    28.39  14.9
2      B     7.94  10.2
3      C    -2.75  16.3
4      D     6.82  11.0
5      E    -0.64   9.4
6      F     0.63  11.4
7      G    18.01  10.4
8      H    12.16  17.6
```

For modeling these data, we use a hierarchical model as proposed by Gelman *et al.* (2003, Section 5.5). We assume a normal distribution for the observed estimate for each school with mean `theta` and inverse-variance `tau.y`. The inverse-variance is given as  $1/\sigma.y^2$  and its prior distribution is uniform on (0,1000). For the mean `theta`, we employ another normal distribution with mean `mu.theta` and inverse-variance `tau.theta`. For their prior distributions, see the following OpenBUGS code:

```
model {
  for (j in 1:J)
  {
    y[j] ~ dnorm (theta[j], tau.y[j])
    theta[j] ~ dnorm (mu.theta, tau.theta)
    tau.y[j] <- pow(sigma.y[j], -2)
  }
  mu.theta ~ dnorm (0.0, 1.0E-6)
  tau.theta <- pow(sigma.theta, -2)
  sigma.theta ~ dunif (0, 1000)
}
```

This model must be stored in a separate file, e.g. 'schools.bug'<sup>2</sup>, in an appropriate directory, say `c:/schools/`. In R the user must prepare the data inputs the `bugs()` function needs. This can be a list containing the name of each data vector, e.g.

```
> J <- nrow(schools)
> y <- schools$estimate
> sigma.y <- schools$sd
> data <- list ("J", "y", "sigma.y")
```

<sup>2</sup>Emacs Speaks Statistics (ESS) by Rossini, Heiberger, Sparapani, Mächler, and Hornik (2004), a package available with Gnu Emacs (Stallmann 1999), recognizes and properly formats Bugs model files that have the .bug extension.



Using these data and the model file, we can run an MCMC simulation to get estimates for `theta`, `mu.theta` and `sigma.theta`. Before running, the user must decide how many chains to be run (`n.chain = 3`) for how many iterations (`n.iter = 1000`). If the length of burn-in is not specified, `n.burnin = floor(n.iter/2)` is used, that is, 500 in this example. Additionally, the user must specify initial values for the chains, for example by writing a function. This can be done by

```
> inits <- function(){
+   list(theta = rnorm(J, 0, 100), mu.theta = rnorm(1, 0, 100),
+        sigma.theta = runif(1, 0, 100))
+ }
```

Now, the user can start the MCMC simulation by typing

```
> schools.sim <- bugs(data, inits, model.file = "c:/schools/schools.bug",
+   parameters = c("theta", "mu.theta", "sigma.theta"),
+   n.chains = 3, n.iter = 1000)
```

in R. For other available arguments, see Appendix A.

The results in objects `schools.sim` can conveniently be printed by `print(schools.sim)`. The generic function `print()` calls the print method for an object of class `bugs` provided by **R2OpenBUGS**. For this example, you will get something like

```
> print(schools.sim)
Inference for Bugs model at "c:/schools/schools.bug"
3 chains, each with 1000 iterations (first 500 discarded)
n.sims = 1500 iterations saved
```

	mean	sd	2.5%	25%	50%	75%	97.5%	Rhat	n.eff
theta[1]	11.1	9.1	-3.0	5.0	10.0	16.0	31.8	1.1	39
theta[2]	7.6	6.6	-4.7	3.3	7.8	11.6	21.1	1.1	42
theta[3]	5.7	8.4	-12.5	0.6	6.1	10.8	21.8	1.0	150
theta[4]	7.1	7.0	-6.6	2.7	7.2	11.5	21.0	1.1	42
theta[5]	5.1	6.8	-9.5	0.7	5.2	9.7	18.1	1.0	83
theta[6]	5.7	7.3	-9.7	1.0	6.2	10.2	20.0	1.0	56
theta[7]	10.4	7.3	-2.1	5.3	9.8	15.3	25.5	1.1	27
theta[8]	8.3	8.4	-6.6	2.8	8.1	12.7	26.2	1.0	64
mu.theta	7.6	5.9	-3.0	3.7	8.0	11.0	19.5	1.1	35
sigma.theta	6.7	5.6	0.3	2.8	5.1	9.2	21.2	1.1	46
deviance	60.8	2.5	57.0	59.1	60.2	62.1	66.6	1.0	170

```
pD = 3 and DIC = 63.8 (using the rule, pD = var(deviance)/2)
```

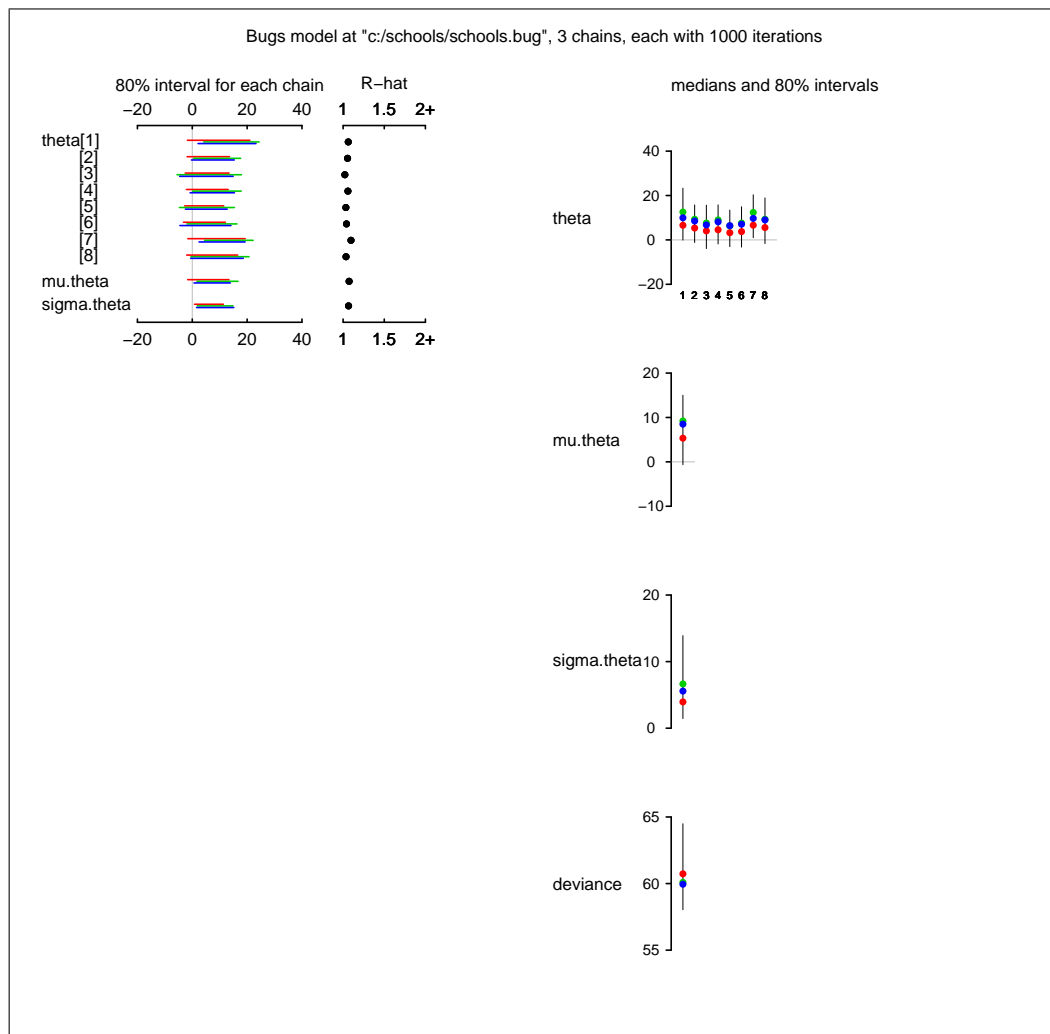
For each parameter, `n.eff` is a crude measure of effective sample size, and `Rhat` is the potential scale reduction factor (at convergence, `Rhat=1`). `DIC` is an estimate of expected predictive error (lower deviance is better).

Additionally, the user can generate a plot of the results by typing `plot(schools.sim)`. The resulting plot is given in Figure 4. In this plot, the left column shows a quick summary of inference and convergence ( $\hat{R}$  is close to 1.0 for all parameters, indicating good mixing of the three chains and thus approximate convergence); and the right column shows inferences for each set of parameters. As can be seen in the right column, **R2OpenBUGS** uses the parameter names in **OpenBUGS** to structure the output into scalar, vector, and arrays of parameters, in addition to storing the parameters as a long vector.

For the interpretation of these results see Gelman *et al.* (2003, Section 5.5).

## 4.2. Leukaemia registration data

The leukaemia registration data (see Section 2.2) are used to show data modeling and output reading into R using the **coda** package. A simple model for these data looks as follows:

Figure 4: Plot produced by **R2OpenBUGS** package for the schools example.

```
model{
  beta.0 ~ dgamma(a.0, tau.0)
  beta.benz ~ dgamma(a.benz, tau.benz)
  a.0 <- 0.575
  tau.0 <- a.0*2
  a.benz <- 0.575
  tau.benz <- a.benz*2
}
```



```

for (i in 1:I)
{
  count[i] ~ dpois(lambda[i])
  lambda[i] <- p[i]*expect[i]
  for (j in 1:J)
  {
    prop[j,i] <- gamma[j,i]*(benz[j] - benzbar)
  }
  p[i]<- beta.0 + beta.benz*sum(prop[,i])
}
}

```

Here `count` denotes the number of observed incidences of childhood leukaemia in ward `i`. These are assumed to be Poisson distributed with mean `lambda` depending on the number of expected cases `expect` in ward `i` and an area-specific risk rate `p`. For calculation of this area specific risk rate we use an intercept `beta.0` and a term depending on the weighted sum of benzene emissions `benz` in each grid cell `j`. The weights are chosen proportional to the amount of area that ward `i` and grid cell `j` have in common.

In R we can define all these data and then initialize the model. The data needed for this example are

`benzbar`: arithmetic mean of all benzene values,

`benz`: a vector containing benzene emissions of all 2132 grid cells,

`expect`: expected number of cases of childhood leukaemia in each of the 873 wards,

`count`: observed number of childhood leukaemia in these wards,

`gamma`: a  $2132 \times 873$  matrix containing the amount of area each grid cell and each ward have in common,

`J`: total number of grid cells, i.e. 2132, and

`I`: total number of ward cells, i.e. 873.

The parameters we want to store are regression coefficients `beta.0` and `beta.benz` as well as `p`, the area specific relative risk compared to the reference rate. This reference rate was used to calculate the expected number of cases in each ward.

Since we want to use the `coda` package for reading the data into OpenBUGS, we specify `codaPkg = TRUE` in the `bugs()` call:

```

> data <- list(benzbar = mean(benz), benz = benz, expect = expect,
+   count = count, gamma = gamma, J = J, I = I)
> parameters <- c("beta.0", "beta.benz", "p")
> inits1 <- list(beta.0 = 1, beta.benz = 1)
> inits2 <- list(beta.0 = 0.5, beta.benz = 0.5)
> inits <- list(inits1, inits2)
> model <- bugs(data, inits, parameters, model.file = "c:/model.bug",
+   n.chains = 2, n.iter = 8000, n.burnin = 5000, n.thin = 1,
+   codaPkg = TRUE )

```

Starting with, e.g.,

```
> library("coda")
> codaobject <- read.bugs(model)
> plot(codaobject)
```

it is now possible to use the **coda** package for output analyses.

## Acknowledgments

The work of Uwe Ligges has been supported by the Deutsche Forschungsgemeinschaft, Sonderforschungsbereich 475. The work of Andrew Gelman has been supported by the U.S. National Science Foundation.

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## A. Help page for the function bugs()

This help page has been shortened.

---

bugs

*Run OpenBUGS from R*

---

### Description

The **bugs** function takes data and starting values as input. It automatically writes a **OpenBUGS** script, calls the model, and saves the simulations for easy access in R.

### Usage

```
bugs(data, inits, parameters.to.save, n.iter, model.file="model.txt",
      n.chains=3, n.burnin=floor(n.iter / 2), n.thin=1,
      saveExec=FALSE, restart=FALSE,
      debug=FALSE, DIC=TRUE, digits=5, codaPkg=FALSE,
      OpenBUGS.pgm=NULL, working.directory=NULL,
      clearWD=FALSE, useWINE=FALSE, WINE=NULL,
      newWINE=TRUE, WINEPATH=NULL, bugs.seed=1, summary.only=FALSE,
      save.history=(.Platform$OS.type == "windows" | useWINE==TRUE),
      over.relax = FALSE)
```

### Arguments

<b>data</b>	either a named list (names corresponding to variable names in the <code>model.file</code> ) of the data for the <b>OpenBUGS</b> model, <i>or</i> a vector or list of the names of the data objects used by the model. If <b>data</b> is a one element character vector (such as <code>"data.txt"</code> ), it is assumed that data have already been written to the working directory into that file, e.g. by the function <code>bugs.data</code> .
<b>inits</b>	a list with <code>n.chains</code> elements; each element of the list is itself a list of starting values for the <b>OpenBUGS</b> model, <i>or</i> a function creating (possibly random) initial values. Alternatively, if <code>inits=NULL</code> , initial values are generated by <b>OpenBUGS</b> . If <b>inits</b> is a character vector with <code>n.chains</code> elements, it is assumed that inits have already been written to the working directory into those files, e.g. by the function <code>bugs.inits</code> .
<b>parameters.to.save</b>	character vector of the names of the parameters to save which should be monitored
<b>model.file</b>	File containing the model written in <b>OpenBUGS</b> code. The extension must be <code>‘.txt’</code> . The old convention allowing <code>model.file</code> to be named <code>‘.bug’</code> has been eliminated because the new <b>OpenBUGS</b> feature that allows the program image to be saved and later restarted uses the <code>.bug</code> extension for the saved images. Alternatively, <code>model.file</code> can be an R function that contains a BUGS model that is written to a temporary model file (see <code>tempfile</code> ) using <code>write.model</code> .
<b>n.chains</b>	number of Markov chains (default: 3)
<b>n.iter</b>	number of total iterations per chain (including burn in; default: 2000)
<b>n.burnin</b>	length of burn in, i.e. number of iterations to discard at the beginning. Default is <code>n.iter/2</code> , that is, discarding the first half of the simulations.
<b>n.thin</b>	Thinning rate. Must be a positive integer. The default is <code>n.thin = 1</code> . The thinning is implemented in the OpenBUGS update phase, so thinned samples are never stored, and they are not counted in <code>n.burnin</code> or <code>n.iter</code> . Setting <code>n.thin=2</code> , doubles the number of iterations OpenBUGS performs, but does not change <code>n.iter</code> or <code>n.burnin</code> . Thinning implemented in this manner is not captured in summaries created by packages such as <b>coda</b> .

<code>saveExec</code>	If TRUE, a re-startable image of the OpenBUGS execution is saved with <code>basename(model.file)</code> and extension <code>.bug</code> in the working directory, which must be specified. The <code>.bug</code> files can be large, so users should monitor them carefully and remove them when not needed.
<code>restart</code>	If TRUE, execution resumes with the final status from the previous execution stored in the <code>.bug</code> file in the working directory. If <code>n.burnin=0</code> , additional iterations are performed and all iterations since the previous burnin are used (including those from past executions). If <code>n.burnin&gt;0</code> , a new burnin is performed, and the previous iterations are discarded, but execution continues from the status at the end of the previous execution. When <code>restart=TRUE</code> , only <code>n.burnin</code> , <code>n.iter</code> , and <code>saveExec</code> inputs should be changed from the call creating the <code>.bug</code> file, otherwise failed or erratic results may be produced. Note the default has <code>n.burnin&gt;0</code> .
<code>debug</code>	if FALSE (default), <b>OpenBUGS</b> is closed automatically when the script has finished running, otherwise <b>OpenBUGS</b> remains open for further investigation. The debug option is not available for linux execution.
<code>DIC</code>	logical; if TRUE (default), compute deviance, pD, and DIC. This is done in <b>OpenBUGS</b> directly using the rule $pD = \bar{D} - \hat{D}$ . If there are less iterations than required for the adaptive phase, the rule $pD = \text{var}(\text{deviance}) / 2$ is used.
<code>digits</code>	number of significant digits used for <b>OpenBUGS</b> input, see <code>formatC</code>
<code>codaPkg</code>	logical; if FALSE (default) a <code>bugs</code> object is returned, if TRUE file names of <b>OpenBUGS</b> output are returned for easy access by the <code>coda</code> package through function <code>read.bugs</code> . A <code>bugs</code> object can be converted to an <code>mcmc.list</code> object as used by the <code>coda</code> package with the method <code>as.mcmc.list</code> (for which a method is provided by <code>R2OpenBUGS</code> ).
<code>OpenBUGS.pgm</code>	For Windows or WINE execution, the full path to the OpenBUGS executable. For linux execution, the full path to the OpenBUGS shell script (not required if OpenBUGS is in the user's PATH variable). If NULL (unset) and the environment variable <code>OpenBUGS_PATH</code> is set the latter will be used as the default. If NULL (unset), the environment variable <code>OpenBUGS_PATH</code> is unset and the global option <code>R2OpenBUGS.pgm</code> is not NULL the latter will be used as the default. If nothing of the former is set and OS is Windows, the most recent OpenBUGS version registered in the Windows registry will be used as the default.
<code>working.directory</code>	sets working directory during execution of this function; <b>OpenBUGS</b> ' in- and output will be stored in this directory; if NULL, a temporary working directory via <code>tempdir</code> is used.
<code>clearWD</code>	logical; indicating whether the files <code>'data.txt'</code> , <code>'inits[1:n.chains].txt'</code> , <code>'log.odc'</code> , <code>'codaIndex.txt'</code> , and <code>'coda[1:nchains].txt'</code> should be removed after <b>OpenBUGS</b> has finished. If set to TRUE, this argument is only respected if <code>codaPkg=FALSE</code> .
<code>useWINE</code>	logical; attempt to use the Wine emulator to run <b>OpenBUGS</b> . Default is FALSE. If WINE is used, the arguments <code>OpenBUGS.pgm</code> and <code>working.directory</code> must be given in form of Linux paths rather than Windows paths (if not NULL).
<code>WINE</code>	Character, path to <code>'wine'</code> binary file, it is tried hard (by a guess and the utilities <code>which</code> and <code>locate</code> ) to get the information automatically if not given.
<code>newWINE</code>	Use new versions of Wine that have <code>'winepath'</code> utility
<code>WINEPATH</code>	Character, path to <code>'winepath'</code> binary file, it is tried hard (by a guess and the utilities <code>which</code> and <code>locate</code> ) to get the information automatically if not given.
<code>bugs.seed</code>	Random seed for <b>OpenBUGS</b> . Must be an integer between 1-14. Seed specification changed between WinBUGS and OpenBUGS; see the OpenBUGS documentation for details.
<code>summary.only</code>	If TRUE, only a parameter summary for very quick analyses is given, temporary created files are not removed in that case.
<code>save.history</code>	If TRUE (the default), trace plots are generated at the end.
<code>over.relax</code>	If TRUE, over-relaxed form of MCMC is used if available from OpenBUGS.

## Details

To run:

1. Write a **BUGS** model in an ASCII file (hint: use `write.model`).
2. Go into R.
3. Prepare the inputs for the `bugs` function and run it (see Example section).
4. An **OpenBUGS** window will pop up and R will freeze up. The model will now run in **OpenBUGS**. It might take awhile. You will see things happening in the Log window within **OpenBUGS**. When **OpenBUGS** is done, its window will close and R will work again.
5. If an error message appears, re-run with `debug=TRUE`.

BUGS version support:

- **OpenBUGS**  $\geq 3.2.1$

Operation system support:

- **MS Windows** no problem
- **Linux, intel processors** GUI display and graphics not available.
- **Mac OS X** and **Unix** in general possible with Wine emulation via `useWINE=TRUE`

If `useWINE=TRUE` is used, all paths (such as `working.directory` and `model.file`, must be given in native (Unix) style, but `OpenBUGS.pgm` can be given in Windows path style (e.g. “c:/Program Files/OpenBUGS/”) or native (Unix) style (e.g. “/path/to/wine/folder/dosdevices/c:/Program Files/OpenBUGS/OpenBUGS/”).

## Value

If `codaPkg=TRUE` the returned values are the names of coda output files written by **OpenBUGS** containing the Markov Chain Monte Carlo output in the CODA format. This is useful for direct access with `read.bugs`.

If `codaPkg=FALSE`, the following values are returned:

<code>n.chains</code>	see Section ‘Arguments’
<code>n.iter</code>	see Section ‘Arguments’
<code>n.burnin</code>	see Section ‘Arguments’
<code>n.thin</code>	see Section ‘Arguments’
<code>n.keep</code>	number of iterations kept per chain (equal to $(n.iter - n.burnin) / n.thin$ )
<code>n.sims</code>	number of posterior simulations (equal to <code>n.chains * n.keep</code> )
<code>sims.array</code>	3-way array of simulation output, with dimensions <code>n.keep</code> , <code>n.chains</code> , and length of combined parameter vector
<code>sims.list</code>	list of simulated parameters: for each scalar parameter, a vector of length <code>n.sims</code> for each vector parameter, a 2-way array of simulations, for each matrix parameter, a 3-way array of simulations, etc. (for convenience, the <code>n.keep*n.chains</code> simulations in <code>sims.matrix</code> and <code>sims.list</code> (but NOT <code>sims.array</code> ) have been randomly permuted)
<code>sims.matrix</code>	matrix of simulation output, with <code>n.chains*n.keep</code> rows and one column for each element of each saved parameter (for convenience, the <code>n.keep*n.chains</code> simulations in <code>sims.matrix</code> and <code>sims.list</code> (but NOT <code>sims.array</code> ) have been randomly permuted)
<code>summary</code>	summary statistics and convergence information for each saved parameter.
<code>mean</code>	a list of the estimated parameter means
<code>sd</code>	a list of the estimated parameter standard deviations
<code>median</code>	a list of the estimated parameter medians
<code>root.short</code>	names of argument <code>parameters.to.save</code> and “deviance”
<code>long.short</code>	indexes; programming stuff
<code>dimension.short</code>	dimension of <code>indexes.short</code>

<code>indexes.short</code>	indexes of <code>root.short</code>
<code>last.values</code>	list of simulations from the most recent iteration; they can be used as starting points if you wish to run <b>OpenBUGS</b> for further iterations
<code>pD</code>	an estimate of the effective number of parameters, for calculations see the section “Arguments”.
<code>DIC</code>	<code>mean(deviance) + pD</code>

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### References

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### See Also

`print.bugs`, `plot.bugs`, as well as **coda** and **BRugs** packages

### Examples

```
# An example model file is given in:
model.file <- system.file(package="R2OpenBUGS", "model", "schools.txt")
# Let's take a look:
file.show(model.file)

# Some example data (see ?schools for details):
data(schools)
schools

J <- nrow(schools)
y <- schools$estimate
sigma.y <- schools$sd
data <- list ("J", "y", "sigma.y")
inits <- function(){
  list(theta=rnorm(J, 0, 100), mu.theta=rnorm(1, 0, 100),
        sigma.theta=runif(1, 0, 100))
}
## or alternatively something like:
# inits <- list(
#   list(theta=rnorm(J, 0, 90), mu.theta=rnorm(1, 0, 90),
#         sigma.theta=runif(1, 0, 90)),
#   list(theta=rnorm(J, 0, 100), mu.theta=rnorm(1, 0, 100),
#         sigma.theta=runif(1, 0, 100)),
#   list(theta=rnorm(J, 0, 110), mu.theta=rnorm(1, 0, 110),
#         sigma.theta=runif(1, 0, 110)))

parameters <- c("theta", "mu.theta", "sigma.theta")

## Not run:
## You may need to specify "OpenBUGS.pgm"
```



```
## also you need write access in the working directory:
schools.sim <- bugs(data, inits, parameters, model.file,
  n.chains=3, n.iter=5000)
print(schools.sim)
plot(schools.sim)

## End(Not run)
```