bayes: glm — Bayesian generalized linear models

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Description

bayes: glm fits a Bayesian generalized linear model to outcomes of different types such as continuous, binary, count, and so on; see [BAYES] bayes and [R] glm for details.

Quick start

Bayesian generalized linear model of y on x1 and x2, using the Gaussian family and log link and using default normal priors for regression coefficients

```
bayes: glm y x1 x2, family(gaussian) link(log)
```

Use a standard deviation of 10 instead of 100 for the default normal priors

```
bayes, normalprior(10): glm y x1 x2, family(gaussian) link(log)
```

Use uniform priors for the slopes and a normal prior for the intercept

```
bayes, prior({y: x1 x2}, uniform(-10,10)) ///
prior({y:_cons}, normal(0,10)): ///
glm y x1 x2, family(gaussian) link(log)
```

Save simulation results to simdata.dta, and use a random-number seed for reproducibility

```
bayes, saving(simdata) rseed(123): ///
glm y x1 x2, family(gaussian) link(log)
```

Specify 20,000 Markov chain Monte Carlo (MCMC) samples, set length of the burn-in period to 5,000, and request that a dot be displayed every 500 simulations

```
bayes, mcmcsize(20000) burnin(5000) dots(500): ///
glm y x1 x2, family(gaussian) link(log)
```

In the above, request that the 90% highest posterior density (HPD) credible interval be displayed instead of the default 95% equal-tailed credible interval

```
bayes, clevel (90) hpd
```

Fit a logit model and display results as odds ratios

```
bayes: glm z x1 x2, family(binomial) eform
```

Display odds ratios on replay

```
bayes, eform
```

Also see Quick start in [BAYES] bayes and Quick start in [R] glm.

Menu

Statistics > Generalized linear models > Bayesian generalized linear models (GLM)

Syntax

```
\texttt{bayes} \ [ \ \textit{, bayesopts} \ ] : \texttt{glm} \ \textit{depvar} \ [ \ \textit{indepvars} \ ] \ [ \textit{if} \ ] \ [ \ \textit{in} \ ] \ [ \ \textit{weight} \ ] \ [ \ \textit{, options} \ ]
```

	Description		
Model <u>family(familyname)</u>	distribution of <i>depvar</i> ; default is family(gaussian) link function; default is canonical link for family() specified		
<u>l</u> ink(<i>linkname</i>)			
Model 2			
<u>nocons</u> tant	suppress constant term		
exposure(varname)	include ln(<i>varname</i>) in model with coefficient constrained to 1 include <i>varname</i> in model with coefficient constrained to 1		
<pre>offset(varname) asis</pre>	retain perfect predictor variables		
mu(varname)	use <i>varname</i> as the initial estimate for the mean of <i>depvar</i>		
<u>ini</u> t(varname)	synonym for mu(varname)		
Reporting			
<u>ef</u> orm	report exponentiated coefficients		
display_options	control spacing, line width, and base and empty cells		
<u>le</u> vel(#)	set credible level; default is level(95)		
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For a detailed description of option bayesopts	ons, see Options in [R] glm. Description		
bayesopts			
bayesopts Priors	Description specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) prior for model parameters; this option may be repeated		
bayesopts Priors * normalprior(#)	Description specify standard deviation of default normal priors for regression coefficients; default is normalprior(100)		
bayesopts Priors * normalprior(#) prior(priorspec) dryrun Simulation	Specify standard deviation of default normal priors for regression coefficients; default is normalprior (100) prior for model parameters; this option may be repeated show model summary without estimation		
bayesopts Priors * normalprior(#) prior(priorspec) dryrun Simulation nchains(#)	specify standard deviation of default normal priors for regression coefficients; default is normalprior (100) prior for model parameters; this option may be repeated show model summary without estimation number of chains; default is to simulate one chain		
bayesopts Priors * normalprior(#) prior(priorspec) dryrun Simulation nchains(#) mcmcsize(#)	specify standard deviation of default normal priors for regression coefficients; default is normalprior (100) prior for model parameters; this option may be repeated show model summary without estimation number of chains; default is to simulate one chain MCMC sample size; default is mcmcsize (10000)		
bayesopts * normalprior(#) prior(priorspec) dryrun Simulation nchains(#) mcmcsize(#) burnin(#)	specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) prior for model parameters; this option may be repeated show model summary without estimation number of chains; default is to simulate one chain MCMC sample size; default is mcmcsize(10000) burn-in period; default is burnin(2500)		
bayesopts Priors * normalprior(#) prior(priorspec) dryrun Simulation nchains(#) mcmcsize(#)	specify standard deviation of default normal priors for regression coefficients; default is normalprior (100) prior for model parameters; this option may be repeated show model summary without estimation number of chains; default is to simulate one chain MCMC sample size; default is mcmcsize (10000)		
bayesopts *normalprior(#) prior(priorspec) dryrun Simulation nchains(#) mcmcsize(#) burnin(#) thinning(#)	specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) prior for model parameters; this option may be repeated show model summary without estimation number of chains; default is to simulate one chain MCMC sample size; default is mcmcsize(10000) burn-in period; default is burnin(2500) thinning interval; default is thinning(1)		
bayesopts Priors * normalprior(#) prior(priorspec) dryrun Simulation nchains(#) mcmcsize(#) burnin(#) thinning(#) rseed(#) exclude(paramref) Blocking	specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) prior for model parameters; this option may be repeated show model summary without estimation number of chains; default is to simulate one chain MCMC sample size; default is mcmcsize(10000) burn-in period; default is burnin(2500) thinning interval; default is thinning(1) random-number seed specify model parameters to be excluded from the simulation results		
bayesopts * normalprior(#) prior(priorspec) dryrun Simulation nchains(#) mcmcsize(#) burnin(#) thinning(#) rseed(#) exclude(paramref) Blocking * blocksize(#)	specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) prior for model parameters; this option may be repeated show model summary without estimation number of chains; default is to simulate one chain MCMC sample size; default is mcmcsize(10000) burn-in period; default is burnin(2500) thinning interval; default is thinning(1) random-number seed specify model parameters to be excluded from the simulation results maximum block size; default is blocksize(50)		
bayesopts Priors * normalprior(#) prior(priorspec) dryrum Simulation nchains(#) mcmcsize(#) burnin(#) thinning(#) rseed(#) exclude(paramref) Blocking	specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) prior for model parameters; this option may be repeated show model summary without estimation number of chains; default is to simulate one chain MCMC sample size; default is mcmcsize(10000) burn-in period; default is burnin(2500) thinning interval; default is thinning(1) random-number seed specify model parameters to be excluded from the simulation results maximum block size; default is blocksize(50)		

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Initialization	
<pre>init ial(initspec)</pre>	specify initial values for model parameters with a single chain
<pre>init#(initspec)</pre>	specify initial values for #th chain; requires nchains()
<pre>initall(initspec)</pre>	specify initial values for all chains; requires nchains()
<u>nomleinit</u> ial	suppress the use of maximum likelihood estimates as starting values
<u>initrand</u> om	specify random initial values
<u>initsumm</u> ary	display initial values used for simulation
* <u>noi</u> sily	display output from the estimation command during initialization
Adaptation	
adaptation(adaptopts)	control the adaptive MCMC procedure
scale(#)	initial multiplier for scale factor; default is scale (2.38)
\underline{cov} ariance (cov)	initial proposal covariance; default is the identity matrix
Reporting	
<pre>clevel(#)</pre>	set credible interval level; default is clevel(95)
hpd	display HPD credible intervals instead of the default equal-tailed credible intervals
<pre>eform[(string)]</pre>	report exponentiated coefficients and, optionally, label as string
batch(#)	specify length of block for batch-means calculations; default is batch(0)
<pre>saving(filename[, replace])</pre>	save simulation results to filename.dta
nomodelsummary	suppress model summary
chainsdetail	display detailed simulation summary for each chain
[no]dots	suppress dots or display dots every 100 iterations and iteration numbers every 1,000 iterations; default is nodots
$\mathtt{dots}(\#[\ ,\ \mathtt{every}(\#)\])$	display dots as simulation is performed
[no]show(paramref)	specify model parameters to be excluded from or included in the output
<u>notab</u> le	suppress estimation table
<u>nohead</u> er	suppress output header
title(string)	display string as title above the table of parameter estimates
display_options	control spacing, line width, and base and empty cells

Advanced

search(search_options) control the search for feasible initial values

corrlag(#) specify maximum autocorrelation lag; default varies

corrtol(#) specify autocorrelation tolerance; default is corrtol(0.01)

* Starred options are specific to the bayes prefix; other options are common between bayes and bayesmh.

Options prior() and block() may be repeated.

priorspec and paramref are defined in [BAYES] bayesmh.

paramref may contain factor variables; see [U] 11.4.3 Factor variables.

collect is allowed; see [U] 11.1.10 Prefix commands.

See [U] 20 Estimation and postestimation commands for more capabilities of estimation commands.

Model parameters are regression coefficients {depvar: indepvars}. Use the dryrun option to see the definitions of model parameters prior to estimation.

For a detailed description of *bayesopts*, see *Options* in [BAYES] bayes.

Remarks and examples

For a general introduction to Bayesian analysis, see [BAYES] Intro. For a general introduction to Bayesian estimation using an adaptive Metropolis-Hastings algorithm, see [BAYES] bayesmh. For remarks and examples specific to the bayes prefix, see [BAYES] bayes. For details about the estimation command, see [R] glm.

For a simple example of the bayes prefix, see Introductory example in [BAYES] bayes. Also see Generalized linear model in [BAYES] bayes.

bayes: glm does not estimate the scale parameter but uses a fixed value as provided by the glm command. If you want to fit a GLM and estimate the scale parameter, use bayes: meglm without specifying random effects.

Stored results

See Stored results in [BAYES] bayes.

Methods and formulas

See Methods and formulas in [BAYES] bayesmh.

Also see

```
[BAYES] bayes — Bayesian regression models using the bayes prefix
[R] glm — Generalized linear models
[BAYES] Bayesian postestimation — Postestimation tools after Bayesian estimation
[BAYES] Bayesian estimation — Bayesian estimation commands
[BAYES] Bayesian commands — Introduction to commands for Bayesian analysis
[BAYES] Intro — Introduction to Bayesian analysis
[BAYES] Glossary
```

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