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## Description

`bayes: hetprobit` fits a Bayesian heteroskedastic probit regression to a binary outcome; see [\[BAYES\] bayes](#) and [\[R\] hetprobit](#) for details.

## Quick start

Bayesian heteroskedastic probit regression of  $y$  on  $x_1$  and  $x_2$ , using  $z_1$  to model the variance and using default normal priors for regression coefficients and log-variance coefficients

```
bayes: hetprobit y x1 x2, het(z1)
```

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

```
bayes, normalprior(10): hetprobit y x1 x2, het(z1)
```

Use uniform priors for the slopes and a normal prior for the intercept of the main regression

```
bayes, prior({y: x1 x2}, uniform(-10,10)) ///  
prior({y: _cons}, normal(0,10)): hetprobit y x1 x2, het(z1)
```

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

```
bayes, saving(simdata) rseed(123): hetprobit y x1 x2, het(z1)
```

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, mcmcsz(20000) burnin(5000) dots(500): hetprobit y x1 x2, het(z1)
```

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
```

Also see [Quick start](#) in [\[BAYES\] bayes](#) and [Quick start](#) in [\[R\] hetprobit](#).

## Menu

Statistics > Binary outcomes > Bayesian regression > Heteroskedastic probit regression

## Syntax

```
bayes [ , bayesopts ] : hetprobit depvar [indepvars] [if] [in] [weight] ,
    het(varlist [ , offset(varnameo) ]) [options]
```

<i>options</i>	Description
<b>Model</b>	
* <i>het</i> ( <i>varlist</i> [...])	independent variables to model the variance and possible offset variable
<i>noconstant</i>	suppress constant term
<i>offset</i> ( <i>varname</i> )	include <i>varname</i> in model with coefficient constrained to 1
<i>asis</i>	retain perfect predictor variables
<b>Reporting</b>	
<i>display_options</i>	control spacing, line width, and base and empty cells
<i>level</i> (#)	set credible level; default is <i>level</i> (95)

\**het*() is required. The full specification is *het*(*varlist* [ , *offset*(*varname*<sub>o</sub>) ]).

*indepvars* and *varlist* may contain factor variables; see [U] 11.4.3 **Factor variables**.

*depvar*, *indepvars*, and *varlist* may contain time-series operators; see [U] 11.4.4 **Time-series varlists**.

*fwweights* are allowed; see [U] 11.1.6 **weight**.

*bayes*: *hetprobit*, *level*() is equivalent to *bayes*, *clevel*(): *hetprobit*.

For a detailed description of *options*, see *Options* in [R] **hetprobit**.

<i>bayesopts</i>	Description
<b>Priors</b>	
* <i>normalprior</i> (#)	specify standard deviation of default normal priors for regression coefficients and log-variance coefficients; default is <i>normalprior</i> (100)
<i>prior</i> ( <i>priorspec</i> )	prior for model parameters; this option may be repeated
<i>dryrun</i>	show model summary without estimation
<b>Simulation</b>	
<i>nchains</i> (#)	number of chains; default is to simulate one chain
<i>mcmcsize</i> (#)	MCMC sample size; default is <i>mcmcsize</i> (10000)
<i>burnin</i> (#)	burn-in period; default is <i>burnin</i> (2500)
<i>thinning</i> (#)	thinning interval; default is <i>thinning</i> (1)
<i>rseed</i> (#)	random-number seed
<i>exclude</i> ( <i>paramref</i> )	specify model parameters to be excluded from the simulation results
<b>Blocking</b>	
* <i>blocksize</i> (#)	maximum block size; default is <i>blocksize</i> (50)
<i>block</i> ( <i>paramref</i> [ , <i>blockopts</i> ])	specify a block of model parameters; this option may be repeated
<i>blocksummary</i>	display block summary
* <i>noblocking</i>	do not block parameters by default

#### Initialization

<code>initial(<i>initspec</i>)</code>	specify initial values for model parameters with a single chain
<code>init#(<i>initspec</i>)</code>	specify initial values for #th chain; requires <code>nchains()</code>
<code>initall(<i>initspec</i>)</code>	specify initial values for all chains; requires <code>nchains()</code>
<code>nomleinitial</code>	suppress the use of maximum likelihood estimates as starting values
<code>initrandom</code>	specify random initial values
<code>initsummary</code>	display initial values used for simulation
* <code>noisily</code>	display output from the estimation command during initialization

#### Adaptation

<code>adaptation(<i>adaptopts</i>)</code>	control the adaptive MCMC procedure
<code>scale(#)</code>	initial multiplier for scale factor; default is <code>scale(2.38)</code>
<code>covariance(cov)</code>	initial proposal covariance; default is the identity matrix

#### Reporting

<code>clevel(#)</code>	set credible interval level; default is <code>clevel(95)</code>
<code>hpd</code>	display HPD credible intervals instead of the default equal-tailed credible intervals
<code>eform[ (<i>string</i>) ]</code>	report exponentiated coefficients and, optionally, label as <i>string</i>
<code>batch(#)</code>	specify length of block for batch-means calculations; default is <code>batch(0)</code>
<code>saving(filename[ , replace ])</code>	save simulation results to <i>filename.dta</i>
<code>nomodelsummary</code>	suppress model summary
<code>chainsdetail</code>	display detailed simulation summary for each chain
<code>[no]dots</code>	suppress dots or display dots every 100 iterations and iteration numbers every 1,000 iterations; default is <code>nodots</code>
<code>dots#[ , every(#) ]</code>	display dots as simulation is performed
<code>[no]show(<i>paramref</i>)</code>	specify model parameters to be excluded from or included in the output
<code>notable</code>	suppress estimation table
<code>noheader</code>	suppress output header
<code>title(<i>string</i>)</code>	display <i>string</i> as title above the table of parameter estimates
<code>display_options</code>	control spacing, line width, and base and empty cells

#### Advanced

<code>search(<i>search_options</i>)</code>	control the search for feasible initial values
<code>corrlag(#)</code>	specify maximum autocorrelation lag; default varies
<code>corrtol(#)</code>	specify autocorrelation tolerance; default is <code>corrtol(0.01)</code>

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\* Starred options are specific to the `bayer` prefix; other options are common between `bayer` 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}` for the main regression and `{lnsigma: varlist}` for the log-variance equation. Use the `dryrun` option to see the definitions of model parameters prior to estimation.

For a detailed description of *bayeropts*, see *Options* in [BAYES] `bayer`.

## 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] [hetprobit](#).

For a simple example of the bayes prefix, see *Introductory example* in [BAYES] [bayes](#).

## 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] [hetprobit](#) — Heteroskedastic probit model

[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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