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## Description

`bayes: heckprobit` fits a Bayesian sample-selection probit regression to a partially observed binary outcome; see [\[BAYES\] bayes](#) and [\[R\] heckprobit](#) for details.

## Quick start

Bayesian sample-selection probit regression of  $y$  on  $x_1$  and  $x_2$ , using  $z_1$  and  $z_2$  to model selection and using default normal priors for regression coefficients and atanh-correlation

```
bayes: heckprobit y x1 x2, select(z1 z2)
```

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

```
bayes, normalprior(10): heckprobit y x1 x2, select(z1 z2)
```

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)): heckprobit y x1 x2, select(z1 z2)
```

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

```
bayes, saving(simdata) rseed(123):, ///  
heckprobit y x1 x2, select(z1 z2)
```

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):, ///  
heckprobit y x1 x2, select(z1 z2)
```

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 [\[R\] heckprobit](#).

## Menu

Statistics > Binary outcomes > Bayesian regression > Probit model with sample selection

## Syntax

```
bayes [ , bayesopts ] : heckprobit depvar indepvars [ if ] [ in ] [ weight ] ,
  select( [ depvars = ] varlists [ , noconstant offset(varnameo) ] ) [ options ]
```

*options*

Description

### Model

\* select() specify selection equation: dependent and independent variables; whether to have constant term and offset variable

noconstant suppress constant term

offset(*varname*) include *varname* in model with coefficient constrained to 1

### Reporting

*display\_options* control spacing, line width, and base and empty cells

level(#) set credible level; default is level(95)

\*select() is required.

The full specification is select( [ *depvar*<sub>s</sub> = ] *varlist*<sub>s</sub> [ , noconstant offset(*varname*<sub>o</sub>) ] ).

*indepvars* and *varlist*<sub>s</sub> may contain factor variables; see [U] 11.4.3 Factor variables.

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

fweights are allowed; see [U] 11.1.6 weight.

bayes: heckprobit, level() is equivalent to bayes, clevel(): heckprobit.

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

*bayesopts*

Description

### Priors

\* normalprior(#) specify standard deviation of default normal priors for regression coefficients and atanh-correlation; default is normalprior(100)

prior(*priorspec*) prior for model parameters; this option may be repeated

dryrun show model summary without estimation

### Simulation

nchains(#) number of chains; default is to simulate one chain

mcmcsize(#) MCMC sample size; default is mcmcsize(10000)

burnin(#) burn-in period; default is burnin(2500)

thinning(#) thinning interval; default is thinning(1)

rseed(#) random-number seed

exclude(*paramref*) specify model parameters to be excluded from the simulation results

### Blocking

\* blocksize(#) maximum block size; default is blocksize(50)

block(*paramref* [ , *blockopts* ]) specify a block of model parameters; this option may be repeated

blocksummary display block summary

\* noblocking 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 `{select: varlist_s}` for the selection equation, and atanh-transformed correlation `{athrho}`. 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] [heckprobit](#).

For a simple example of the bayes prefix, see *Introductory example* in [BAYES] [bayes](#). Also see *Heckman selection model* 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] [heckprobit](#) — Probit model with sample selection

[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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