# Title

bayes: truncreg — Bayesian truncated regression

Description	Quick start	Menu
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#### Syntax Also see

# Description

bayes: truncreg fits a Bayesian truncated linear regression to a continuous outcome; see [BAYES] bayes and [R] truncreg for details.

# Quick start

- Bayesian truncated linear regression of y on x1 and x2, using a lower truncation limit of 17 and using default normal priors for regression coefficients and default inverse-gamma prior for the variance bayes: truncreg y x1 x2, ll(17)
- Use a standard deviation of 10 instead of 100 for the default normal priors bayes, normalprior(10): truncreg y x1 x2, ll(17)
- Use a shape of 1 and a scale of 2 instead of values of 0.01 for the default inverse-gamma prior bayes, igammaprior(1 2): truncreg y x1 x2, ll(17)
- 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)): truncreg y x1 x2, ll(17)
- Save simulation results to simdata.dta, and use a random-number seed for reproducibility bayes, saving(simdata) rseed(123):, /// truncreg y x1 x2, ll(17)
- 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):, /// truncreg y x1 x2, ll(17)
- 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] truncreg.

## Menu

Statistics > Linear models and related > Bayesian regression > Truncated regression

# **Syntax**

<u>blocksumm</u>ary

\* noblocking

options	Description	
Model		
<u>nocons</u> tant	suppress constant term	
ll( <i>varname</i>   #)	left-truncation variable or limit	
ul( <i>varname</i>   #)	right-truncation variable or limit	
<u>off</u> set( <i>varname</i> )	include varname in model with coefficient constrained to 1	
Reporting		
display_options	control spacing, line width, and base and empty cells	
<u>l</u> evel(#)	set credible level; default is level(95)	
	is equivalent to bayes, clevel(): truncreg. options, see Options in [R] truncreg. Description	
pavesopis	Description	
Priors		
Priors	specify standard deviation of default normal priors for regression	
Priors	specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) specify shape and scale of default inverse-gamma prior for	
Priors formalprior(#) figammaprior(# #)	<pre>specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) specify shape and scale of default inverse-gamma prior for variance; default is igammaprior(0.01 0.01)</pre>	
Priors * normalprior(#) * igammaprior(# #) prior(priorspec)	<pre>specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) specify shape and scale of default inverse-gamma prior for variance; default is igammaprior(0.01 0.01) prior for model parameters; this option may be repeated</pre>	
Priors <u>normalpr</u> ior(#) <u>igammapr</u> ior(# #) prior( <i>priorspec</i> ) dryrun	<pre>specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) specify shape and scale of default inverse-gamma prior for variance; default is igammaprior(0.01 0.01)</pre>	
Priors normalprior(#) igammaprior(# #) prior( <i>priorspec</i> ) dryrun Simulation	<pre>specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) specify shape and scale of default inverse-gamma prior for variance; default is igammaprior(0.01 0.01) prior for model parameters; this option may be repeated show model summary without estimation</pre>	
Priors normalprior(#) igammaprior(# #) prior(priorspec) dryrun Simulation nchains(#)	<pre>specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) specify shape and scale of default inverse-gamma prior for variance; default is igammaprior(0.01 0.01) prior for model parameters; this option may be repeated show model summary without estimation number of chains; default is to simulate one chain</pre>	
Priors normalprior(#) igammaprior(# #) prior(priorspec) dryrun Simulation nchains(#) <u>mcmcs</u> ize(#)	<pre>specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) specify shape and scale of default inverse-gamma prior for variance; default is igammaprior(0.01 0.01) 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)</pre>	
Priors normalprior(#) figammaprior(# #) prior(priorspec) dryrun Simulation nchains(#) <u>mcmcsize(#)</u> <u>burn</u> in(#)	<pre>specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) specify shape and scale of default inverse-gamma prior for variance; default is igammaprior(0.01 0.01) 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)</pre>	
Priors fnormalprior(#) figammaprior(# #) prior(priorspec) dryrun Simulation nchains(#) <u>mcmcs</u> ize(#) <u>burn</u> in(#) <u>thin</u> ning(#)	<pre>specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) specify shape and scale of default inverse-gamma prior for variance; default is igammaprior(0.01 0.01) 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)</pre>	
Priors * normalprior(#) * igammaprior(# #) prior(priorspec) dryrun Simulation nchains(#) mcmcsize(#) burnin(#) thinning(#) rseed(#)	<pre>specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) specify shape and scale of default inverse-gamma prior for variance; default is igammaprior(0.01 0.01) 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</pre>	
Priors * normalprior(#) * <u>igammapr</u> ior(# #) prior( <i>priorspec</i> ) dryrun Simulation nchains(#) <u>mcmcs</u> ize(#) <u>burn</u> in(#) <u>thin</u> ning(#)	<pre>specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) specify shape and scale of default inverse-gamma prior for variance; default is igammaprior(0.01 0.01) 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</pre>	
Priors * normalprior(#) * igammaprior(# #) prior(priorspec) dryrun Simulation nchains(#) <u>mcmcsize(#)</u> <u>burn</u> in(#) <u>thin</u> ning(#) rseed(#) <u>excl</u> ude(paramref) Blocking	<pre>specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) specify shape and scale of default inverse-gamma prior for variance; default is igammaprior(0.01 0.01) 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 resu</pre>	
Priors * normalprior(#) * igammaprior(# #) prior(priorspec) dryrun Simulation nchains(#) mcmcsize(#) burnin(#) thinning(#) rseed(#) exclude(paramref) Blocking * blocksize(#)	<pre>specify standard deviation of default normal priors for regression coefficients; default is normalprior(100) specify shape and scale of default inverse-gamma prior for variance; default is igammaprior(0.01 0.01) 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)</pre>	

display block summary

do not block parameters by default

Initialization	
<pre>initial(initspec)</pre>	specify initial values for model parameters with a single chain
<pre>init#(initspec)</pre>	specify initial values for #th chain; requires nchains()
initall( <i>initspec</i> )	specify initial values for all chains; requires nchains()
nomleinitial	suppress the use of maximum likelihood estimates as starting values
initrandom	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( <i>adaptopts</i> )	control the adaptive MCMC procedure
<u>sc</u> ale(#)	initial multiplier for scale factor; default is scale(2.38)
<pre>covariance(cov)</pre>	initial proposal covariance; default is the identity matrix
Reporting	
<u>clev</u> el(#)	set credible interval level; default is clevel(95)
hpd	display HPD credible intervals instead of the default equal-tailed credible intervals
<u>ef</u> orm ( <i>string</i> )	report exponentiated coefficients and, optionally, label as string
batch(#)	<pre>specify length of block for batch-means calculations; default is batch(0)</pre>
<pre>saving(filename[, replace])</pre>	save simulation results to <i>filename</i> .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
dots(# $[, every(#)]$ )	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
<pre>title(string)</pre>	display string as title above the table of parameter estimates
display_options	control spacing, line width, and base and empty cells
Advanced	
<pre>search(search_options)</pre>	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} and variance {sigma2}. 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**

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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] **truncreg**.

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<sup>+</sup>

[R] truncreg — Truncated regression

[BAYES] Bayesian postestimation — Postestimation tools for bayesmh and the bayes prefix

[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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