

bayes: dsgenl — Bayesian nonlinear dynamic stochastic general equilibrium models

[Description](#)[Quick start](#)[Menu](#)[Syntax](#)[Remarks and examples](#)[Stored results](#)[Methods and formulas](#)[Also see](#)

Description

`bayes: dsgenl` fits a Bayesian nonlinear dynamic stochastic general equilibrium (DSGE) model to continuous multivariate time series; see [\[BAYES\] bayes](#) and [\[DSGE\] dsgenl](#) for details.

Quick start

Nonlinear DSGE model in which observed variable y depends on unobserved state z

```
bayes, prior({rho}, uniform(0,1)) prior({alpha}, beta(5,5)): ///
dsgenl (y = z^{alpha}) (ln(F.z) = {rho}*ln(z)),          ///
exostate(z) observed(y)
```

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

```
bayes, prior({rho}, uniform(0,1)) prior({alpha}, beta(5,5)): ///
rseed(17) saving(bdsgenlsim):                          ///
dsgenl (y = z^{alpha}) (ln(F.z) = {rho}*ln(z)),          ///
exostate(z) observed(y)
```

Specify 20,000 Markov chain Monte Carlo (MCMC) samples, and set length of burn-in period to 5,000

```
bayes, prior({rho}, uniform(0,1)) prior({alpha}, beta(5,5)): ///
mcmcsize(20000) burnin(5000):                          ///
dsgenl (y = z^{alpha}) (ln(F.z) = {rho}*ln(z)),          ///
exostate(z) observed(y)
```

Estimate parameters of a four-equation production model. Priors for α , β , and ρ are given by beta distributions with means 0.3, 0.9, and 0.5, respectively

```
bayes, prior({alpha}, beta(3,7))                        ///
prior({beta}, beta(9,1))                               ///
prior({rho}, beta(7,7)) :                             ///
dsgenl (1/c = {alpha}*{beta}*(1/F.c)*(F.y/F.k))        ///
(y = z*k^{alpha}) (F.k = y - c)                       ///
(ln(F.z) = {rho}*ln(z)) ,                             ///
exostate(z) endostate(k) observed(y) unobserved(c)
```

In the above, request that a 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](#).

Menu

Statistics > Multivariate time series > Bayesian models > Nonlinear DSGE models

Syntax

```
bayes, prior(userparams, ...) [bayesopts] : dsge1 (eqn_list) [if] [in] [, options]
```

<i>options</i>	Description
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Model

* <i>observed(string)</i>	list observed control variables
<i>unobserved(string)</i>	list unobserved control variables
* <i>exostate(string)</i>	list exogenous state variables
<i>endostate(string)</i>	list endogenous state variables
<i>linearapprox</i>	take a linear, rather than log-linear, approximation
<i>level(#)</i>	set credible level; default is <i>level(95)</i>
<i>noidencheck</i>	do not check for parameter identification; implied
<i>solve</i>	return model solution at initial values; implied

**observed()* and *exostate()* are required.

bayes: dsge1, level() is equivalent to *bayes, clevel(): dsge1*.

For a detailed description of *options*, see [Options](#) in [\[DSGE\] dsge1](#).

Options *level()*, *noidencheck*, and *stable* do not appear on the dialog box.

<i>bayesopts</i>	Description
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Priors

* <i>igammaprior(# #)</i>	specify shape and scale of default inverse-gamma prior for standard deviations of shocks; default is <i>igammaprior(0.01 0.01)</i>
<i>prior(priorspec)</i>	prior for model parameters; this option may be repeated and is required for all user-defined parameters <i>userparams</i>
<i>dryrun</i>	show model summary without estimation

Simulation

<i>nchains(#)</i>	number of chains; default is to simulate one chain
<i>mcmcsize(#)</i>	MCMC sample size; default is <i>mcmcsize(10000)</i>
<i>burnin(#)</i>	burn-in period; default is <i>burnin(2500)</i>
<i>thinning(#)</i>	thinning interval; default is <i>thinning(1)</i>
<i>rseed(#)</i>	random-number seed
<i>exclude(paramref)</i>	specify model parameters to be excluded from the simulation results

Blocking

<i>block(paramref [, blockopts])</i>	specify a block of model parameters; this option may be repeated
<i>blocksummary</i>	display block summary

Initialization

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

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>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 <code>filename.dta</code>
<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(paramref)</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(string)</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(search_options)</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>

*Starred options are specific to the `bayes` prefix; other options are common between `bayes` and `bayesmh`.
`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 user-defined parameters `userparams` and standard deviations of shocks `{sd(e.exogstate)}`. 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`.

`nomleinitial` is assumed. Default parameter values are set to means of priors.

Remarks and examples

[stata.com](http://www.stata.com)

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 [DSGE] `dsgenl`.

For a simple example of the `bayes` prefix, see *Introductory example* in [BAYES] `bayes`. For an introduction to and examples of Bayesian DSGEs, see [DSGE] `Intro 9` and [DSGE] `Intro 9b`.

Stored results

See *Stored results* in [BAYES] `bayes`. Also see *Stored results* in [DSGE] `dsgenl`.

Methods and formulas

See *Methods and formulas* in [DSGE] **dsge** and [DSGE] **Intro 9**. See *Methods and formulas* in [BAYES] **bayesmh**.

Also see

[BAYES] **bayes: dsge postestimation** — Postestimation tools for bayes: dsge and bayes: dsngen

[BAYES] **bayes** — Bayesian regression models using the bayes prefix

[DSGE] **dsngen** — Nonlinear dynamic stochastic general equilibrium models

[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**