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## Description

`bayes: mlogit` fits a Bayesian multinomial logistic regression to a categorical outcome; see [\[BAYES\] bayes](#) and [\[R\] mlogit](#) for details.

## Quick start

Bayesian multinomial logistic regression of `y` on `x1` and `x2`, using default normal priors for regression coefficients

```
bayes: mlogit y x1 x2
```

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

```
bayes, normalprior(10): mlogit y x1 x2
```

Use uniform priors for the slopes and a normal prior for the intercept for the category 2

```
bayes, prior({2: x1 x2}, uniform(-10,10)) ///  
prior({2: _cons}, normal(0,10)): mlogit y x1 x2
```

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

```
bayes, saving(simdata) rseed(123): mlogit y x1 x2
```

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): mlogit y x1 x2
```

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
```

Display relative-risk ratios instead of coefficients

```
bayes: mlogit y x1 x2, rrr
```

Display relative-risk ratios on replay

```
bayes, rrr
```

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

## Menu

Statistics > Categorical outcomes > Bayesian regression > Multinomial logistic regression

# Syntax

```
bayes [ , bayesopts ] : mlogit depvar [ indepvars ] [ if ] [ in ] [ weight ] [ , options ]
```

<i>options</i>	Description
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## Model

<u>noconstant</u>	suppress constant term
<u>baseoutcome</u> (#)	value of <i>depvar</i> that will be the base outcome

## Reporting

<u>rrr</u>	report relative-risk ratios
<u>display_options</u>	control spacing, line width, and base and empty cells
<u>level</u> (#)	set credible level; default is level(95)

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

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

fweights are allowed; see [U] 11.1.6 weight.

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

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

<i>bayesopts</i>	Description
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## Priors

* <u>normalprior</u> (#)	specify standard deviation of default normal priors for regression coefficients; default is normalprior(100)
<u>prior</u> ( <i>priorspec</i> )	prior for model parameters; this option may be repeated
<u>dryrun</u>	show model summary without estimation

## Simulation

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

## Blocking

* <u>blocksize</u> (#)	maximum block size; default is blocksize(50)
<u>block</u> ( <i>paramref</i> [ , <i>blockopts</i> ])	specify a block of model parameters; this option may be repeated
<u>blocksummary</u>	display block summary
* <u>noblocking</u>	do not block parameters by default

## Initialization

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

## Adaptation

<code>adaptation(adaptopts)</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>rrr</code>	report relative-risk ratios
<code>eform(string)</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(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>corrto1(#)</code>	specify autocorrelation tolerance; default is <code>corrto1(0.01)</code>

\* Starred options are specific to the `bayer` prefix; other options are common between `bayer` and `bayermh`.

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

`priorspec` and `paramref` are defined in [BAYES] `bayermh`.

`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  $\{outcome_1 : indepvars\}$ ,  $\{outcome_2 : indepvars\}$ , and so on, where  $outcome_{\#}$ 's are the values of the dependent variable or the value labels of the dependent variable if they exist. 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] `bayermh`. For remarks and examples specific to the `bayer` prefix, see [BAYES] `bayer`. For details about the estimation command, see [R] `mlogit`.

For a simple example of the `bayer` prefix, see *Introductory example* in [BAYES] `bayer`. Also see *Multinomial logistic regression* in [BAYES] `bayer`.

## Stored results

See *Stored results* in [\[BAYES\] bayes](#).

## Methods and formulas

See *Methods and formulas* in [\[BAYES\] bayesmh](#).

## Reference

Parish, W. J., A. Aldridge, and M. van Hasselt. 2024. [A Bayesian method for addressing multinomial misclassification with applications for alcohol epidemiological modeling](#). *Stata Journal* 24: 113–137.

## Also see

[\[BAYES\] bayes](#) — Bayesian regression models using the bayes prefix

[\[R\] mlogit](#) — Multinomial (polytomous) logistic regression

[\[BAYES\] 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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