Glossary

- **a posteriori**. In the context of Bayesian analysis, we use a posteriori to mean "after the sample is observed". For example, a posteriori information is any information obtained after the data sample is observed. See *posterior distribution, posterior*.
- **a priori**. In the context of Bayesian analysis, we use a priori to mean "before the sample is observed". For example, a priori information is any information obtained before the data sample is observed. In a Bayesian model, a priori information about model parameters is specified by prior distributions.
- **acceptance rate**. In the context of the MH algorithm, acceptance rate is the fraction of the proposed samples that is accepted. The optimal acceptance rate depends on the properties of the target distribution and is not known in general. If the target distribution is normal, however, the optimal acceptance rate is known to be 0.44 for univariate distributions and 0.234 for multivariate distributions.
- **adaptation**. In the context of the MH algorithm, adaptation refers to the process of tuning or adapting the proposal distribution to optimize the MCMC sampling. Typically, adaptation is performed periodically during the MCMC sampling. The bayesmh command performs adaptation every # of iterations as specified in option adaptation(every(#)) for a maximum of adaptation(maxiter()) iterations. In a continuous-adaptation regimes, the adaptation lasts during the entire process of the MCMC sampling. See [BAYES] bayesmh.
- adaptation period. Adaptation period includes all MH adaptive iterations. It equals the length of the adaptation interval, as specified by adaptation(every()), times the maximum number of adaptations, adaptation(maxiter()).
- **adaptive iteration**. In the adaptive MH algorithm, adaptive iterations are iterations during which adaptation is performed.
- Akaike information criterion, AIC. Akaike information criterion (AIC) is an information-based modelselection criterion. It is given by the formula $-2 \times \log$ likelihood + 2k, where k is the number of parameters. AIC favors simpler models by penalizing for the number of model parameters. It does not, however, account for the sample size. As a result, the AIC penalization diminishes as the sample size increases, as does its ability to guard against overparameterization.
- **batch means**. Batch means are means obtained from batches of sample values of equal size. Batch means provide an alternative method for estimating MCMC standard errors (MCSE). The batch size is usually chosen to minimize the correlation between different batches of means.
- **Bayes factor**. Bayes factor is given by the ratio of the marginal likelihoods of two models, M_1 and M_2 . It is a widely used criterion for Bayesian model comparison. Bayes factor is used in calculating the posterior odds ratio of model M_1 versus M_2 ,

$$\frac{P(M_1|\mathbf{y})}{P(M_2|\mathbf{y})} = \frac{P(\mathbf{y}|M_1)}{P(\mathbf{y}|M_2)} \frac{P(M_1)}{P(M_2)}$$

where $P(M_i|\mathbf{y})$ is a posterior probability of model M_i , and $P(M_i)$ is a prior probability of model M_i . When the two models are equally likely, that is, when $P(M_1) = P(M_2)$, the Bayes factor equals the posterior odds ratio of the two models.

Bayes's theorem. The Bayes's theorem is a formal method for relating conditional probability statements. For two (random) events X and Y, the Bayes's theorem states that

$$P(X|Y) \propto P(Y|X)P(X)$$

that is, the probability of X conditional on Y is proportional to the probability of X and the probability of Y conditional on X. In Bayesian analysis, the Bayes's theorem is used for combining prior information about model parameters and evidence from the observed data to form the posterior distribution.

Bayesian analysis. Bayesian analysis is a statistical methodology that considers model parameters to be random quantities and estimates their posterior distribution by combining prior knowledge about parameters with the evidence from the observed data sample. Prior knowledge about parameters is described by prior distributions and evidence from the observed data is incorporated through a likelihood model. Using the Bayes's theorem, the prior distribution and the likelihood model are combined to form the posterior distribution of model parameters. The posterior distribution is then used for parameter inference, hypothesis testing, and prediction.

- **Bayesian estimation**. Bayesian estimation consists of fitting Bayesian models and estimating their parameters based on the resulting posterior distribution. Bayesian estimation in Stata can be done using the convenient bayes prefix or the more general bayesmh command. See [BAYES] **Bayesian estimation** for details.
- Bayesian estimation results. Estimation results obtained after the bayes prefix or the bayesmh command.
- **Bayesian hypothesis testing**. Bayesian hypothesis testing computes probabilities of hypotheses conditional on the observed data. In contrast to the frequentist hypothesis testing, the Bayesian hypothesis testing computes the actual probability of a hypothesis *H* by using the Bayes's theorem,

$$P(H|\mathbf{y}) \propto P(\mathbf{y}|H)P(H)$$

where **y** is the observed data, $P(\mathbf{y}|H)$ is the marginal likelihood of **y** given *H*, and P(H) is the prior probability of *H*. Two different hypotheses, H_1 and H_2 , can be compared by simply comparing $P(H_1|\mathbf{y})$ to $P(H_2|\mathbf{y})$.

- **Bayesian information criterion, BIC**. The Bayesian information criterion (BIC), also known as Schwarz criterion, is an information based criterion used for model selection in classical statistics. It is given by the formula $-2 \times \log$ likelihood $+ k \times \ln n$, where k is the number of parameters and n is the sample size. BIC favors simpler, in terms of complexity, models and it is more conservative than AIC.
- Bayesian lasso prior. See global-local shrinkage prior.
- **Bayesian model checking**. In Bayesian statistics, model checking refers to testing likelihood and prior model adequacy in the context of a research problem and observed data. A simple sanity check may include verifying that posterior inference produces results that are reasonable in the context of the problem. More substantive checks may include analysis of the sensitivity of Bayesian inference to changes in likelihood and prior distribution specifications. See *posterior predictive checking*.
- **Bayesian predictions**. Bayesian predictions are samples from the posterior predictive distribution of outcome variables and functions of these samples and, optionally, model parameters. Examples of Bayesian predictions include replicated data, out-of-sample predictions, and test statistics of simulated outcomes.
- **Bayesian variable selection**. Bayesian variable selection uses special priors (global-local shrinkage priors and spike-and-slab priors) for regression coefficients to regulate the effects of predictors on an outcome of interest in a regression model. These priors and the information from the observed data are used to shrink regression coefficients toward zero based on how well predictors explain the outcome. Unlike some classical variable-selection methods that either include or exclude each predictor,

Bayesian variable selection uses all predictors but downweights the effects of weak predictors. It also produces coefficient estimates that account for variable-selection uncertainty. See *Remarks and* examples of [BAYES] **bayesselect**.

- **blocking**. In the context of the MH algorithm, blocking refers to the process of separating model parameters into different subsets or blocks to be sampled independently of each other. MH algorithm generates proposals and applies the acceptance–rejection rule sequentially for each block. It is recommended that correlated parameters are kept in one block. Separating less-correlated or independent model parameters in different blocks may improve the mixing of the MH algorithm.
- **burn-in period**. The burn-in period is the number of iterations it takes for an MCMC sequence to reach stationarity.
- central posterior interval. See equal-tailed credible interval.
- conditional conjugacy. See semiconjugate prior.
- **conjugate prior**. A prior distribution is conjugate for a family of likelihood distributions if the prior and posterior distributions belong to the same family of distributions. For example, the gamma distribution is a conjugate prior for the Poisson likelihood. Conjugacy may provide an efficient way of sampling from posterior distributions and is used in Gibbs sampling.
- continuous parameters. Continuous parameters are parameters with continuous prior distributions.
- **credible interval**. In Bayesian analysis, the credible interval of a scalar model parameter is an interval from the domain of the marginal posterior distribution of that parameter. Two types of credible intervals are typically used in practice: equal-tailed credible intervals and HPD credible intervals.
- **credible level**. The credible level is a probability level between 0% and 100% used for calculating credible intervals in Bayesian analysis. For example, a 95% credible interval for a scalar parameter is an interval the parameter belongs to with the probability of 95%.
- **cross-variable**. A cross-variable is a Stata term to refer to the dependent variable used as a lagged regressor in the VAR model in an outcome equation that is not its own. For instance, in a VAR model with two dependent variables (y1 and y2) and two lags,

$$\begin{split} \mathbf{y1} &= a_{11}\mathbf{L}.\mathbf{y1} + a_{12}\mathbf{L}2.\mathbf{y1} + a_{21}\mathbf{L}.\mathbf{y2} + a_{22}\mathbf{L}2.\mathbf{y2} + a_0 + u_1 \\ \mathbf{y2} &= b_{11}\mathbf{L}.\mathbf{y1} + b_{12}\mathbf{L}2.\mathbf{y1} + b_{21}\mathbf{L}.\mathbf{y2} + b_{22}\mathbf{L}2.\mathbf{y2} + b_0 + u_2 \end{split}$$

y2 is the cross-variable in the first outcome equation (y1), and y1 is the cross-variable in the second outcome equation (y2). Note that y1 and y2 are self-variables in the first and second equations, respectively. Cross-variables are always endogenous variables. We also often refer to the coefficients of cross-variables as cross-variable first-lag coefficients (a_{21} and b_{11}), cross-variable second-lag coefficients (a_{22} and b_{12}), and so on. See [BAYES] bayes: var.

- cross-variable first-lag coefficients. Regression coefficients in a VAR model that correspond to first lags ([U] 11.4.4 Time-series varlists) of cross-variables. See [BAYES] bayes: var.
- cross-variable tightness parameter. A cross-variable tightness parameter is a parameter, λ_2 , that controls the tightness of the Minnesota prior distribution by controlling the prior variance for the cross-variable coefficients. It is specified in the Minnesota prior option crosstight(). See Methods and formulas of [BAYES] bayes: var for details.
- **cusum plot, CUSUM plot.** The cusum (CUSUM) plot of an MCMC sample is a plot of cumulative sums of the differences between sample values and their overall mean against the iteration number. Cusum plots are useful graphical summaries for detecting early drifts in MCMC samples.

- deviance information criterion, DIC. The deviance information criterion (DIC) is an information based criterion used for Bayesian model selection. It is an analog of AIC and is given by the formula $D(\bar{\theta}) + 2 \times p_D$, where $D(\bar{\theta})$ is the deviance at the sample mean and p_D is the effective complexity, a quantity equivalent to the number of parameters in the model. Models with smaller DIC are preferred.
- **diminishing adaptation**. Diminishing adaptation of the adaptive algorithm is the type of adaptation in which the amount of adaptation decreases with the size of the MCMC chain.
- discrete parameters. Discrete parameters are parameters with discrete prior distributions.
- effective sample size, ESS. Effective sample size (ESS) is the MCMC sample size T adjusted for the autocorrelation in the sample. It represents the number of independent observations in an MCMC sample. ESS is used instead of T in calculating MCSE. Small ESS relative to T indicates high autocorrelation and consequently poor mixing of the chain.
- efficiency. In the context of MCMC, efficiency is a term used for assessing the mixing quality of an MCMC procedure. Efficient MCMC algorithms are able to explore posterior domains in less time (using fewer iterations). Efficiency is typically quantified by the sample autocorrelation and effective sample size. An MCMC procedure that generates samples with low autocorrelation and consequently high ESS is more efficient.
- endogenous variable. In the context of VAR, an endogenous variable is a dependent variable included in the model as a regressor with a lag operator; see [U] 11.4.4 Time-series variates. Also see endogenous variable in [TS] Glossary for a general definition.
- equal-tailed credible interval. An equal-tailed credible interval is a credible interval defined in such a way that both tails of the marginal posterior distribution have the same probability. A $\{100 \times (1-\alpha)\}$ % equal-tailed credible interval is defined by the $\alpha/2$ th and $\{(1-\alpha)/2\}$ th quantiles of the marginal posterior distribution.
- exogenous variable. In the context of VAR, an exogenous variable is an independent variable (regressor) included in the model. Also see *exogenous variable* in [TS] **Glossary** for a general definition.
- exogenous variable tightness parameter. An exogenous variable tightness parameter is a parameter, λ_4 , that controls the tightness of the Minnesota prior distribution by controlling the prior variance for the exogenous variable coefficients. It is specified in the Minnesota prior option exogtight(). See Methods and formulas of [BAYES] bayes: var for details.
- **feasible initial value**. An initial-value vector is feasible if it corresponds to a state with a positive posterior probability.
- fixed effects. See fixed-effects parameters.
- **fixed-effects parameters**. In the Bayesian context, the term "fixed effects" or "fixed-effects parameters" is a misnomer, because all model parameters are inherently random. We use this term in the context of Bayesian multilevel models to refer to regression model parameters and to distinguish them from the random-effects parameters. You can think of fixed-effects parameters as parameters modeling population averaged or marginal relationship of the response and the variables of interest.
- **frequentist analysis.** Frequentist analysis is a form of statistical analysis where model parameters are considered to be unknown but fixed constants and the observed data are viewed as a repeatable random sample. Inference is based on the sampling distribution of the data.
- **full conditionals**. A full conditional is the probability distribution of a random variate conditioned on all other random variates in a joint probability model. Full conditional distributions are used in Gibbs sampling.

full Gibbs sampling. See Gibbs sampling, Gibbs sampler.

- Gelman-Rubin convergence diagnostic, Gelman-Rubin convergence statistic. Gelman-Rubin convergence diagnostic assesses MCMC convergence by analyzing differences between multiple Markov chains. The convergence is assessed by comparing the estimated between-chains and within-chain variances for each model parameter. Large differences between these variances indicate nonconvergence. See [BAYES] bayesstats grubin.
- **Gibbs sampling, Gibbs sampler**. Gibbs sampling is an MCMC method, according to which each random variable from a joint probability model is sampled according to its full conditional distribution.
- global–local shrinkage prior. A global–local shrinkage prior is one type of prior used for regression coefficients in Bayesian variable selection. The prior for coefficient β_j is normal with mean zero and standard deviation $\tau \lambda_j$, where τ is a global shrinkage parameter and λ_j is a local shrinkage parameter. The priors used for τ and λ_j 's depend on the type of the global–local shrinkage prior. For the horseshoe prior, these parameters have half-Cauchy distributions. For the Bayesian lasso prior, λ_j 's have a Rayleigh distribution instead of half-Cauchy. See *Global–local shrinkage priors* in *Methods and formulas* of [BAYES] bayesselect.
- **highest posterior density credible interval, HPD credible interval**. The highest posterior density (HPD) credible interval is a type of a credible interval with the highest marginal posterior density. An HPD interval has the shortest width among all other credible intervals. For some multimodal marginal distributions, HPD may not exists. See *highest posterior density region*, *HPD region*.
- **highest posterior density region, HPD region**. The highest posterior density (HPD) region for model parameters has the highest marginal posterior probability among all domain regions. Unlike an HPD credible interval, an HPD region always exist.
- horseshoe prior. See global-local shrinkage prior.
- **hybrid MH sampling, hybrid MH sampler**. A hybrid MH sampler is an MCMC method in which some blocks of parameters are updated using the MH algorithms and other blocks are updated using Gibbs sampling.
- **hyperparameter**. In Bayesian analysis, hyperparameter is a parameter of a prior distribution, in contrast to a model parameter.
- **hyperprior**. In Bayesian analysis, hyperprior is a prior distribution of hyperparameters. See *hyperparameter*.
- **improper prior**. A prior is said to be improper if it does not integrate to a finite number. Uniform distributions over unbounded intervals are improper. Improper priors may still yield proper posterior distributions. When using improper priors, however, one has to make sure that the resulting posterior distribution is proper for Bayesian inference to be invalid.
- inclusion coefficients. Inclusion coefficients γ_j 's describe the importance of predictors in Bayesian variable selection. For global-local shrinkage priors, $\gamma_j = (\lambda_j^2/\lambda_0^2)/(1 + \lambda_j^2/\lambda_0^2)$, where λ_j is a local shrinkage parameter associated with coefficient β_j and λ_0 is a prespecified positive scale value. For spike-and-slab priors, which are mixtures of two priors, γ_j 's are the binary indicators. For these priors, the inclusion coefficients can be interpreted as inclusion probabilities. Predictors with inclusion coefficients greater than 0.5 may be considered important predictors. See *Remarks and examples* and *Methods and formulas* of [BAYES] bayesselect.
- **independent a posteriori**. Parameters are considered independent a posteriori if their marginal posterior distributions are independent; that is, their joint posterior distribution is the product of their individual marginal posterior distributions.

- independent a priori. Parameters are considered independent a priori if their prior distributions are independent; that is, their joint prior distribution is the product of their individual marginal prior distributions.
- **informative prior**. An informative prior is a prior distribution that has substantial influence on the posterior distribution.
- in-sample predictions. See replicated outcome.
- interval hypothesis testing. Interval hypothesis testing performs interval hypothesis tests for model parameters and functions of model parameters.
- interval test. In Bayesian analysis, an interval test applied to a scalar model parameter calculates the marginal posterior probability for the parameter to belong to the specified interval.
- Jeffreys prior. The Jeffreys prior of a vector of model parameters θ is proportional to the square root of the determinant of its Fisher information matrix $I(\theta)$. Jeffreys priors are locally uniform and, by definition, agree with the likelihood function. Jeffreys priors are considered noninformative priors that have minimal impact on the posterior distribution.
- lag coefficient. In the context of time-series regression analysis, a lag coefficient is a regression coefficient that corresponds to a variable included in the regression model with a lag operator; see [U] 11.4.4 Time-series varlists.
- **lag-decay parameter**. A lag-decay parameter is a parameter, λ_3 , that controls the tightness of the Minnesota prior distribution by controlling the prior variance as a function of a lag for all endogenous variable coefficients. It is specified in the Minnesota prior option lagdecay(). See *Methods and formulas* of [BAYES] **bayes: var** for details.
- **marginal distribution**. In Bayesian context, a distribution of the data after integrating out parameters from the joint distribution of the parameters and the data.
- marginal likelihood. In the context of Bayesian model comparison, a marginalized over model parameters $\boldsymbol{\theta}$ likelihood of data \mathbf{y} for a given model M, $P(\mathbf{y}|M) = m(\mathbf{y}) = \int P(\mathbf{y}|\boldsymbol{\theta}, M)P(\boldsymbol{\theta}|M)d\boldsymbol{\theta}$. Also see Bayes factor.
- **marginal posterior distribution**. In Bayesian context, a marginal posterior distribution is a distribution resulting from integrating out all but one parameter from the joint posterior distribution.
- **Markov chain**. Markov chain is a random process that generates sequences of random vectors (or states) and satisfies the Markov property: the next state depends only on the current state and not on any of the previous states. MCMC is the most common methodology for simulating Markov chains.
- **matrix model parameter**. A matrix model parameter is any model parameter that is a matrix. Matrix elements, however, are viewed as scalar model parameters.

Matrix model parameters are defined and referred to within the bayesmh command as {param, matrix} or {eqname:param, matrix} with the equation name eqname. For example, {Sigma, matrix} and {Scale:Omega, matrix} are matrix model parameters. Individual matrix elements cannot be referred to within the bayesmh command, but they can be referred within postestimation commands accepting parameters. For example, to refer to the individual elements of the defined above, say, 2 × 2 matrices, use {Sigma_1_1}, {Sigma_2_1}, {Sigma_1_2}, {Sigma_2_2} and {Scale:Omega_1_1}, {Scale:Omega_2_1}, {Scale:Omega_2_2}, respectively. See [BAYES] bayesmh.

matrix parameter. See matrix model parameter.

- **MCMC, Markov chain Monte Carlo**. MCMC is a class of simulation-based methods for generating samples from probability distributions. Any MCMC algorithm simulates a Markov chain with a target distribution as its stationary or equilibrium distribution. The precision of MCMC algorithms increases with the number of iterations. The lack of a stopping rule and convergence rule, however, makes it difficult to determine for how long to run MCMC. The time needed to converge to the target distribution within a prespecified error is referred to as mixing time. Better MCMC algorithms have faster mixing times. Some of the popular MCMC algorithms are random-walk Metropolis, Metropolis–Hastings, and Gibbs sampling.
- MCMC replicates. An MCMC sample of simulated outcomes.
- **MCMC sample**. An MCMC sample is obtained from MCMC sampling. An MCMC sample approximates a target distribution and is used for summarizing this distribution.
- MCMC sample size. MCMC sample size is the size of the MCMC sample. It is specified in bayesmh's option mcmcsize(); see [BAYES] bayesmh.
- **MCMC sampling, MCMC sampler**. MCMC sampling is an MCMC algorithm that generates samples from a target probability distribution.
- **MCMC standard error, MCSE** MCSE is the standard error of the posterior mean estimate. It is defined as the standard deviation divided by the square root of ESS. MCSEs are analogs of standard errors in frequentist statistics and measure the accuracy of the simulated MCMC sample.
- Metropolis–Hastings (MH) sampling, MH sampler. A Metropolis–Hastings (MH) sampler is an MCMC method for simulating probability distributions. According to this method, at each step of the Markov chain, a new proposal state is generated from the current state according to a prespecified proposal distribution. Based on the current and new state, an acceptance probability is calculated and then used to accept or reject the proposed state. Important characteristics of MH sampling is the acceptance rate and mixing time. The MH algorithm is very general and can be applied to an arbitrary target distribution. However, its efficiency is limited, in terms of mixing time, and decreases as the dimension of the target distribution increases. Gibbs sampling, when available, can provide much more efficient sampling than MH sampling.
- **Minnesota prior**. In Bayesian VAR models, Minnesota priors are used as priors for regression coefficients. A Minnesota prior is a multivariate normal distribution with a special mean vector and covariance matrix. The mean vector contains all zeroes except the values corresponding to self-variable first-lag coefficients, which are set to 1. The covariance matrix can be fixed or can be a product of a fixed matrix and a matrix model parameter as in the case of a conjugate Minnesota prior. The Minnesota prior assumes that, a priori, each univariate time series in the model is a random walk.
- **Minnesota factor covariance**. In Bayesian VAR models with a conjugate Minnesota prior, the factor covariance matrix is used to form the covariance of the multivariate normal prior for regression co-efficients. The latter is defined as the Kronecker product of the unknown covariance matrix of error terms with the Minnesota factor covariance.
- **mixing of Markov chain**. Mixing refers to the rate at which a Markov chain traverses the parameter space. It is a property of the Markov chain that is different from convergence. Poor mixing indicates a slow rate at which the chain explores the stationary distribution and will require more iterations to provide inference at a given precision. Poor (slow) mixing is typically a result of high correlation between model parameters or of weakly-defined model specifications.
- **model hypothesis testing**. Model hypothesis testing tests hypotheses about models by computing model posterior probabilities.

- model parameter. A model parameter refers to any (random) parameter in a Bayesian model. Model parameters can be scalars or matrices. Examples of model parameters as defined in bayesmh are {mu}, {scale:s}, {Sigma,matrix}, and {Scale:Omega,matrix}. See [BAYES] bayesmh and, specifically, Declaring model parameters and Referring to model parameters in that entry. Also see Different ways of specifying model parameters in [BAYES] Bayesian postestimation.
- **model posterior probability**. Model posterior probability is probability of a model *M* computed conditional on the observed data **y**,

$$P(M|\mathbf{y}) = P(M)P(\mathbf{y}|M) = P(M)m(\mathbf{y})$$

where P(M) is the prior probability of a model M and $m(\mathbf{y})$ is the marginal likelihood under model M.

- **noninformative prior**. A noninformative prior is a prior with negligible influence on the posterior distribution. See, for example, *Jeffreys prior*.
- objective prior. See noninformative prior.
- **one-at-a-time MCMC sampling.** A one-at-a-time MCMC sample is an MCMC sampling procedure in which random variables are sampled individually, one at a time. For example, in Gibbs sampling, individual variates are sampled one at a time, conditionally on the most recent values of the rest of the variates.
- out-of-sample predictions. Predictions of future observations; see simulated outcome.
- **overdispersed initial value**. An overdispersed initial value is obtained from a distribution that is overdispersed or has larger variability relative to the true marginal posterior distribution. Overdispersed initial values are used with multiple Markov chains for diagnosing MCMC convergence. Also see *Specifying initial values* in [BAYES] **bayesmh**.
- **posterior distribution, posterior**. A posterior distribution is a probability distribution of model parameters conditional on observed data. The posterior distribution is determined by the likelihood of the parameters and their prior distribution. For a parameter vector $\boldsymbol{\theta}$ and data y, the posterior distribution is given by

$$P(\boldsymbol{\theta}|\mathbf{y}) = \frac{P(\boldsymbol{\theta})P(\mathbf{y}|\boldsymbol{\theta})}{P(\mathbf{y})}$$

where $P(\theta)$ is the prior distribution, $P(\mathbf{y}|\theta)$ is the model likelihood, and $P(\mathbf{y})$ is the marginal distribution for \mathbf{y} . Bayesian inference is based on a posterior distribution.

- posterior independence. See independent a posteriori.
- posterior interval. See credible interval.
- **posterior odds**. Posterior odds for θ_1 compared with θ_2 is the ratio of posterior density evaluated at θ_1 and θ_2 under a given model,

$$\frac{p(\boldsymbol{\theta}_1|\mathbf{y})}{p(\boldsymbol{\theta}_2|\mathbf{y})} = \frac{p(\boldsymbol{\theta}_1)}{p(\boldsymbol{\theta}_2)} \frac{p(\mathbf{y}|\boldsymbol{\theta}_1)}{p(\mathbf{y}|\boldsymbol{\theta}_2)}$$

In other words, posterior odds are prior odds times the likelihood ratio.

posterior predictive checking. Posterior predictive checking is a methodology for assessing goodness of fit of a Bayesian model using replicated data simulated from the posterior predictive distribution of the model. For example, graphical diagnostics of the replicated residuals may be used to check the distributional assumptions of the model error terms. A more formal and systematic approach uses test quantities and test statistics to measure discrepancies between replicated data and observed data. Test

statistics such as a mean, minimum, and maximum can be used to compare different aspects of the observed data distribution with those of the replicated-data distribution. Posterior predictive *p*-values, also called Bayesian *p*-values, computed for test quantities and test statistics are used to quantify the discrepancy between the observed and replicated data. Also see *Bayesian model checking*.

- **posterior predictive distribution**. Posterior predictive distribution is a distribution of unobserved (future) data conditional on observed data. Posterior predictive distribution is derived by marginalizing the likelihood function with respect to the posterior distribution of model parameters.
- **posterior predictive p-value**. Posterior predictive *p*-value, also called a Bayesian *p*-value, is the probability that a test quantity (or statistic) computed for the replicated data is greater or equal to the test quantity computed for the observed data. Posterior predictive *p*-values are used in posterior predictive checking. *p*-values less than 0.05 or greater than 0.95 typically indicate model misfit (Gelman et al. 2014).
- predictive distribution. See prior predictive distribution and posterior predictive distribution.
- **predictive inference**. In Bayesian statistics, predictive inference is inference about unobserved (future) data conditionally on past data and prior knowledge of model parameters. Predictive inference is based on prior predictive or posterior predictive distribution of model parameters.
- **predictive outcome**. Predictive outcome \tilde{y} is a value or a set of values simulated from a posterior predictive distribution $p(\tilde{y}|y)$ of a Bayesian model (Gelman et al. 2014). In contrast with replicated outcome, predictive outcomes may use the values of independent variables that are different from those used to fit the model. Also see *simulated outcome*.
- **prior distribution, prior**. In Bayesian statistics, prior distributions are probability distributions of model parameters formed based on some a priori knowledge about parameters. Prior distributions are independent of the observed data.
- prior independence. See independent a priori.
- **prior odds**. Prior odds for θ_1 compared with θ_2 is the ratio of prior density evaluated at θ_1 and θ_2 under a given model, $p(\theta_1)/p(\theta_2)$. Also see *posterior odds*.
- **prior predictive distribution**. Prior predictive distribution is a distribution of unobserved (future) data derived by marginalizing the likelihood function with respect to the prior distribution of model parameters. Also see *marginal distribution*.
- **prior tightness**. A prior tightness is controlled by a tightness parameter, which is typically a multiplier for the prior variance. The smaller the value of this parameter, the smaller the prior variance, and the "tighter" (more highly concentrated) the prior around the prior mean. See [BAYES] **bayes: var**.
- **probability of unit circle inclusion**. In the context of Bayesian VAR, this is a posterior probability that all moduli of eigenvalues of a companion matrix lie within the unit circle. The higher this probability, the more likely the stability condition is met for the considered Bayesian VAR model.
- **proposal distribution**. In the context of the MH algorithm, a proposal distribution is used for defining the transition steps of the Markov chain. In the standard random-walk Metropolis algorithm, the proposal distribution is a multivariate normal distribution with zero mean and adaptable covariance matrix.
- **pseudoconvergence**. A Markov chain may appear to converge when in fact it did not. We refer to this phenomenon as pseudoconvergence. Pseudoconvergence is typically caused by multimodality of the stationary distribution, in which case the chain may fail to traverse the weakly connected regions of the distribution space. A common way to detect pseudoconvergence is to run multiple chains using different starting values and to verify that all of the chain converge to the same target distribution.

random effects. See random-effects parameters.

- random-effects linear form. A linear form representing a random-effects variable that can be used in substitutable expressions.
- **random-effects parameters**. In the context of Bayesian multilevel models, random-effects parameters are parameters associated with a random-effects variable. Random-effects parameters are assumed to be conditionally independent across levels of the random-effects variable given all other model parameters. Often, random-effects parameters are assumed to be normally distributed with a zero mean and an unknown variance–covariance matrix.
- **random-effects variable**. A variable identifying the group structure for the random effects at a specific level of hierarchy.
- reference prior. See noninformative prior.
- **replicated data**. Replicated data, **y**^{rep}, are data that could be observed if the experiment that produced the observed data, **y**^{obs}, were replicated using the same model and the same values of independent variables that generated **y**^{obs}. See Gelman et al. (2014, 145), [BAYES] **bayespredict**, and [BAYES] **bayesstats ppvalues**.
- **replicated outcome**. Replicated outcome is a special case of a simulated outcome that is generated using the same values of independent variables as those used to fit the model. Also see *replicated data*.
- scalar model parameter. A scalar model parameter is any model parameter that is a scalar. For example, {mean} and {hape:alpha} are scalar parameters, as declared by the bayesmh command. Elements of matrix model parameters are viewed as scalar model parameters. For example, for a 2 × 2 matrix parameter {Sigma,matrix}, individual elements {Sigma_1_1}, {Sigma_2_1}, {Sigma_1_2}, and {Sigma_2_2} are scalar parameters. If a matrix parameter contains a label, the label should be included in the specification of individual elements as well. See [BAYES] bayesmh.

scalar parameter. See scalar model parameter.

self-variable. A self-variable is a Stata term to refer to the dependent variable used as a lagged regressor in the VAR model in its own outcome equation. For instance, in a VAR model with two dependent variables (y1 and y2) and two lags,

$$\begin{split} \mathbf{y1} &= a_{11}\mathbf{L}.\mathbf{y1} + a_{12}\mathbf{L}2.\mathbf{y1} + a_{21}\mathbf{L}.\mathbf{y2} + a_{22}\mathbf{L}2.\mathbf{y2} + a_0 + u_1 \\ \mathbf{y2} &= b_{11}\mathbf{L}.\mathbf{y1} + b_{12}\mathbf{L}2.\mathbf{y1} + b_{21}\mathbf{L}.\mathbf{y2} + b_{22}\mathbf{L}2.\mathbf{y2} + b_0 + u_2 \end{split}$$

y1 is the self-variable in the first outcome equation (y1), and y2 is the self-variable in the second outcome equation (y2). Note that y2 and y1 are cross-variables in the first and second equations, respectively. Self-variables are always endogenous variables. We also often refer to the coefficients of self-variables as self-variable first-lag coefficients (a_{11} and b_{21}), self-variable second-lag coefficients (a_{12} and b_{22}), and so on. See [BAYES] bayes: var.

- self-variable first-lag coefficients. Self-variable first-lag coefficients are regression coefficients in a VAR model that correspond to first lags ([U] 11.4.4 Time-series varlists) of self-variables. See [BAYES] bayes: var.
- self-variable tightness parameter. A self-variable tightness parameter is a parameter, λ_1 , that controls the tightness of the Minnesota prior distribution by controlling the prior variance for the self-variable coefficients. It is specified in the Minnesota prior option selftight(). See Methods and formulas of [BAYES] bayes: var for details.

- **semiconjugate prior**. A prior distribution is semiconjugate for a family of likelihood distributions if the prior and (full) conditional posterior distributions belong to the same family of distributions. For semiconjugacy to hold, parameters must typically be independent a priori; that is, their joint prior distribution must be the product of the individual marginal prior distributions. For example, the normal prior distribution for a mean parameter of a normal data distribution with an unknown variance (which is assumed to be independent of the mean a priori) is a semiconjugate prior. Semiconjugacy may provide an efficient way of sampling from posterior distributions and is used in Gibbs sampling.
- **simulated outcome**. In Bayesian predictive inference, simulated outcomes are samples from the posterior predictive distribution. In the context of bayespredict, we define a simulated outcome as a $T \times n$ matrix of new outcome values simulated from the posterior predictive distribution, $p(\tilde{\mathbf{y}}|\mathbf{y})$, for a particular outcome variable \mathbf{y} , where T is the MCMC sample size and n is the number of observations.
- **spike-and-slab prior**. A spike-and-slab prior is one type of prior used for regression coefficients in Bayesian variable selection. The prior for coefficient β_j is a mixture of two normal distributions with different standard deviations or a mixture of two Laplace distributions with different scales. The mixing components γ_j 's are independent Bernoulli random variables with probability of success θ , where θ has a beta prior distribution. See *Spike-and-slab priors* in *Methods and formulas* of [BAYES] **bayess-elect**.
- **stationary distribution**. Stationary distribution of a stochastic process is a joint distribution that does not change over time. In the context of MCMC, stationary distribution is the target probability distribution to which the Markov chain converges. When MCMC is used for simulating a Bayesian model, the stationary distribution is the target joint posterior distribution of model parameters.
- subjective prior. See informative prior.
- subsampling the chain. See thinning.
- **sufficient statistic**. Sufficient statistic for a parameter of a parametric likelihood model is any function of the sample that contains all the information about the model parameter.
- test quantity. In Bayesian predictive inference, test quantity is any function of a simulated outcome, \mathbf{y}^{sim} , and model parameters $\boldsymbol{\theta}$. It is estimated by sampling from the joint posterior distribution $p(\mathbf{y}^{\text{sim}}, \boldsymbol{\theta})$. A test quantity that depends only on \mathbf{y}^{sim} is called a test statistic. Test quantities are used in posterior predictive checking to assess model fit.
- test statistic. A special case of a test quantity that depends only on the data.
- **thinning**. Thinning is a way of reducing autocorrelation in the MCMC sample by subsampling the MCMC chain every prespecified number of iterations determined by the thinning interval. For example, the thinning interval of 1 corresponds to using the entire MCMC sample; the thinning interval of 2 corresponds to using every other sample value; and the thinning interval of 3 corresponds to using values from iterations 1, 4, 7, 10, and so on. Thinning should be applied with caution when used to reduce autocorrelation because it may not always be the most appropriate way of improving the precision of estimates.

tightness. See prior tightness.

tightness parameter. A tightness parameter is a parameter, typically a multiplier for the prior variance, that controls the tightness of the prior. The smaller the value of this parameter, the smaller the prior variance, and the "tighter" (more highly concentrated) the prior around the prior mean. See *self-variable tightness parameter*, cross-variable tightness parameter, lad-decay parameter, and exogenous variable tightness parameter. Also see Methods and formula in [BAYES] **bayes: var**.

vague prior. See noninformative prior.

valid initial state. See feasible initial value.

vanishing adaptation. See diminishing adaptation.

VAR, vector autoregression. See VAR in [TS] Glossary.

variable selection. See Bayesian variable selection.

Zellner's g-prior. Zellner's *g*-prior is a form of a weakly informative prior for the regression coefficients in a linear model. It accounts for the correlation between the predictor variables and controls the impact of the prior of the regression coefficients on the posterior with parameter *g*. For example, g = 1 means that prior weight is 50% and $g \rightarrow \infty$ means diffuse prior.

Reference

Gelman, A., J. B. Carlin, H. S. Stern, D. B. Dunson, A. Vehtari, and D. B. Rubin. 2014. *Bayesian Data Analysis.* 3rd ed. Boca Raton, FL: Chapman and Hall/CRC.

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