

# 1 Model and Likelihood

Consider a linear  $K$ -factor model for  $D$  assets of the form

$$y_{it} = \alpha_d + \mathbf{f}_t' \boldsymbol{\beta}_d + \varepsilon_{it}$$

where  $d = 1, \dots, D$  and  $t = 1, \dots, T$  and  $\mathbf{f}_t' = (f_{t1}, \dots, f_{tK})$  is a  $K \times 1$  vector. This is a special case of the seemingly unrelated regression (SUR) model in which the regressors are *identical* across equations. Stacking observations for a given time period across assets, define  $\mathbf{y}_t' = (y_{1t}, \dots, y_{Dt})$  and analogously  $\boldsymbol{\varepsilon}_t' = (\varepsilon_{t1}, \dots, \varepsilon_{tD})$ . Now let  $\mathbf{x}_t' = (1, \mathbf{f}_t')$  and  $\boldsymbol{\gamma}_d' = (\alpha_d, \boldsymbol{\beta}_d')$  so we have

$$\mathbf{y}_t = X_t \boldsymbol{\gamma} + \boldsymbol{\varepsilon}_t$$

where  $X_t = I_D \otimes \mathbf{x}_t'$  and  $\boldsymbol{\gamma}' = (\boldsymbol{\gamma}_1', \dots, \boldsymbol{\gamma}_D')$ . Now, suppose that

$$\boldsymbol{\varepsilon}_t | \mathbf{x}_t \sim \text{iid } \mathcal{N}_D(0, \Omega)$$

Let  $Y_T$  denote the full data sample, i.e.  $\{\mathbf{y}_t, \mathbf{x}_t\}_{t=1}^T$ . Then the likelihood is

$$\pi(Y_T | \boldsymbol{\gamma}, \Omega^{-1}) \propto |\Omega^{-1}|^{T/2} \exp \left[ -\frac{1}{2} \sum_{t=1}^T (\mathbf{y}_t - X_t \boldsymbol{\gamma})' \Omega^{-1} (\mathbf{y}_t - X_t \boldsymbol{\gamma}) \right]$$

where we parameterize this problem in terms of the  $D \times D$  *precision* matrix  $\Omega^{-1}$  and the  $p \times 1$  vector of regression coefficients  $\boldsymbol{\gamma}$ , where  $p = D(K + 1)$ .

# 2 Prior and Posterior Distribution

To complete the model we specify the following prior distribution

$$\pi(\boldsymbol{\gamma}, \Omega^{-1}) = \mathcal{N}_p(\boldsymbol{\gamma} | \boldsymbol{\gamma}_0, G_0) \mathcal{W}_D(\Omega^{-1} | \rho_0, R_0)$$

This prior is conditionally conjugate with the normal likelihood. In particular, we have  $\boldsymbol{\gamma} | \Omega^{-1}, Y_T \sim \mathcal{N}_p(\bar{\boldsymbol{\gamma}}, G_T)$  where

$$\begin{aligned} G_T &= \left[ G_0^{-1} + \sum_{t=1}^T X_t' \Omega^{-1} X_t \right]^{-1} \\ \bar{\boldsymbol{\gamma}} &= G_T \left[ G_0^{-1} \boldsymbol{\gamma}_0 + \sum_{t=1}^T X_t' \Omega^{-1} \mathbf{y}_t \right] \end{aligned}$$

and  $\Omega^{-1}|Y_T \sim \mathcal{W}_D(\rho_0 + T, R_T)$  where

$$R_T = \left[ R_0^{-1} + \sum_{t=1}^T (\mathbf{y}_t - X_t \boldsymbol{\gamma}) (\mathbf{y}_t - X_t \boldsymbol{\gamma})' \right]^{-1}$$

### 3 MCMC

Using the full set of conditional posteriors, given in the preceding section, we can simulate from the joint posterior for this model using a Gibbs sampler:

1. Select a starting value  $\Omega^{-1(0)}$  for the precision matrix.
2. Draw  $\boldsymbol{\gamma}^{(1)} \sim \mathcal{N}(\bar{\boldsymbol{\gamma}}^{(1)}, G_T^{(1)})$  where

$$\begin{aligned} G_T^{(1)} &= \left[ G_0^{-1} + \sum_{t=1}^T X_t' \Omega^{-1(0)} X_t \right]^{-1} \\ \bar{\boldsymbol{\gamma}}^{(1)} &= G_T^{(1)} \left[ G_0^{-1} \boldsymbol{\gamma}_0 + \sum_{t=1}^T X_t' \Omega^{-1(0)} \mathbf{y}_t \right] \end{aligned}$$

3. Draw  $\Omega^{-1(1)} \sim \mathcal{W}_D(\rho_T, R_T^{(1)})$  where

$$R_T^{(1)} = \left[ R_0^{-1} + \sum_{t=1}^T (\mathbf{y}_t - X_t \boldsymbol{\gamma}^{(1)}) (\mathbf{y}_t - X_t \boldsymbol{\gamma}^{(1)})' \right]^{-1}$$

4. Repeat the preceding two steps a total of  $G$  times. In the  $g$ th iteration:

- (i) Draw  $\boldsymbol{\gamma}^{(g)} \sim \mathcal{N}(\bar{\boldsymbol{\gamma}}^{(g)}, G_T^{(g)})$  where

$$\begin{aligned} G_T^{(g)} &= \left[ G_0^{-1} + \sum_{t=1}^T X_t' \Omega^{-1(g-1)} X_t \right]^{-1} \\ \bar{\boldsymbol{\gamma}}^{(g)} &= G_T^{(g)} \left[ G_0^{-1} \boldsymbol{\gamma}_0 + \sum_{t=1}^T X_t' \Omega^{-1(g-1)} \mathbf{y}_t \right] \end{aligned}$$

(ii) Draw  $\Omega^{-1(g)} \sim \mathcal{W}_D \left( \rho_T, R_T^{(g)} \right)$  where

$$R_T^{(g)} = \left[ R_0^{-1} + \sum_{t=1}^T (\mathbf{y}_t - X_t \boldsymbol{\gamma}^{(g)}) (\mathbf{y}_t - X_t \boldsymbol{\gamma}^{(g)})' \right]^{-1}$$

5. Discard the first  $B$  draws.

Note that in iteration  $g$ ,  $G_T^{(g)}$  and  $\tilde{\boldsymbol{\gamma}}^{(g)}$  are calculated using  $\Omega^{-1(g-1)}$  while  $R_T^{(g)}$  is calculated using  $\boldsymbol{\gamma}^{(0)}$ . This is because we choose to initialize the sample with a starting value  $\Omega^{-1(0)}$  for the precision matrix rather than for the vector of regression coefficients.

## 4 Numerical Details

### 4.1 Evaluating the MV Normal Density

As one of the steps in the calculation of the marginal likelihood (see below) we will need to repeatedly evaluate the log of a multivariate normal density at a fixed set of parameter values. Let  $Z$  be a  $p \times n$  matrix, each of whose columns is a point  $\mathbf{z}$  at which we wish to evaluate  $\log \mathcal{N}_p(\mathbf{z}|\mu, \Sigma)$  where  $\mu$  is the mean vector and  $\Sigma$  the covariance matrix of a multivariate normal. Because our problem is parameterizes in terms of the *precision* matrix rather than the covariance matrix, the calculations given here assume that we are given  $\Sigma^{-1}$  rather than  $\Sigma$ . In terms of the precision matrix, the log of the MV normal density is given by

$$\log \mathcal{N}_p(\mathbf{z}|\mu, \Sigma^{-1}) = -\frac{p}{2} \log(2\pi) + \frac{1}{2} \log |\Sigma^{-1}| - \frac{1}{2} (\mathbf{z} - \mu)' \Sigma^{-1} (\mathbf{z} - \mu)$$

Now let  $R$  be the Cholesky factor of  $\Sigma^{-1}$  so that  $\Sigma^{-1} = R'R$  and define  $\tilde{\mathbf{z}} = \mathbf{z} - \mu$  and  $\mathbf{v} = R\tilde{\mathbf{z}}$ . Using these definitions,

$$(\mathbf{z} - \mu)' \Sigma^{-1} (\mathbf{z} - \mu) = (R\tilde{\mathbf{z}})'(R\tilde{\mathbf{z}}) = \mathbf{v}'\mathbf{v}$$

and, letting  $R_{ii}$  denote the  $i$ th diagonal element of  $R$ ,

$$\begin{aligned}\log |\Sigma^{-1}| &= \log |R'R| = \log (|R'| \cdot |R|) = \log |R'| + \log |R| \\ &= 2 \log |R| = 2 \cdot \sum_{i=1}^p \log R_{ii}\end{aligned}$$

since  $|A| = |A'|$ ,  $|AB| = |A| \cdot |B|$  and the determinant of a triangular matrix equals the product of its diagonal elements. Thus, we have

$$\log \mathcal{N}_p(\mathbf{z}|\mu, \Sigma^{-1}) = -\frac{p}{2} \log(2\pi) + \text{trace}[\log(\text{diag}\{R\})] - \frac{1}{2} \mathbf{v}'\mathbf{v}$$

The only term in the preceding expression that depends on  $\mathbf{z}$  is  $\mathbf{v}'\mathbf{v}$ . We can calculate this term simultaneously for all columns of  $Z$  as follows. First let  $\tilde{Z}$  denote the result of subtracting the vector  $\mu$  from each column of  $Z$ , i.e.  $\tilde{Z} = Z - \mu \mathbf{1}'_n$ . To calculate  $\mathbf{v}'\mathbf{v}$  for each column of  $Z$  we simply square the elements of  $R\tilde{Z}$  and take the column sums of the resulting matrix.

## 4.2 Efficient Calculation of $R_T$

In the second step of each iteration we compute  $\left(R_0^{-1} + \sum_{t=1}^T \hat{\mathbf{e}}_t \hat{\mathbf{e}}_t'\right)^{-1}$  where  $\hat{\mathbf{e}}_t = \mathbf{y}_t - X_t \boldsymbol{\gamma}$ . Since  $R_0$  is simply the prior scale matrix for  $\Omega^{-1}$  and hence remains unchanged during the iterations, we can pre-compute it and store the result before starting the sampler. Since  $X_t$  is a sparse matrix, there is a much more efficient and compact way to compute the sum of outer products of residuals. Define:

$$\tilde{Y} = \begin{bmatrix} \mathbf{y}'_1 \\ \vdots \\ \mathbf{y}'_T \end{bmatrix}, \quad \tilde{X} = \begin{bmatrix} \mathbf{x}'_1 \\ \vdots \\ \mathbf{x}'_T \end{bmatrix}, \quad \Gamma = \begin{bmatrix} \gamma_1 & \cdots & \gamma_D \end{bmatrix}, \quad \hat{\mathbf{e}} = \begin{bmatrix} \hat{\mathbf{e}}'_1 \\ \vdots \\ \hat{\mathbf{e}}'_T \end{bmatrix}$$

so that  $\hat{\mathbf{e}} = \tilde{Y} - \tilde{X}\Gamma$ . Note that the vector of regression coefficients  $\boldsymbol{\gamma}$  is the vec of the *matrix* of regression coefficients  $\Gamma$ . Thus, expressed in terms of dense matrix operations

$$R_T^{-1} = R_0^{-1} + (\tilde{Y} - \tilde{X}\Gamma)'(\tilde{Y} - \tilde{X}\Gamma)$$

The final step is to invert this sum (which is positive definite) to calculate  $R_T$ . Note that the Matrix Inversion Lemma (Sherman-Morrison-Woodbury Formula) does *not* simplify this calculation unless  $D > T$ .

### 4.3 Efficient Calculation of $G_T$

Because we parameterize our multivariate normal sampler in terms of the *precision* matrix rather than the covariance matrix, we work with the *inverse* of  $G_T$ , namely

$$G_T^{-1} = G_0^{-1} + \sum_{t=1}^T X_t' \Omega^{-1} X_t$$

Since it is simply the prior precision matrix for the vector  $\gamma$  of regression coefficients we can pre-compute  $G_0$  (assuming that we elicit a prior in terms of the covariance matrix). Now, the sum over  $X_t' \Omega^{-1} X_t$  can in fact be simplified using the properties of the Kronecker product.<sup>1</sup> Recall that  $X_t = I_D \otimes \mathbf{x}_t'$ . Since  $(A \otimes B)' = A' \otimes B'$ ,

$$X_t' \Omega^{-1} X_t = (I_D \otimes \mathbf{x}_t) \Omega^{-1} X_t$$

Since  $\Omega^{-1} X_t = (\Omega^{-1} X_t) \otimes 1$ ,  $\Omega^{-1} = \Omega^{-1} \otimes 1$ , and  $(A \otimes B)(C \otimes D) = AC \otimes BD$ , provided that everything is conformable, we have

$$\begin{aligned} (I_D \otimes \mathbf{x}_t) \Omega^{-1} X_t &= (I_D \otimes \mathbf{x}_t) (\Omega^{-1} X_t \otimes 1) \\ &= \Omega^{-1} X_t \otimes \mathbf{x}_t = [\Omega^{-1} (I_D \otimes \mathbf{x}_t')] \otimes \mathbf{x}_t \\ &= [(\Omega^{-1} \otimes 1) (I_D \otimes \mathbf{x}_t')] \otimes \mathbf{x}_t \\ &= \Omega^{-1} \otimes \mathbf{x}_t' \otimes \mathbf{x}_t \end{aligned}$$

Finally, since  $A \otimes (B + C) = A \otimes B + A \otimes C$ ,

$$\sum_{t=1}^T X_t' \Omega^{-1} X_t = \sum_{t=1}^T \Omega^{-1} \otimes \mathbf{x}_t' \otimes \mathbf{x}_t = \Omega^{-1} \otimes \left( \sum_{t=1}^T \mathbf{x}_t' \otimes \mathbf{x}_t \right)$$

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<sup>1</sup>See, e.g., Horn and Johnson (1994) Chapter 4.2.

This is an extremely useful simplification: because  $\sum_{t=1}^T \mathbf{x}'_t \otimes \mathbf{x}_t$  involves neither  $\Omega^{-1}$  nor  $\boldsymbol{\gamma}$ , only the data, we can pre-compute this quantity. In fact, there is one final simplification that makes this quantity even simpler. By writing out the definition of the Kronecker Product, we see that  $\mathbf{x}'_t \otimes \mathbf{x}_t = \mathbf{x}_t \mathbf{x}'_t$  and hence

$$\sum_{t=1}^T X'_t \Omega^{-1} X_t = \Omega^{-1} \otimes \left( \sum_{t=1}^T \mathbf{x}_t \mathbf{x}'_t \right) = \Omega^{-1} \otimes \tilde{X}' \tilde{X}$$

where  $\tilde{X}' = \begin{bmatrix} \mathbf{x}_1 & \cdots & \mathbf{x}_T \end{bmatrix}$ . Thus, we have  $G_T^{-1} = G_0^{-1} + \Omega^{-1} \otimes \tilde{X}' \tilde{X}$ .

#### 4.4 Efficient Calculation of $\bar{\boldsymbol{\gamma}}$

The vector  $\bar{\boldsymbol{\gamma}}$  is constructed from several pieces. The first is  $G_0^{-1} \boldsymbol{\gamma}_0$ , the solution to the linear system  $G_0 \mathbf{v} = \boldsymbol{\gamma}_0$ . Since this piece depends only on the prior, we can pre-compute it. The next piece is the sum  $\sum_{t=1}^T X'_t \Omega^{-1} \mathbf{y}_t$ . We noted above,  $X_t$  is sparse so there is a more efficient way to compute this quantity. Indeed, while this is far from obvious at first glance, it is possible to *factor*  $\Omega^{-1}$  outside of the sum using some clever matrix operations, allowing us to drastically reduce the computational complexity of the sampler. To accomplish this simplification we combine the definition of  $X_t$  as  $I_D \otimes \mathbf{x}'_t$  with two properties of the Kronecker Product, namely:

$$(A \otimes B) (C \otimes D) = AC \otimes BD$$

which holds provided that the respective matrices are conformable and

$$\text{vec}(AB) = (B' \otimes I_k) \text{vec}(A)$$

where  $A$  is  $k \times \ell$  and  $B$  is  $\ell \times m$ . Applying the first property twice in succession followed by the second property, we find that

$$\begin{aligned}
X_t' \Omega^{-1} \mathbf{y}_t &= (I_D \otimes \mathbf{x}_t')' \Omega^{-1} \mathbf{y}_t = (I_D \otimes \mathbf{x}_t) \Omega^{-1} \mathbf{y}_t \\
&= (I_D \otimes \mathbf{x}_t) (\Omega^{-1} \mathbf{y}_t \otimes 1) = I_D \Omega^{-1} \mathbf{y}_t \otimes \mathbf{x}_t 1 \\
&= \Omega^{-1} \mathbf{y}_t \otimes \mathbf{x}_t = [(\Omega^{-1} \mathbf{y}_t) \ 1] \otimes [I_{K+1} \mathbf{x}_t] \\
&= (\Omega^{-1} \mathbf{y}_t \otimes I_{K+1}) (1 \otimes \mathbf{x}_t) \\
&= ([\Omega^{-1} \mathbf{y}_t] \otimes I_{K+1}) \mathbf{x}_t = \left( [\mathbf{y}_t' \Omega^{-1}]' \otimes I_{K+1} \right) \text{vec}(\mathbf{x}_t) \\
&= \text{vec}(\mathbf{x}_t \mathbf{y}_t' \Omega^{-1})
\end{aligned}$$

where we have used the fact that  $\text{vec}(\mathbf{x}_t) = \mathbf{x}_t$ . Finally, since we can interchange the vec summation operations,

$$\begin{aligned}
\sum_{t=1}^T \text{vec}(\mathbf{x}_t \mathbf{y}_t' \Omega^{-1}) &= \text{vec} \left[ \sum_{t=1}^T \mathbf{x}_t \mathbf{y}_t' \Omega^{-1} \right] = \text{vec} \left[ \left( \sum_{t=1}^T \mathbf{x}_t \mathbf{y}_t' \right) \Omega^{-1} \right] \\
&= \text{vec}(\tilde{X}' \tilde{Y} \Omega^{-1})
\end{aligned}$$

where, as above,

$$\tilde{Y} = \begin{bmatrix} \mathbf{y}_1' \\ \vdots \\ \mathbf{y}_T' \end{bmatrix}, \quad \tilde{X} = \begin{bmatrix} \mathbf{x}_1' \\ \vdots \\ \mathbf{x}_T' \end{bmatrix}$$

Thus we see that

$$\bar{\boldsymbol{\gamma}} = G_T \left[ G_0^{-1} \boldsymbol{\gamma}_0 + \text{vec}(\tilde{X}' \tilde{Y} \Omega^{-1}) \right]$$

Because it does not change between iterations, we can pre-compute the product  $\tilde{X}' \tilde{Y}$ . The only term that remains to be addressed is  $G_T$ . Because our normal sampler is parameterized in terms of the precision matrix rather than the covariance matrix we calculated  $G_T^{-1}$  rather than  $G_T$  above. Rather than inverting it in this step, which is a very bad idea given its size, we notice that our expression for  $\bar{\boldsymbol{\gamma}}$  takes the form  $\mathbf{v} = A^{-1} \mathbf{b}$ . Therefore,

$$\bar{\boldsymbol{\gamma}} = \text{solve} \left[ G_T^{-1}, G_0^{-1} \boldsymbol{\gamma}_0 + \text{vec}(\tilde{X}' \tilde{Y} \Omega^{-1}) \right]$$

## 5 Calculating the Marginal likelihood

We calculate the marginal likelihood using the method of Chib (1995). Let  $\theta$  denote the full collection of parameters. By Bayes' Rule

$$\pi(\theta|Y_T) = \frac{\pi(\theta)f(Y_T|\theta)}{f(Y_T)}$$

where  $f(Y_T)$  is the marginal likelihood, aka the marginal data density, aka the evidence. This identity holds true for *any* value of  $\theta$ . In particular it holds at the posterior mean  $\theta^*$ . Solving for  $f(Y_T)$  and evaluating the result at  $\theta^*$ , we have

$$f(Y_T) = \frac{\pi(\theta^*)f(Y_T|\theta^*)}{\pi(\theta^*|Y_T)}$$

Thus, we can express the *log* marginal likelihood as

$$\log f(Y_T) = \log \pi(\theta^*) + \log f(Y_T|\theta^*) - \pi(\theta^*|Y_T)$$

Specializing this to the SUR model considered above,

$$\log f(Y_T) = \log \pi(\boldsymbol{\gamma}^*) + \log \pi(\Omega^{-1*}) + \log f(Y_T|\boldsymbol{\gamma}^*, \Omega^{-1*}) - \pi(\boldsymbol{\gamma}^*, \Omega^{-1*}|Y_T)$$

since our priors over  $\boldsymbol{\gamma}$  and  $\Omega^{-1}$  are independent. The Chib (1995) method approximates  $\log f(Y_T)$  by evaluating each of the terms on the right-hand-side of the preceding expression using the output of the Gibbs sampler.

**The Contribution of the Prior** Evaluating the first two terms,  $\log \pi(\boldsymbol{\gamma}^*)$  and  $\log \pi(\Omega^{-1*})$ , is easy: these are simply the priors for  $\boldsymbol{\gamma}$  and  $\Omega^{-1}$  evaluated at the posterior means. We take the sample average of the Gibbs draws to approximate  $\boldsymbol{\gamma}^*$  and  $\Omega^{-1*}$  and evaluate the Normal and Wishart distributions at these points, with parameters given by the prior:

$$\begin{aligned} \pi(\boldsymbol{\gamma}^*) &= \mathcal{N}_p(\boldsymbol{\gamma}^*|\boldsymbol{\gamma}_0, G_0) \\ \pi(\Omega^{-1*}) &= \mathcal{W}_D(\Omega^{-1*}|\rho_0, R_0) \end{aligned}$$



**The Contribution of the Likelihood** Above we assumed a normal distribution for the regression errors, specifically,  $\boldsymbol{\varepsilon}_t | \mathbf{x}_t \sim \text{iid } \mathcal{N}_D(0, \Omega)$ . From the regression specification it follows that  $\mathbf{y}_t \sim \text{iid } \mathcal{N}_D(X_t \boldsymbol{\gamma}, \Omega)$  and thus the log likelihood evaluated at the posterior mean is

$$\log f(Y_T | \boldsymbol{\gamma}^*, \Omega^{-1*}) = \sum_{t=1}^T \log \mathcal{N}_D(\mathbf{y}_t | X_t \boldsymbol{\gamma}^*, \Omega^{-1*})$$

parameterized in terms of the precision matrix rather than the covariance matrix. Equivalently, but more conveniently, we may write

$$\log f(Y_T | \boldsymbol{\gamma}^*, \Omega^{-1*}) = \sum_{t=1}^T \log \mathcal{N}_D(\mathbf{y}_t - X_t \boldsymbol{\gamma}^* | \mathbf{0}, \Omega^{-1*})$$

The advantage of this version of the likelihood is that the parameters of the normal density are constant over  $t$ , allowing us to exploit the efficient algorithm for repeatedly evaluating a MV normal density with fixed parameters, described above. Note that we can simultaneously calculate all of the arguments for the normal density as  $\tilde{Y} - \tilde{X} \boldsymbol{\gamma}^*$  using the notation defined above in the description of efficient calculations for  $R_T$ .

**The Contribution of the Posterior** To evaluate the third term, we factorize the joint posterior as the product of a conditional and marginal, namely:

$$\pi(\boldsymbol{\gamma}^*, \Omega^{-1*} | Y_T) = \pi(\boldsymbol{\gamma}^* | \Omega^{-1*}, Y_T) \times \pi(\Omega^{-1*} | Y_T)$$

Because we have *analytical expressions* for the conditional posteriors in this model we can evaluate the first term in the product immediately. We have  $\boldsymbol{\gamma} | \Omega^{-1} \sim \mathcal{N}_p(\boldsymbol{\gamma} | \bar{\boldsymbol{\gamma}}, G_T)$  where  $G_T$  and  $\bar{\boldsymbol{\gamma}}$  depend only on the prior, the data, and  $\Omega^{-1}$ . To perform the required calculation, we simply evaluate the normal density at  $\boldsymbol{\gamma}^*$  and evaluate  $G_T$  and  $\bar{\boldsymbol{\gamma}}$  at  $\Omega^{-1*}$ , that is:

$$\pi(\boldsymbol{\gamma}^* | \Omega^{-1*}, Y_T) = \mathcal{N}_p(\boldsymbol{\gamma}^* | \bar{\boldsymbol{\gamma}}^*, G_T^*)$$

where

$$\begin{aligned} G_T^* &= \left[ G_0^{-1} + \sum_{t=1}^T X_t' \Omega^{-1*} X_t \right]^{-1} = \left[ G_0^{-1} + \Omega^{-1*} \otimes \tilde{X}' \tilde{X} \right]^{-1} \\ \bar{\gamma}^* &= G_T^* \left[ G_0^{-1} \gamma_0 + \sum_{t=1}^T X_t' \Omega^{-1*} \mathbf{y}_t \right] = G_T^* \left[ G_0^{-1} \gamma_0 + \text{vec} \left( \tilde{X}' \tilde{Y} \Omega^{-1*} \right) \right] \end{aligned}$$

The evaluation of the second term in the product that gives the contribution of the posterior to the marginal likelihood is a bit more involved. We write

$$\begin{aligned} \pi(\Omega^{-1*} | Y_T) &= \int \pi(\gamma, \Omega^{-1*} | Y_T) d\gamma \\ &= \int \pi(\Omega^{-1*} | \gamma, Y_T) \pi(\gamma | Y_T) d\gamma \end{aligned}$$

and approximate the second integral using the draws from the Gibbs sampler:

$$\begin{aligned} \pi(\Omega^{-1*} | Y_T) &\approx \frac{1}{G} \sum_{g=1}^G \pi(\Omega^{-1*} | \gamma^{(g)}, Y_T) \\ &= \frac{1}{G} \sum_{g=1}^G \mathcal{W}_D(\Omega^{-1*} | \rho_T, R_T^{(g)}) \end{aligned}$$

where

$$\begin{aligned} R_T^{(g)} &= \left[ R_0^{-1} + \sum_{t=1}^T (\mathbf{y}_t - X_t \gamma^{(g)}) (\mathbf{y}_t - X_t \gamma^{(g)})' \right]^{-1} \\ &= \left[ R_0^{-1} + (\tilde{Y} - \tilde{X} \Gamma^{(g)})' (\tilde{Y} - \tilde{X} \Gamma^{(g)}) \right]^{-1} \end{aligned}$$

## 6 Prediction

Suppose we are interested in predicting the cross-section of returns  $y_{n+1}$  at time  $(n+1)$ . The Bayes prediction density of these returns, conditioned on the data  $Y_{n+1}$  and the factors  $f_{n+1}$ , is given by

$$p(y_{n+1} | Y_n, f_{n+1}) = \int_{\gamma, \Omega^{-1}} \mathcal{N}_d(y_{n+1} | X_{n+1} \gamma, \Omega) d\pi(\gamma, \Omega^{-1} | Y_n)$$

which is estimated by the ergodic Monte Carlo average

$$p(y_{n+1}|Y_n, f_{n+1}) = \frac{1}{G} \sum_{g=1}^G \mathcal{N}_d(y_{n+1}|X_{n+1}\gamma^{(g)}, \Omega^{(g)})$$

with the MCMC draws  $\{\gamma^{(g)}, \Omega^{(g)}\}$  from the posterior distribution.

## 7 Student-t errors

Suppose now that the errors are distributed as multivariate-t

$$\varepsilon_t \sim t_{d,\nu}(0, \Omega)$$

so that

$$\begin{aligned} E(\varepsilon_t) &= 0, \quad \nu > 1 \\ Var(\varepsilon_t) &= \frac{\nu}{\nu - 2} \Omega, \quad \nu > 2 \end{aligned}$$

The analysis of this model utilizes the hierarchical representation

$$\begin{aligned} \varepsilon_t | \lambda_t &\sim N(0, \lambda_t^{-1} \Omega) \\ \lambda_t &\sim G\left(\frac{\nu}{2}, \frac{\nu}{2}\right) \end{aligned}$$

which means that conditioned on  $(\nu, \{\lambda_t\})$ , the results presented for the Gaussian model can be applied with minor modifications. The MCMC sampling is completed with the sampling of  $(\nu, \{\lambda_t\})$ .

Following Albert and Chib (1993), let us assume that the support of  $\nu$  is the set of values  $\{\nu_j\}_{j=1}^J$ , for example,  $\{4, 6, 8, 10, 12, 14, 16\}$  and that a priori

$$\Pr(\nu = \nu_j) = q_j$$

Then, simple calculations show that

$$\gamma|Y_n, \Omega^{-1}, \nu, \{\lambda_t\} \sim \mathcal{N}_{d+p}(\hat{\gamma}_\lambda, G_{n,\lambda})$$

where

$$\hat{\gamma}_\lambda = G_{n,\lambda} \left( G_0^{-1} \gamma_0 + \sum_{t=1}^n \lambda_t X_t' \Omega^{-1} y_t \right)$$

$$G_{n,\lambda} = \left( G_0^{-1} + \sum_{t=1}^n \lambda_t X_t' \Omega^{-1} X_t \right)^{-1}$$

and

$$\Omega^{-1} | Y_n, \gamma, \nu, \{\lambda_t\} \sim \mathcal{W}_d \left( \rho_0 + n, \left( R_0^{-1} + \sum_{t=1}^n \lambda_t (y_t - X_t \gamma) (y_t - X_t \gamma)' \right)^{-1} \right)$$

Moreover,

$$\Pr(\nu = \nu_j | Y_n, \gamma, \Omega^{-1}) \propto q_j \prod_{t=1}^n t_{d,\nu_j}(y_t | X_t \gamma, \Omega)$$

and

$$\lambda_t | Y_n, \gamma, \nu \sim G \left( \frac{\nu + d}{2}, \frac{\nu + (y_t - X_t \gamma)' (y_t - X_t \gamma)}{2} \right)$$

One sweep of the MCMC sampling is completed by sampling these four distributions in this order.

## 7.1 Marginal Likelihood

The Chib (1995) method can again be applied to find the log marginal likelihood as

$$\log \Pr(\nu^*) + \log \pi(\gamma^*) + \log \pi(\Omega^{-1*}) + \sum_{t=1}^n \log t_{d,\nu^*}(y_t | X_t \gamma^*, \Omega^*) - \log \pi(\nu^*, \gamma^*, \Omega^{-1*} | Y_n)$$

where  $\nu^*$  is the posterior mode (which is easily computed from the sampled values), the last term is calculated as

$$\Pr(\nu^* | Y_n) \times \pi(\Omega^{-1*} | Y_n, \nu^*) \times \pi(\gamma^* | Y_n, \Omega^{-1*}, \nu^*)$$

in which the first term is obtained from the posterior frequency distribution of  $\nu$ , the second term is obtained from a reduced run in which  $\nu$  is fixed at  $\nu^*$  and the remaining three distributions are sampled and the draws

$$\left\{ \gamma^{(g)}, \lambda_t^{(g)} \right\}_{g=1}^G$$

from this reduced MCMC run are used to calculate  $\pi(\Omega^{-1*}|Y_n, \nu^*)$  as

$$\frac{1}{G} \sum_{g=1}^G \mathcal{W}_d \left( \Omega^{-1*} | \rho_0 + n, \left( R_0^{-1} + \sum_{t=1}^n \lambda_t^{(g)} (y_t - X_t \gamma^{(g)}) (y_t - X_t \gamma^{(g)})' \right)^{-1} \right)$$

and the final term  $\pi(\gamma^*|Y_n, \Omega^{-1*})$  is obtained from a second reduced run in which  $\nu$  is fixed at  $\nu^*$  and  $\Omega^{-1}$  is fixed at  $\Omega^{-1*}$  and the draws

$$\left\{ \lambda_t^{(g)} \right\}$$

from this reduced run are used to give

$$\pi(\gamma^*|Y_n, \Omega^{-1*}, \nu^*) = \frac{1}{G} \sum_{g=1}^G \mathcal{N}_{d+p} \left( \gamma^* | \hat{\gamma}_{\lambda^{(g)}}^*, G_{n, \lambda^{(g)}}^* \right)$$

where

$$\begin{aligned} \hat{\gamma}_{\lambda^{(g)}}^* &= G_{n, \lambda^{(g)}}^* \left( G_0^{-1} \gamma_0 + \sum_{t=1}^n \lambda_t^{(g)} X_t' \Omega^{-1*} y_t \right) \\ G_{n, \lambda^{(g)}}^* &= \left( G_0^{-1} + \sum_{t=1}^n \lambda_t^{(g)} X_t' \Omega^{-1*} X_t \right)^{-1} \end{aligned}$$