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# A Partially Confirmatory Approach to Scale Development With the Bayesian Lasso

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#### Abstract

The exploratory and confirmatory approaches of factor analysis lie on two ends of a continuum of substantive input for scale development. Recent advancements in Bayesian regularization methods enable more flexibility in covering a wide range of the substantive continuum. Based on the Bayesian Lasso (least absolute shrinkage and selection operator) methods for the regression model and covariance matrix, this research proposes a partially confirmatory approach to address the loading and residual structures at the same time. With at least one specified loading per item, a one-step procedure can be applied to figure out both structures simultaneously. With a few specified loadings per factor, a two-step procedure is preferred to capture the model configuration correctly. In both cases, the Bayesian hierarchical formulation is implemented using Markov Chain Monte Carlo estimation with different Lasso or regular priors. Both simulated and real data sets were analyzed to evaluate the validity, robustness, and practical usefulness of the proposed approach across different situations.

#### Translational Abstract

Exploratory factor analysis (EFA) and confirmatory factor analysis (CFA) are traditional approaches for development of psychological scales, and each has pros and cons. EFA is driven by data, whereas CFA is driven by theory. However, different amounts of substantive knowledge are available to analyze the factor structure of a scale, so the approaches become two ends of a continuum. Although it is desirable to cover a wide range of the continuum in practice, there are psychometric challenges in combining EFA and CFA within one framework. Relying on Bayesian Lasso methods and sound statistical modeling, this research proposes a partially confirmatory factor analysis (PCFA) approach to address the challenges. With a two-step procedure, the proposed approach can offer more flexibility to scale development, especially when there are multiple factors and many items. Using both simulation experiments and real data sets related to mathematics learning and humor style questionnaires, this article shows that PCFA can help researchers better develop scales with both exploratory and confirmatory elements across different settings.

Keywords: factor analysis, Bayesian Lasso, partially confirmatory, Lasso loading, residual covariance

Factor analysis (FA) is a statistical technique widely used for developing psychological scales with multiple factors and many items. The major purpose is to explain item responses in term of

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Correspondence concerning this article should be addressed to Jinsong Chen or Junhao Pan, Department of Psychology, Sun Yat-Sen University, 132 Waihuan Dong Road, Higher Education Mega Center, Guangzhou 510006, China. E-mail: jinsong.chen@live.com or panjunh@mail.sysu .edu.cn the structure of the factorial loadings and residuals. With FA, there are two typical approaches of scale development: exploratory (EFA; Jennrich & Sampson, 1966) and confirmatory (CFA; Jöreskog, 1969). EFA is data driven with little substantive knowledge or constraint on the loading structure, whereas CFA is theory driven with strong knowledge or constraint on the structure. In general, however, different amounts of substantive input can be available, making the two approaches two ends of a continuum. To be more flexible on the substantive continuum, one traditional practice is to apply post hoc model modification (PMM) based on modification indices (Kaplan, 1990; Sörbom, 1989) to free up constraints on the basis of the confirmatory approach. However, PMM is theoretically suboptimal and will be problematic when many constraints need to be released (Bentler, 2007; MacCallum, 1986; Steiger, 1990).

With the recent development of Bayesian regularization methods such as Bayesian ridge regression (Hsiang, 1975), Bayesian least absolute shrinkage and selection operator or Lasso (Hans, 2009; Park & Casella, 2008), and spike-and-slab prior (O'Hara & Sillanpää, 2009), it is increasingly possible to achieve more flexibility on the substantive continuum in scale development using FA. Specifically, cross-loadings can be regularized via the Bayesian structural equation modeling (BSEM) approach with the ridge regression prior (B. O. Muthén & Asparouhov, 2012) or spikeand-slab prior (Lu, Chow, & Loken, 2016). Regularization of off-diagonal elements in the residual variance-covariance matrix can be conducted via the Bayesian covariance Lasso for CFA (BLCCFA) approach (Pan, Ip, & Dubé, 2017). Studies have also demonstrated the superiority of Bayesian methods over PMM for identifying significant cross-loadings or residual covariance in corresponding settings. The methods all offer variable selection with different norms of penalty on the parameters of interests, Bayesian hierarchical modeling with the Markov Chain Monte Carlo (MCMC; Gilks, Richardson, & Spiegelhalter, 1996) algorithm, and simultaneous estimation of the shrinkage parameters with proper priors.

On the basis of previous literature, this research proposes a partially confirmatory factor analysis (PCFA) approach with the Bayesian Lasso method to address the loading and residual covariance structures. Addressing both structures at the same time is challenging, and some constraints can be used. With the constraint of one specified loading per item, a one-step procedure can be applied to capture the loading and residual covariance structures simultaneously. Otherwise, an additional step with a few specified loadings per factor and diagonal residual matrix can be used to satisfy the constraints before applying the above step, making it a two-step procedure. In both cases, the MCMC estimation is implemented based on derivations with different Lasso or regular priors on the loading and residual structures. Both the Gibbs sampler (Casella & George, 1992; Geman & Geman, 1984) and Metropolis-Hastings algorithm (Hastings, 1970; Metropolis, Rosenbluth, Rosenbluth, Teller, & Teller, 1953) are used in the MCMC estimation. Note that the one-step procedure reduces to the BLCCFA approach when all loadings are specified or fixed, whereas the first step in the two-step procedure is similar to the BSEM approach but with different regularization methods. With the two procedures, the proposed approach can cover a wide range of the substantive continuum and be flexible and useful for scale development, especially with three or more factors and many items. Both simulated and real data sets were analyzed to evaluate the validity, robustness, and practical usefulness of the proposed approach across different situations.

#### The Partially Confirmatory Framework

Suppose there are *J* items in a psychological scale with *N* respondents, and item responses  $\mathbf{y}_1, \mathbf{y}_2, \ldots, \mathbf{y}_N$  are independently and randomly distributed observations. In a general FA model, each  $\mathbf{y}_i = (\mathbf{y}_{i1}, \mathbf{y}_{i2}, \ldots, \mathbf{y}_{iJ})^T$  satisfies the following equation:

$$\mathbf{y}_i = \mathbf{\mu} + \mathbf{\Lambda} \boldsymbol{\omega}_i + \boldsymbol{\varepsilon}_i, i = 1, 2, \dots, N, \tag{1}$$

where model parameters include the  $J \times 1$  intercept vector  $\mu$ (usually of little interest during scale development),  $J \times K$  loading matrix  $\Lambda = (\lambda_{jk})$ ,  $K \times K$  factorial covariance matrix  $\Phi$ , and  $J \times$ J residual covariance matrix  $\Psi = (\psi_{jj'})$ , with  $K \times 1$  latent factors  $\omega_i \sim N_K[0, \Phi]$  and  $J \times 1$  residuals  $\varepsilon_i \sim N_J[0, \Psi]$ . To determine the scale of the latent factors, one can either fix the factorial variance or fix one loading per factor. In this research, all factorial variances are fixed to one, which means  $\Phi$  is a correlation matrix. Loadings in  $\Lambda$  can be fixed to zero (based on substantive knowledge), specified as free to estimate (also based on substantive knowledge), or unspecified (learned through regularization). With different combinations of zero-fixed, specified, and unspecified loadings, different amounts of substantive input can be transformed into different structures of the loading matrix of  $\Lambda$ .

In EFA, both the number of factors and structure of  $\Lambda$  are unknown and of primary interest. This means all loadings are unspecified. Moreover,  $\Psi$  is modeled as a diagonal matrix, and the accuracy of loading and  $\Phi$  estimates is secondary. The model is usually underidentified and cannot be estimated with traditional algorithms such as the maximum likelihood estimation. Instead, purely data-driven techniques such as principal component analysis, subjective criteria such as eigenvalue, and rotation are adopted to obtain a small number of factors and a simple structure for the loading matrix. In CFA, the number of factors and the structure of  $\Lambda$  are given based on theory-driven hypothesis. This means that all loadings are either zero or specified. Free elements in  $\Lambda$  and  $\Phi$  are of primary interest.  $\Psi$  can be nondiagonal, with usually a few nonzero off-diagonal elements, which are nuisance parameters imposing challenges for model estimation and fitting. Under certain conditions, the model is overidentified and can be estimated with traditional methods (see, e.g., Kaplan, 2009).

The framework becomes essentially exploratory when all loadings are unspecified, and confirmatory when all loadings are either zero-fixed or specified. More generally, developers can designate some loadings as zero or specified while keeping others as unspecified in the framework, which means the loading matrix is partially specified. A typical PCFA model with two factors and 10 items can be found in Figure 1, where there is one specified and unspecified loading per item. Note that all residuals are allowed to correlate with each other in the figure. However, two thorny issues need to be addressed before the framework can be useful: the accommodation of both specified and unspecified loadings, and the identification of nonzero off-diagonal elements or residual correlations in  $\Psi$ . The Bayesian Lasso can address both issues.

#### The Bayesian Lasso Approach

#### **Bayesian Regression and Covariance Lasso**

The issues of both unspecified loadings and a residual covariance matrix can be addressed based on variable selection with the Bayesian Lasso. For a regression model with *p* predictors  $\boldsymbol{\beta} = (\beta_1, \dots, \beta_p)^T$  and responses

$$\mathbf{y} = \boldsymbol{\mu} \mathbf{1}_n + \mathbf{X} \boldsymbol{\beta} + \boldsymbol{\varepsilon}, \tag{2}$$

the Lasso estimates are least squares estimates with  $L_1$  norm penalty (Tibshirani, 1996):

$$\hat{\boldsymbol{\beta}} = \underset{\boldsymbol{\beta}}{\operatorname{argmin}} \left[ (\tilde{\mathbf{y}} - \mathbf{X}\boldsymbol{\beta})^T (\tilde{\mathbf{y}} - \mathbf{X}\boldsymbol{\beta}) + \delta \sum_{j=1}^p |\beta_j| \right]$$
(3)

where **X** is the standardized regressors,  $\tilde{\mathbf{y}} = \mathbf{y} - \bar{\mathbf{y}}\mathbf{1}_n$ ,  $\boldsymbol{\varepsilon}$  is the residual term, and  $\delta \ge 0$  is the tuning parameter that determines the amount of shrinkage. The Bayesian Lasso estimates accomplish the same goals as the frequentist counterpart by choosing appropriate forms of prior distributions that play the role of the penalty in Equation 3. In a fully Bayesian approach with a hier-



Figure 1. A C-step example in PCFA with two factors and 10 items.

archical formulation (Park & Casella, 2008), hierarchical priors can be applied and, as a result, the priors of the predictors follow the double exponential (i.e., Laplace) distributions with the shrinkage parameter. With hyperpriors and the MCMC algorithm, the shrinkage parameter can be estimated simultaneously with other parameters.

The Bayesian regression Lasso has been extended to address covariance structure with sparse off-diagonal terms bounded away from zero (Khondker, Zhu, Chu, Lin, & Ibrahim, 2013; Wang, 2012). Specifically, the  $L_1$  norm penalty is applied to the inverse of the covariance matrix with the graphical Lasso. Double exponential priors with similar shrinkage parameters are used for the off-diagonal elements while independent exponential priors can be used for the diagonal terms. With hyperparameters for the priors and an adapted MCMC algorithm, the Bayesian covariance Lasso is successfully implemented to address a sparse residual covariance structure under the CFA context (Pan et al., 2017).

#### The Confirmatory Steps

In PCFA, the loadings in one item can be zero, specified, or unspecified. For a specific item, the latent factors and loading vector in Equation 1 can be rearranged so that the zero or specified loadings and unspecified loadings can be partitioned. Denote the  $K \times 1$  loading vector for item j as  $\Lambda_j = \begin{pmatrix} X_j \\ \Lambda_j \end{pmatrix}$ , where  $\Lambda_j^T$  is the *j*th row of  $\Lambda$  and j = 1, 2, ..., J. The  $K' \times 1$  vector  $\Lambda'_j$  and  $K'' \times 1$ vector  $\Lambda''_j$  are the zero or specified and unspecified parts, respectively. All unspecified loadings are estimated with Bayesian Lasso and are called Lasso loadings in this research. Similar to Bayesian regression Lasso, different priors can be assigned to Lasso loadings in hierarchical form, as

$$\begin{aligned}
\Lambda_{j}^{"} &\sim N_{K''}(\mathbf{0}, \mathbf{D}_{\tau_{j}}), \\
\mathbf{D}_{\tau_{j}} &= diag(\tau_{j1}^{2}, \dots, \tau_{jK''}^{2}), \\
\tau_{jk}^{2} &\sim Gamma\left(1, \frac{\delta_{l}^{2}}{2}\right),
\end{aligned} \tag{4}$$

where k = 1, 2, ..., K'' and  $\delta_i$  is the shrinkage parameter. Note that an individual parameter can be set to shrink each item *j*, but the difference of regularization is trivial. Here one shrinkage parameter is used for simplicity of discussion. The hierarchical priors form a double exponential distribution with the  $L_1$  norm penalty, which can maintain conjugacy of the posterior distribution. The explicit form of the full conditional distribution can be obtained, as shown in Appendix A, which can be directly sampled from using the Gibbs sampler. Hyperpriors can be assigned to  $\delta_l$  as  $\delta_l^2 \sim Gamma(a_l, b_l)$  and the common choices for hyperparameters are 1 for  $a_l$  and a small  $b_l$ . For specified loadings, conjugate prior distribution  $\Lambda'_j \sim N_{K'}(\Lambda_{0j}, \mathbf{H}_{0j})$  can be assigned with  $\Lambda_{0j}$  and  $\mathbf{H}_{0j}$  as hyperparameters, which can also lead to an explicit full conditional distribution (Lee, 2007). If needed, a conjugate prior for the intercept vector can be assigned as  $\boldsymbol{\mu} \sim N_j(\boldsymbol{\mu}_0, \mathbf{H}_{\mu 0})$ , with  $\boldsymbol{\mu}_0$  and  $\mathbf{H}_{\mu 0}$  as hyperparameters and an explicit full conditional distribution for sampling.

The residual covariance matrix in a partially confirmatory model is not necessarily diagonal. The entire  $\Psi$  can be modeled as a sparse structure on off-diagonal elements with the Bayesian covariance Lasso. Denote  $\Sigma = \Psi^{-1} = (\sigma_{jj'})_{J \times J'}$ . Double exponential priors can be assigned to the off-diagonal elements, which has the form  $\frac{\delta_i}{2} \exp(-\delta_s | \sigma_{jj'} |)$  with j < j', whereas independent exponential priors can be assigned to the diagonal elements with the form  $\frac{\delta_i}{2} \exp(-\frac{\delta_i}{2} \sigma_{jj})$  (Wang, 2012). The full conditional distribution can be obtained and directly sampled from using a block Gibbs sampler. Hyperpriors can be assigned to the shrinkage parameter  $\delta_s$  as  $\delta_s \sim Gamma(a_s, b_s)$  and the common choices for hyperparameters are 1 for  $a_s$  and a small  $b_s$ . The related full conditional distributions are presented in Appendix C.

The factorial correlation matrix is also estimated by adapting the relevant MCMC algorithms (Liu, 2008; Liu & Daniels, 2006). However, an explicit form of the full conditional distribution for the factorial correlation matrix  $\mathbf{\Phi}$  cannot be easily obtained and directly sampled from. Instead, one can first sample from the original covariance matrix  $\mathbf{\Phi}^*$  with conjugate prior  $\mathbf{\Phi}^* \sim Inv-Wishart(S_0, \mathbf{v}_0)$  and then transform the sampled  $\mathbf{\Phi}^*$  into corresponding  $\mathbf{\Phi}$  with the reduction function  $\mathbf{\Phi} = \mathbf{D}^{-1}\mathbf{\Phi}^*\mathbf{D}^{-1}$ , where **D** is an expansion parameter matrix. The provisional  $\mathbf{\Phi}$  is accepted based on a Metropolis-Hastings acceptance probability, which is computed by comparing the determinant of the correlation matrix obtained from the previous draw and current draw. Liu (2008) and Liu and Daniels (2006) provide more details on how to sample the correlation matrix  $\mathbf{\Phi}$ .

It is challenging to estimate a partially specified loading matrix and a residual covariance matrix with off-diagonal elements simultaneously. The PCFA model can be especially problematic when there are too many unspecified or Lasso loadings. Specifically, we found that parameter estimates are unstable or difficult to converge if there is no specified loading for many items. This research imposed a constraint of having at least one specified loading per item, although it might be fine if there are a few items without any specified loading. The constraint is meaningful substantively since there is usually a target factor for every item during scale development. The specified loading can be the major loading under the BSEM context (B. O. Muthén & Asparouhov, 2012) or any other loading that the developer is confident to specify. With the constraint of one specified loading per item, the formulation tends to be confirmatory and is called the confirmatory step (C-step) in PCFA. Figure 1 gives an example of the C-step with two factors, 10 items, and one specified loading per item. Note that item residuals are allowed to completely correlate with each other under the Bayesian Lasso covariance estimation. When all loadings are specified, the C-step is essentially the BLCCFA (Pan et al., 2017), with a minor difference of fixing the factorial variance rather than having one item per factor for scale indeterminacy.

#### The Exploratory Step

When no loading can be specified for many items, the above constraint can be released by modeling  $\Psi$  as a diagonal matrix or ignoring any residual correlation. The loadings in  $\Lambda$  can be mostly unspecified given there is at least one specified loading per factor to identify the factor. For more stable performance, more specified loadings per factor are preferred. The hierarchical priors of the unspecified loadings conditional on the diagonal residual terms become:

$$\begin{split} \mathbf{\Lambda}_{jj}^{"} | \mathbf{\Psi}_{jj} &\sim N_{K''}(\mathbf{0}, \mathbf{\Psi}_{jj} \mathbf{D}_{\tau_j}), \\ \mathbf{\Psi}_{jj}^{-1} &\sim Gamma(a_{0j}, b_{0j}), \\ \mathbf{D}_{\tau_j} &= \operatorname{diag}(\tau_{j1}^2, \dots, \tau_{jK''}^2), \\ \tau_{jk}^2 &\sim Gamma\left(1, \frac{\delta_l^2}{2}\right), \end{split}$$
(5)

where  $\psi_{jj}$  is the *j*th diagonal term of  $\Psi$ , and the priors for the shrinkage parameters  $\delta_i$  are similar to the C-step. The explicit full conditional distribution can be obtained (see Appendix B) and directly sampled from using the Gibbs sampler. Other parameter estimates are the same as the C-step above, except that there is no off-diagonal element in the residual covariance matrix. This is

similar to the BSEM approach, but with the Lasso method rather than the ridge regression or spike-and-slab priors (Lu et al., 2016; B. O. Muthén & Asparouhov, 2012). Parameter estimates are deemed biased when the true residual structure is not diagonal, but it would suffice to identify at least one specified loading per item for the C-step. Because the formulation tends to be exploratory especially compared with the previous step, it is called the exploratory step (E-step) in PCFA. Figure 2 gives an example of the E-step with two factors, 10 items, and two specified loadings per factor. Note that there is no off-diagonal element in the residual structure.

#### Markov Chain Monte Carlo Estimation

Graphical representations of the hierarchical structure of the C-step and E-step are presented in Figures 3 and 4, respectively. The figures show how the item responses  $Y_{ii}$  can be modeled using latent factors, loadings, and residuals, which are characterized by their mean, covariance structure, and shrinkage parameters in a hierarchical form. The Bayesian hierarchical structure can be estimated using MCMC, which is a simulation-based algorithm that iteratively resamples from the probability distributions based on a stochastic process of Markov chains (Gill, 2002). Under mild regularity conditions, the Markov chains will converge to a stationary posterior distribution after a sufficiently large number of iterations called the burn-in period. Two types of MCMC are used: the Gibbs sampler (Casella & George, 1992), which is easier to implement but requires the explicit full conditional distributions, and the Metropolis-Hastings algorithm (Chib & Greenberg, 1995), which relies only on the joint posterior distribution. Except for the factorial correlation matrix  $\Phi$ , which is estimated using the Metropolis-Hastings algorithm as described above, all model parameters are estimated using the Gibbs sampler or block Gibbs sampler.

The procedure for sampling the parameters of interest from their full conditional distributions has six steps:

Step 1: Draw  $\Omega$  from  $p(\Omega | \mathbf{Y}, \boldsymbol{\mu}, \boldsymbol{\Lambda}, \boldsymbol{\Phi}, \boldsymbol{\Psi})$ ;

Step 2: Draw  $\mathbf{A}_{j}^{'}$  from  $p(\mathbf{A}_{j}^{'} | \mathbf{Y}, \mathbf{\Lambda}_{-j}, \mathbf{\Omega}, \boldsymbol{\mu}, \boldsymbol{\Phi}, \boldsymbol{\Psi}, \boldsymbol{\delta}_{lj})$  for j = 1 to J; Step 3: Draw  $\mathbf{A}_{j}^{'}$  from  $p(\mathbf{A}_{j}^{'} | \mathbf{Y}, \mathbf{\Lambda}_{-j}, \mathbf{\Omega}, \boldsymbol{\mu}, \boldsymbol{\Phi}, \boldsymbol{\Psi})$  for j = 1 to J; Step 4: Draw  $\boldsymbol{\mu}$  from  $p(\boldsymbol{\mu} | \mathbf{Y}, \mathbf{\Omega}, \mathbf{\Lambda}, \boldsymbol{\Phi}, \boldsymbol{\Psi})$ , if needed;

Step 5: Draw  $\Phi$  from  $p(\Phi | \mathbf{Y}, \Omega, \Lambda, \mu, \Psi)$ ;

Step 6: Draw  $\Sigma$  from  $p(\Sigma | \mathbf{Y}, \boldsymbol{\mu}, \boldsymbol{\Omega}, \boldsymbol{\Lambda}, \boldsymbol{\Phi}, \delta_s)$  for the C-step, or  $p(\Sigma | \mathbf{Y}, \boldsymbol{\mu}, \boldsymbol{\Omega}, \boldsymbol{\Lambda}, \boldsymbol{\Phi})$  for the E-step, and compute  $\Psi = \Sigma^{-1}$ .

Multiple chains with different initial values can be run to monitor the convergence of the algorithm. After the burn-in period, the



Figure 2. An E-step example in PCFA with two factors and 10 items.



Figure 3. A directed acyclic graph for the C-step.

convergence for the parameters of interest can be determined using the estimated potential scale reduction (EPSR) value (Gelman, 1996). The EPSR value compares the ratio of the weighted average of the within-chain variance and between-chain variance to the within-chain variance. The chains are said to converge to the stationary distribution if this ratio is less than 1.1 (Gelman, Carlin, Stern, & Rubin, 2004). In this research, we found that all chains had reached stationary status within 8,000 iterations and then set the burn-in as 10,000 iterations for all subsequent studies. After the burn-in phase, parameter estimates are inferenced based on 10,000 and 20,000 draws for the simulation and real-data analyses, respectively. The uncertainty of estimates is characterized with the concept of highest posterior density (HPD) intervals, or more specifically the  $100(1 - \alpha)\%$  HPD interval (Box & Tiao, 1973). Both the EPSR value and HPD interval can be obtained with the R package coda (Plummer, Best, Cowles, & Vines, 2006). All programming was conducted on the R platform (R Development Core Team, 2010). Finally,

two sets of prior values were evaluated for sensitivity analysis in a preliminary study, as shown in Table 1, and the difference turned out to be trivial. We adopted the first set of values for subsequent analyses, which was less informative.

#### **Empirical Studies**

#### Study 1: Performance in a Relatively Simple Case

Study 1 evaluated the performance of the E-step and C-step in a relatively simple case with two factors and 10 items, namely, J = 10 and K = 2, through a simulation study. The factorial correlation was  $\phi_{12} = \phi_{21} = .3$ ; the loading matrix was

$$\mathbf{\Lambda}^{T} = \begin{bmatrix} \lambda_{11} & \lambda_{21} & \lambda_{31} & \lambda_{41} & \lambda_{51} & \lambda_{61} & \lambda_{0} & \lambda_{0} & \lambda_{0} \\ \lambda_{0} & \lambda_{0} & \lambda_{0} & \lambda_{52} & \lambda_{62} & \lambda_{72} & \lambda_{82} & \lambda_{92} & \lambda_{10,2} \end{bmatrix},$$
  
with  $\lambda_{11} = \lambda_{21} = \lambda_{31} = \lambda_{41} = \lambda_{72} = \lambda_{82} = \lambda_{92} = \lambda_{10,2} = .7,$   
 $\lambda_{51} = \lambda_{61} = \lambda_{52} = \lambda_{62} = .5,$  and all other loadings  $\lambda_{0} = 0.$  For



Figure 4. A directed acyclic graph for the E-step.

Table 1		
Two Sets	of Prior	Values

Set	$a_{lj}$	b <sub>lj</sub>	$a_s$	b <sub>s</sub>	$a_0$	$b_0$	$\Lambda_{0j}$	$oldsymbol{H}_{0j}$	So	<b>v</b> <sub>0</sub>
1	1	.01	1	.01	1	.01	0	4 <b>I</b>	$I + .1_{od}$	K + 2
2	1	.1	1	.1	1	.1	0	Ι	$I + .5_{od}$	K + 8

*Note.*  $\mathbf{I}$  = identity matrix;  $\mathbf{I} + .1_{od}$  = diagonal elements as 1 and off-diagonal elements as .1;  $\mathbf{I} + .5_{od}$  = diagonal elements as 1 and off-diagonal elements as .5;  $\boldsymbol{\mu}_0$  and  $\mathbf{H}_{\mu 0}$  not set as the intercept is not of interest in this research.

the residual structure in  $\Psi$ , the diagonal elements were  $\psi_{jj} = .51$  for j = 1 to 4 and 7 to 10 and  $\psi_{jj} = .35$  for j = 5 and 6, and the nonzero off-diagonal elements were  $\psi_{12} = \psi_{21} = \psi_{74} = \psi_{47} = .3$  with all other off-diagonal elements  $\psi_0 = 0$ . Two levels of sample sizes were simulated, with N = 250 and 500, and the number of replications was 200 for each sample size.

For the E-step recovery, a PCFA model with two specified loadings per factor was fitted, that is,  $\lambda_{11}$  and  $\lambda_{21}$  for Factor F1 and  $\lambda_{92}$  and  $\lambda_{10,2}$  for Factor F2 (see Figure 2). To clarify, a structure matrix **R** with specified loading denoted by 1 and unspecified loading denoted by -1 can be defined corresponding to the factor loading matrix **A**, as follows:

which implies four specified loadings and 16 Lasso loadings (8 nonzero and 8 zero, i.e.,  $\lambda_0$ ). For the C-step recovery, the PCFA model in Figure 1 with one loading per item was fitted. The equivalent structure matrix is as follows:

which implies 10 specified loadings and 10 Lasso loadings (8 zero and 2 nonzero). Performance assessment involved computing the bias of the parameter estimates (BIAS), the mean of the standard

error estimates (*SE*), the root mean square error (RMSE) between the estimates and the true values, and the percentage of estimates significantly different from zero at  $\alpha = .05$  (SIG%) based on the HPD interval.

Simulation results for the E-step can be found in Table 2. The loading recovery in terms of BIAS and RMSE could be poor when the items were associated with off-diagonal residual elements, but was acceptable otherwise. Poor loading estimates could not be improved with a larger sample size, which suggests a systemic error due to ignoring the nonzero off-diagonal elements. Additional simulation studies revealed that the biases of estimates were also related to the level of residual correlations, the sparsity of the residual covariance matrix, and the magnitude of the crossloadings (see Appendix D for details). Recovery of factorial correlations was also poor with large SE and lower power (i.e., SIG%). However, the Type I error rates for misidentifying zero loadings (i.e.,  $\lambda_0$ ) as significant were almost zero, and the power for detecting significant loadings was close to one. This result suggests that all significant loading estimates will be correctly identified in the E-step and accordingly can be used as specified loadings to satisfy the sufficient constraint for the C-step.

Additional simulation studies were conducted with the same data sets by fitting the model in the E-step with one specified loading per factor (that is,  $\lambda_{11}$  for F1 and  $\lambda_{10,2}$  for F2) or no

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E-Step Results in Study 1

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			N = 1	250		N = 500				
Par	True	BIAS	RMSE	SE	SIG%	BIAS	RMSE	SE	SIG%	
N <sub>11</sub>	0.7	0.210	0.219	0.122	0.962	0.230	0.240	0.105	0.985	
N <sub>21</sub>	0.7	0.220	0.228	0.123	0.962	0.231	0.240	0.105	0.985	
N <sub>31</sub>	0.7	-0.111	0.125	0.096	0.971	-0.107	0.118	0.077	0.990	
N <sub>41</sub>	0.7	-0.130	0.142	0.098	0.976	-0.131	0.141	0.079	0.990	
N <sub>51</sub>	0.5	-0.054	0.079	0.133	0.986	-0.059	0.081	0.118	0.990	
N <sub>61</sub>	0.5	-0.063	0.085	0.133	0.971	-0.060	0.081	0.117	0.980	
N <sub>52</sub>	0.5	-0.002	0.058	0.133	0.981	0.002	0.058	0.118	0.965	
A <sub>62</sub>	0.5	0.002	0.059	0.133	0.976	-0.002	0.057	0.118	0.965	
N <sub>72</sub>	0.7	-0.030	0.068	0.124	0.929	-0.023	0.077	0.099	0.955	
N <sub>82</sub>	0.7	-0.030	0.071	0.122	0.929	-0.025	0.078	0.098	0.955	
۱ <sub>92</sub>	0.7	-0.030	0.075	0.123	0.929	-0.025	0.078	0.098	0.955	
A <sub>10.2</sub>	0.7	-0.035	0.072	0.122	0.929	-0.032	0.082	0.097	0.955	
N <sub>0</sub>	0	-0.006	0.102	0.168	0.002	-0.009	0.110	0.151	0.004	
¢ <sub>12</sub>	0.3	0.039	0.072	0.219	0.176	0.063	0.091	0.204	0.475	
155	0.35	0.051	0.068	0.044	1.000	0.037	0.050	0.030	1.000	
166	0.35	0.050	0.070	0.044	1.000	0.042	0.053	0.030	1.000	
$\psi_{ii}$	0.51	-0.019	0.129	0.055	1.000	-0.028	0.121	0.038	1.000	
$\hat{S}_l$	—	4.148	0.571	2.231	1.000	4.153	0.354	2.199	1.000	

....

Note.  $\psi_{ij}$  averaged across elements from j = 1 to 4 and 7 to 10;  $\lambda_0$  averaged across all zero loading estimates; for  $\delta_i$ , BIAS = mean; RMSE = SD.

specified loading (that is, all loadings were unspecified). The results presented in Table 3 indicate that the Type I error rates decreased and the power increased when the number of specified loadings per factor increased from zero to two. The case of one specified loading per factor was acceptable in general and can serve as the minimum constraint, while two or more specified loadings per factor is preferred for more stable performance in the E-step.

Table 4 summarizes the simulation results for the C-step. The loading estimates were acceptable in general, with those for the first factor relatively worse likely due to the within-factor residual covariance  $\psi_{12}$ . The recovery for nonzero Lasso loadings (i.e.,  $\lambda_{61}$ ,  $\lambda_{52}$ ) was also slightly worse. Recovery of factorial correlation was better than that in the E-step, but the power for the smaller sample size was still low. The Type I error rates for both zero Lasso loadings and zero residual terms were zero, suggesting the conservative nature of the Bayesian Lasso. The power for  $\psi_{12}$  was also close to zero, making it difficult to identify nonzero residual covariance within one factor. The reason was likely because the related  $\lambda_{11}$  and  $\lambda_{21}$  estimates were positively biased. Additional simulation studies revealed that power was also related to the position of the nonzero elements (that is, within-factor or between-factor) and the magnitude of residual correlation (see Appendix D for details). Finally, all estimates improved with a larger sample size.

The estimates of the shrinkage parameters (Tables 2 and 4) were about the same across different sample sizes, suggesting that the penalty was largely independent of the sample size. The increase of  $\delta_l$  from the E-step to the C-step might hint at interference between the  $L_1$  penalties on the loadings and off-diagonal residual elements. To better investigate the impact of the Lasso technique in parameter estimation, one simulated data set was randomly chosen, and the loadings  $\lambda_{10,1}$ ,  $\lambda_{12}$ ,  $\lambda_{61}$ , and  $\lambda_{52}$  were treated as specified and unspecified, respectively. The corresponding posterior distributions (see Figure 5) suggest that, compared with the specified loadings, the unspecified loadings always pushed toward zero with the Lasso prior, and cases with smaller values were more remarkable.

#### Study 2: Performance in a More Complex Case

Study 2 evaluated the performance of the E-step and C-step in a more complex case with three factors and 18 items, namely, J = 18 and K = 3, through a simulation study. The factorial correlations were set as  $\phi_{kk'} = .3$  for k and K' = 1 to 3 and  $k \neq k'$ . The loading matrix was

$$\mathbf{\Lambda}^{T} = \begin{bmatrix} \lambda_{11} \quad \lambda_{21} \quad \lambda_{31} \quad \lambda_{41} \quad \lambda_{51} \quad \lambda_{61} \quad \lambda_{71} \quad \lambda_{0} \quad \lambda_{18,1} \\ \lambda_{0} \quad \lambda_{0} \quad \lambda_{0} \quad \lambda_{0} \quad \lambda_{0} \quad \lambda_{0} \quad \lambda_{22} \quad \lambda_{12,2} \quad \lambda_{12,2} \quad \lambda_{13,2} \quad \lambda_{0} \quad \lambda_{0} \quad \lambda_{0} \quad \lambda_{0} \quad \lambda_{0} \quad \lambda_{0} \\ \lambda_{13} \quad \lambda_{0} \quad \lambda_{12,3} \quad \lambda_{13,3} \quad \lambda_{14,3} \quad \lambda_{15,3} \quad \lambda_{16,3} \quad \lambda_{17,3} \quad \lambda_{18,3} \end{bmatrix}$$

The following loadings were set as .7:  $\lambda_{j1}$  for j = 2 to 5,  $\lambda_{j2}$  for j = 8 to 11, and  $\lambda_{j3}$  for j = 14 to 17; the following loadings were set as .5:  $\lambda_{11}$ ,  $\lambda_{61}$ ,  $\lambda_{71}$ ,  $\lambda_{18,1}$ ,  $\lambda_{62}$ ,  $\lambda_{72}$ ,  $\lambda_{12,2}$ ,  $\lambda_{13}$ ,  $\lambda_{13}$ ,  $\lambda_{12,3}$ ,  $\lambda_{13,3}$ , and  $\lambda_{18,3}$ ; and all other loadings were set as  $\lambda_0 = 0$ . For the residual structure in  $\Psi$ , the diagonal elements were  $\psi_{jj} = .51$  for j = 2 to 5, 8 to 11, and 14 to 17 and  $\psi_{jj} = .35$  for j = 1, 6, 7, 12, 13, and 18. The nonzero off-diagonal elements for the lower triangle were  $\psi_{43} = \psi_{16,15} = \psi_{10,9} = \psi_{85} = \psi_{14,11} = \psi_{17,2} = .3$ 

with a symmetric upper triangle. The first three terms were withinfactor whereas the last three were between-factor. All other offdiagonal elements were set as  $\psi_0 = 0$ . This was a balanced design, as each factor was measured by four items with cross-loadings, two items with within-factor residual covariance, and two items with between-factor residual covariance.

The structure matrix  $\mathbf{R}^{\mathrm{T}}$  for the E-step and C-step were, respectively,

SIG%(0)

N = 500

SIG%(1)

SIG%(2)

Table 3Type I Error Rate and Power of Bayesian Lasso in the E-Step

SIG%(0)

True

Par

λ <sub>11</sub>	0.7	0.270	0.890	0.962	0.330	0.950	0.985
λ <sub>21</sub>	0.7	0.270	0.890	0.962	0.330	0.950	0.985
λ <sub>31</sub>	0.7	0.370	0.900	0.971	0.350	0.960	0.990
$\lambda_{41}$	0.7	0.450	0.910	0.976	0.460	0.970	0.990
$\lambda_{51}$	0.5	0.690	0.690	0.986	0.770	0.750	0.990
λ <sub>61</sub>	0.5	0.640	0.630	0.971	0.760	0.800	0.980
$\lambda_{52}$	0.5	0.770	1.000	0.981	0.750	0.990	0.965
λ <sub>62</sub>	0.5	0.750	0.990	0.976	0.750	0.990	0.965
λ <sub>72</sub>	0.7	0.290	0.890	0.929	0.380	0.960	0.955
λ <sub>82</sub>	0.7	0.290	0.890	0.929	0.370	0.960	0.955
λ <sub>92</sub>	0.7	0.290	0.890	0.929	0.370	0.960	0.955
λ <sub>10,2</sub>	0.7	0.300	0.890	0.929	0.370	0.960	0.955
λ <sub>0</sub>	0	0.270	0.014	0.002	0.401	0.017	0.004

SIG%(2)

N = 250

SIG%(1)

*Note.* SIG%(0), SIG%(1), and SIG%(2) represent the percentage of estimates significantly different from zero at  $\alpha = .05$  with no specified loading, one specified loading, and two specified loadings per factor, respectively.

Table 4			
C-step Results	in	Study	1

			N = 250				N = 500			
Par	True	BIAS	RMSE	SE	SIG%	BIAS	RMSE	SE	SIG%	
λ <sub>11</sub>	0.7	0.081	0.092	0.137	1.000	0.078	0.087	0.123	1.000	
$\lambda_{21}$	0.7	0.084	0.096	0.137	1.000	0.078	0.086	0.123	1.000	
λ <sub>31</sub>	0.7	-0.067	0.081	0.131	1.000	-0.058	0.068	0.115	1.000	
$\lambda_{41}$	0.7	-0.087	0.096	0.130	0.995	-0.082	0.089	0.116	1.000	
$\lambda_{51}$	0.5	-0.039	0.064	0.133	0.995	-0.043	0.059	0.116	0.995	
$\lambda_{61}$	0.5	-0.128	0.139	0.139	0.780	-0.105	0.114	0.125	0.875	
$\lambda_{52}$	0.5	-0.082	0.102	0.151	0.775	-0.061	0.075	0.131	0.910	
$\lambda_{62}^{52}$	0.5	0.013	0.050	0.134	0.985	0.001	0.036	0.117	0.995	
$\lambda_{72}$	0.7	-0.054	0.070	0.137	0.990	-0.046	0.056	0.117	1.000	
$\lambda_{82}$	0.7	-0.029	0.053	0.134	0.990	-0.015	0.035	0.112	1.000	
$\lambda_{02}$	0.7	-0.025	0.052	0.134	0.990	-0.016	0.036	0.112	1.000	
$\lambda_{10,2}$	0.7	-0.027	0.054	0.135	1.000	-0.017	0.036	0.112	1.000	
$\lambda_0$	0	-0.008	0.043	0.111	0.000	-0.008	0.037	0.104	0.000	
φ <sub>12</sub>	0.3	0.069	0.100	0.191	0.560	0.086	0.101	0.172	0.850	
$\psi_{21}$	0.3	-0.109	0.118	0.146	0.005	-0.102	0.110	0.138	0.040	
$\psi_{74}$	0.3	-0.077	0.086	0.098	0.755	-0.070	0.076	0.089	0.915	
$\psi_{55}$	0.35	0.106	0.116	0.114	1.000	0.083	0.091	0.095	1.000	
Ψ66	0.35	0.097	0.109	0.111	1.000	0.082	0.090	0.093	1.000	
$\Psi_{ii}$	0.51	0.025	0.091	0.133	1.000	0.014	0.075	0.118	1.000	
ΨΩ	0	0.023	0.043	0.082	0.000	0.019	0.036	0.071	0.000	
δ		7.584	0.488	2.653	1.000	7.543	0.355	2.528	1.000	
δ	_	2.188	0.125	0.354	1.000	2.276	0.100	0.360	1.000	

*Note.*  $\psi_{jj}$  averaged across elements from j = 1 to 4 and 7 to 10;  $\lambda_0$  averaged across all zero loading estimates; for  $\delta_l$  and  $\delta_s$ , BIAS = mean; RMSE = *SD*.

This implies that there were six specified loadings and 48 Lasso loadings (30 zero and 18 nonzero) for the E-step, and 18 specified loadings and 36 Lasso loadings (30 zero and 6 nonzero) for the C-step. The sample size was N = 250, and other conditions were similar to the above.

Table 5 summarizes the simulation results for the E-step and C-step. The performance of both steps in Study 2 was clearly better than for the simpler Study 1 case above. For the E-step, loading estimates involving residual covariance were worse, but all nonsignificant and significant loadings could be successfully identified. Not surprisingly, the C-step estimates were better than the E-step counterpart. In the C-step, the nonzero Lasso loadings (e.g.,  $\lambda_{71},\lambda_{18,1})$  performed similarly well as specified loadings. Loading estimates with within-factor residual covariance (e.g.,  $\lambda_{31}$ ,  $\lambda_{41}$ ) were still positively biased, but within a reasonable range. The extremely low Type I error rates for both zero Lasso loadings and zero residual terms were desirable. Recovery of factorial correlations and most residual terms was also satisfactory. Of a little concern was the power to identify within-factor residual covariance. In general, the results suggest that the PCFA approach has more satisfactory performance in this more complex case.

Additional simulation studies were conducted across different situations, including more heterogeneous factor loadings, different patterns of the loading and residual covariance matrix, and a different number of specified loadings. In brief, the results were generally similar to or consistent with the findings in these two simulation studies, and details can be found in Appendix D.

#### Study 3: Questionnaire of Mathematics Learning

During scale development, it is common to have some items with unclear factor loading, and this issue is well known within the scope of the PCFA approach. Table 6 lists potential items for a questionnaire addressing anxiety, confidence (i.e., self-efficacy), and utility (valuing) in mathematics learning. A similar background questionnaire can be found in large-scale assessments such as the Trends in International Mathematics and Science Study (TIMSS; International Association for the Evaluation of Educational Achievement, 2011). For content analysis, three mathematics educators were asked to rate if the items measured specific factors clearly, clearly not, or unclearly with 1, -1, or 0, respectively. The educators were encouraged to be more conservative, namely, to give a ranking of 1 or -1 only

and



Figure 5. Posterior distributions of specified and unspecified loadings. See the online article for the color version of this figure.

when they were very confident and to use 0 as the default value. The mean ratings in Table 6 reflect the degree of uncertainty for item-factor fitness from a substantive perspective, which cannot be easily accommodated under the traditional EFA or CFA approach. Under the PCFA approach, one can simply designate those elements with 1 as specified loadings and all others as Lasso loadings.

Responses from 218 Chinese fifth grade students were collected and analyzed using the E-step and C-step procedures in sequence, with estimation results given in Table 7. Significant Lasso loadings identified in the E-step were specified to satisfy the constraint necessary for the C-step. Note that for Item 14, two loadings were found significant with close values in the E-step and accordingly were both specified in the C-step. No significant residual covariance can be identified in the C-step, which can explain the similarity of estimation results between the two steps. One can also triangulate the results with the BLCCFA model by fixing all significant Lasso loadings in the C-step as specified and nonsignificant Lasso loadings as zero. Not surprisingly, the difference was trivial (results not shown to save space, but available upon request).

#### **Study 4: Humor Styles Questionnaire**

When a more confirmatory approach is adopted to develop a psychological scale, the C-step can be used directly. The Humor Styles Questionnaire (HSQ; Martin, Puhlik-Doris, Larsen, Gray, & Weir, 2003) with four factors and 32 items can be found in Appendix E. As each item supposedly loads on only one factor, the development process of the HSQ was oriented toward a confirmatory approach. Although the major loading of each item was confirmed, researchers were concerned about cross-loadings for some items, as the related behaviors tended to be multidimensional (Heintz, 2017).

Data for HSQ from 1,070 respondents were publicly available at https://openpsychometrics.org/\_rawdata/. Dissatisfactory fitting as shown in Table 8 was obtained using standard CFA, with no cross-loading or residual covariance in Mplus (L. K. Muthén & Muthén, 1998–2010) with the robust maximum likelihood estimation. Moreover, the PMM approach was challenging since many cross-loadings and residual covariance terms were suggested based on the modification index. In contrast, the C-step with major and Lasso loadings and residual covariance structure

Table 5				
Simulation	Results	for	Study	2

			E-step				C-step			
Par	True	BIAS	RMSE	SE	SIG%	BIAS	RMSE	SE	SIG%	
λ <sub>11</sub>	0.5	-0.014	0.064	0.093	1.000	-0.010	0.051	0.096	1.000	
$\lambda_{21}$	0.7	-0.056	0.083	0.083	1.000	-0.035	0.053	0.104	1.000	
λ <sub>31</sub>	0.7	0.162	0.169	0.085	1.000	0.082	0.093	0.106	1.000	
$\lambda_{41}$	0.7	0.163	0.168	0.085	1.000	0.084	0.093	0.106	1.000	
$\lambda_{51}$	0.7	-0.069	0.090	0.082	1.000	-0.034	0.054	0.104	1.000	
$\lambda_{61}$	0.5	-0.039	0.068	0.096	1.000	-0.007	0.049	0.096	1.000	
$\lambda_{71}$	0.5	-0.041	0.076	0.095	1.000	-0.056	0.081	0.099	1.000	
λ <sub>18.1</sub>	0.5	-0.018	0.063	0.093	1.000	-0.052	0.076	0.099	0.995	
λ <sub>62</sub>	0.5	-0.033	0.067	0.094	1.000	-0.061	0.084	0.100	1.000	
$\lambda_{72}$	0.5	-0.039	0.073	0.094	1.000	-0.017	0.055	0.097	1.000	
$\lambda_{82}$	0.7	-0.057	0.082	0.083	1.000	-0.042	0.062	0.107	1.000	
$\lambda_{92}$	0.7	0.179	0.184	0.087	1.000	0.080	0.090	0.108	1.000	
$\lambda_{10,2}$	0.7	0.174	0.180	0.086	1.000	0.087	0.096	0.108	1.000	
$\lambda_{11,2}$	0.7	-0.066	0.085	0.083	1.000	-0.038	0.055	0.105	1.000	
λ <sub>12.2</sub>	0.5	-0.039	0.067	0.094	1.000	-0.012	0.049	0.098	1.000	
λ <sub>13.2</sub>	0.5	-0.036	0.071	0.094	1.000	-0.057	0.080	0.101	0.990	
λ <sub>13</sub>	0.5	-0.048	0.081	0.096	1.000	-0.054	0.078	0.101	1.000	
$\lambda_{12,3}$	0.5	-0.046	0.076	0.093	1.000	-0.057	0.078	0.100	0.995	
λ <sub>13.3</sub>	0.5	-0.046	0.075	0.094	1.000	-0.012	0.054	0.097	1.000	
$\lambda_{14,3}$	0.7	-0.089	0.109	0.082	1.000	-0.042	0.061	0.106	1.000	
λ <sub>15.3</sub>	0.7	0.190	0.195	0.087	1.000	0.081	0.091	0.108	1.000	
$\lambda_{16,3}$	0.7	0.191	0.196	0.087	1.000	0.082	0.092	0.108	1.000	
λ <sub>17.3</sub>	0.7	-0.084	0.105	0.083	1.000	-0.041	0.060	0.105	1.000	
λ <sub>18.3</sub>	0.5	-0.053	0.079	0.096	1.000	-0.014	0.053	0.096	1.000	
λ	0	-0.013	0.070	0.107	0.002	-0.007	0.038	0.086	0.000	
$\phi_{kk'}$	0.3	0.027	0.068	0.156	0.612	0.031	0.074	0.135	0.805	
$\Psi_{ii}$	0.35	0.064	0.079	0.043	1.000	0.071	0.084	0.073	1.000	
$\psi_{ii}$	0.51	-0.065	0.149	0.052	1.000	-0.009	0.081	0.105	1.000	
$\psi_w$	0.3	_		_	_	-0.090	0.101	0.106	0.702	
$\psi_b$	0.3	—		—		-0.062	0.071	0.070	1.000	
ψο	0	—	_			0.005	0.028	0.053	0.000	
$\delta_l$		6.059	0.558	1.912	1.000	7.789	0.302	1.468	1.000	
δ	—					2.615	0.130	0.248	1.000	

*Note.* For  $\phi_{kk'}$ , k and K' = 1 to 3 and  $k \neq k'$ ; for  $\psi_{jj}$ , j = 2 to 4, 8 to 12, and 14 to 17; for  $\psi_{ii}$ , i = 1, 6, 7, 11, 12, and 18;  $\psi_w$  averaged across  $\psi_{43}$ ,  $\psi_{16,15}$ , and  $\psi_{10,9}$  (within factor);  $\psi_b$  averaged across  $\psi_{85}$ ,  $\psi_{14,11}$ , and  $\psi_{17,2}$  (between factors); for  $\delta_l$  and  $\delta_s$ , BIAS = mean; RMSE = *SD*.

can be directly applied. The estimation resulted in seven significant cross-loadings and 15 significant residual covariance terms, as shown in Table 9. It appeared that both cross-loadings and residual covariance contributed to model fitness. Note that most significant cross-loadings were around .1 and only six significant residual terms were above .1. This result suggests that many estimates were significant merely due to the sample size.

A BLCCFA model with all significant loadings in the C-step as specified was used for triangulation. It resulted in the same sig-

Table 6

Questionnaire of Mathematics Learning and the Mean Rating of Item-Factor Fitness

#	Content	F1	F2	F3
1	Mathematics makes me nervous	1.00	0.00	-1.00
2	Mathematics exams are easy for me	-0.33	1.00	-1.00
3	I worry about mathematics exams	1.00	0.00	-1.00
4	I am not bothered by mathematics courses	0.33	0.67	-0.67
5	I find mathematics confusing	0.67	0.33	-0.33
6	My brain is empty when solving mathematics problem	0.67	0.33	-0.33
7	Solving difficult mathematics problem is one of my strengths	-0.33	1.00	-1.00
8	I usually do well in mathematics	-0.67	1.00	-1.00
9	I learn things quickly in mathematics	-0.67	0.67	-1.00
10	Mathematics helps me in my daily life	-0.67	-1.00	1.00
11	I think mathematics will help me in my future life	-0.67	0.00	0.67
12	I need to do well in mathematics to get into the university of my choice	-1.00	-0.33	1.00
13	I need to do well in mathematics to get the job I want	-1.00	-0.33	1.00
14	I am willing to take more mathematics courses	-1.00	0.33	0.33

Note. All items employed a 5-point Likert scale, with 1 indicating agree a lot and 5 indicating disagree a lot; F1 = anxiety; F2 = confidence; F3 = utility.

 Table 7

 Parameter Estimates for the Questionnaire of Mathematics Learning

		E-ste	ep			C-step			
Item	F1	F2	F3	RV	F1	F2	F3	RV	
1	$0.708^{*}$	0.056	-0.044	0.522	0.660*	0.024	-0.041	0.587	
2	$-\overline{0.157}$	0.573*	-0.057	0.623	$-\overline{0.116}$	$0.560^{*}$	-0.019	0.657	
3	$0.653^{*}$	$-\overline{0.078}$	0.131	0.603	$0.584^{*}$	$-\overline{0.026}$	0.016	0.676	
4	-0.058	$0.269^{*}$	0.099	0.856	-0.027	0.366*	0.040	0.850	
5	$0.652^{*}$	0.005	-0.044	0.568	$0.627^{*}$	-0.016	-0.025	0.615	
6	$0.607^{*}$	-0.054	-0.064	0.562	0.637*	-0.057	-0.032	0.573	
7	0.007	$0.746^{*}$	-0.070	0.519	-0.008	$0.671^{*}$	-0.010	0.586	
8	0.028	$0.704^{*}$	0.034	0.501	0.013	$0.650^{*}$	0.063	0.565	
9	0.088	0.485*	0.149	0.681	0.051	$0.574^{*}$	0.051	0.681	
10	-0.113	0.064	$0.427^{*}$	0.710	-0.066	0.016	0.534*	0.692	
11	-0.047	0.085	$0.560^{*}$	0.584	-0.006	0.042	0.653*	0.569	
12	-0.066	-0.042	0.593*	0.649	-0.060	-0.008	0.551*	0.690	
13	0.076	-0.006	0.783*	0.440	0.052	0.042	$0.678^{*}$	0.564	
14	-0.042	0.319*	0.306*	0.650	-0.014	0.330*	0.339*	0.654	

*Note.* F1 = anxiety; F2 = confidence; F3 = utility; RV = residual variance. Underscored are specified loadings.

\* Significant loadings.

nificant loadings and 88 significant residual covariance terms, 19 of which were above .1 (see Table 9). The loading estimates were generally close, whereas there was some difference between the residual structures, suggesting the impact of the Lasso loadings.

Three standard CFA models were fitted with the same significant loadings identified in the C-step but with different residual structures: residual covariance above .1 in the C-step (M1), all significant residual covariance in the C-step (M2), and residual covariance above .1 in BLCCFA (M3). As shown in Table 8, all three models were acceptable and the difference was small. Moreover, although M2 was the best model, M1 was not a bad choice considering the small number of residual terms involved. The results suggest that the C-step can offer some flexibility in choosing the final model when the sample size is large.

#### Discussion

A different amount of substantive knowledge can be available during scale development, making the EFA and CFA two ends of a continuum. The partially confirmatory approach with Bayesian Lasso is flexible in addressing a wide range of the substantive continuum that would be challenging otherwise. Specifically, a C-step can be adopted when the constraint of one specified loading per item is satisfied, and an additional E-step is recommended before applying the C-step when only a few loadings per factor can be specified. Note that the C-step reduces to the Bayesian covariance Lasso CFA when there is no Lasso loading, while the E-step is similar to the BSEM without residual covariance. The PCFA approach with Bayesian Lasso can be derived using the Bayesian hierarchical formulation with different Lasso and regular priors by combining the Bayesian regression and covariance Lasso. Moreover, this approach is implemented using MCMC estimation with both a block Gibbs sampler and an adapted Metropolis-Hastings algorithm.

Simulation studies showed that both steps can perform as expected, and the performance was better for the complex case with three factors than for the simple case with two factors. This finding is consistent with the nature of the  $L_1$  norm penalty that can better select variables among many candidates and is meaningful for scale development with many factors, which is not uncommon in social and behavioral science. Study 3 suggested that a conservative approach can be adopted to specify the item-factor connections substantively with the help of both the E-step and C-step of the PCFA approach, allowing the data to give more information. This would be especially valuable when consensus or agreement among content experts is difficult to achieve. The two steps are also especially beneficial when the sample size is relatively small (e.g.,  $\sim$ 200). Practitioners can apply an intermediate stage after qualitative content analysis but before collecting data on a large scale.

Table 8CFA Model Fitness for the Humor Styles Questionnaire

	-							
Model	RMSEA	90%CI	CFI	TLI	SRMR	BIC	df	<sup>#</sup> RCT
M0	0.054	[0.052, 0.057]	0.844	0.831	0.065	96697	458	0
M1	0.040	[0.037, 0.043]	0.917	0.907	0.046	95981	445	6
M2	0.037	[0.035, 0.040]	0.929	0.920	0.044	95898	436	15
M3	0.038	[0.035, 0.041]	0.928	0.918	0.043	95923	432	19

*Note.* RMSEA = root mean square error; CFA = confirmatory factor analysis; M0 = no cross-loading or residual covariance; M1-M3 = all significant loadings identified in the C-step; M1 = residual covariance above .1 in the C-step; M2 = all significant residual covariance in the C-step; M3 = residual covariance above .1 in BLCCFA; #RCT = number of residual covariance term.

Table 9							
Significant	Loadings	and Residua	l Estimates for	r the	Humor	Styles	Questionnaire

		C-	step			BLC	CFA			RCV	
Item	F1	F2	F3	F4	F1	F2	F3	F4	Par	C-step	BLC
1	0.649				0.575				$\psi_{1.13}$		<u>0.116</u>
2		0.612				0.615			$\psi_{13.14}$		0.116
3			0.531				0.637		$\psi_{15,16}$	0.094	0.126
4				0.631				0.593	$\psi_{1,17}$		0.122
5	0.615				0.616				$\psi_{10,18}$	<u>0.146</u>	0.114
6	<u>0.203</u>	0.378			0.242	0.327			$\psi_{11,19}$	0.112	0.109
7	-0.142		0.580		-0.110		0.539		$\psi_{8,20}$	<u>0.153</u>	0.165
8				0.714				0.698	$\psi_{11,21}$		-0.130
9	0.452				0.461				$\psi_{17,21}$		0.142
10		0.726				0.732			$\psi_{11,23}$	-0.079	-0.073
11			0.540				0.483		$\psi_{4,24}$	0.096	0.123
12				0.635				0.610	$\psi_{15,24}$	-0.072	-0.084
13	0.614				0.530				$\psi_{1,25}$		0.124
14		0.674				0.669			$\psi_{5,25}$	-0.085	
15			0.647				0.579		$\psi_{13,25}$	0.147	0.263
16				0.567				0.592	$\psi_{17,25}$		0.130
17	0.697				0.624				$\psi_{21,25}$		0.142
18		0.723				0.690			$\psi_{23,25}$	0.086	0.116
19	0.178		0.452		0.344		0.435		$\psi_{27,28}$	0.092	
20				0.710				0.679	$\psi_{5,29}$	0.110	0.158
21	0.651				0.583				$\psi_{22,29}$	0.069	0.093
22		0.422				0.449			$\psi_{24,29}$		-0.113
23			0.489				0.496		$\psi_{25,29}$		0.110
24				0.483				0.423	$\psi_{6,30}$	0.218	0.235
25	0.634				0.484				$\psi_{22,31}$	0.071	0.097
26		0.666				0.686					
27			0.551				0.549				
28		0.140	<u>0.160</u>	0.249		<u>0.190</u>	0.271	0.221			
29	0.561		0.125	-0.163	0.472		0.128	-0.187			
30		0.447				0.397					
31			0.637				0.574				
32				0.659				0.688			

*Note.* BLCCFA = Bayesian covariance Lasso for CFA; F1 = affiliative humor; F2 = self-enhancing humor; F3 = aggressive humor; F3 = self-defeating humor; RCV = residual covariance; BLC = BLCCFA; Underscored are cross-loadings or residual covariance above .1; only residual covariance significant at the C-step or significant and above .1 in BLCCFA is presented.

Study 4 demonstrated the advantage of the C-step to identify significant cross-loadings and residual covariance when encountering many factors with a large number of items, a task that would be challenging with the traditional PMM approach. Although unnecessary because the constraint of one specified loading per item was well satisfied with the available literature, we tried the E-step with only two major loadings per factor and found that all other major loadings were successfully recovered. Accordingly, a combination of the E-step and C-step can offer more freedom during scale development. The C-step can also offer some flexibility in choosing a succinct model when the sample size is large. Finally, the Bayesian covariance Lasso CFA without any Lasso loading can be used for triangulation, as shown in both studies.

Although the PCFA with Bayesian Lasso approach appears to be promising, some issues can be addressed in future research to make the framework more comprehensive and versatile.

First, the PCFA approach assumes that substantive knowledge about the loading structure is partially available but knowledge about the factors (i.e., number of factors and their operational definitions) is largely available. During scale development, however, it is likely that there is a different amount of knowledge about the factors (e.g., a range of factor numbers or blurred definitions). It would be valuable but also challenging to extend the PCFA to accommodate these scenarios.

Second, it would be useful to equip the proposed framework with powerful means of model comparison to allow comparison of different substantive knowledge. One possible but also challenging means is to rely on Bayes factor for comparison.

Third, categorical data are widely available in social and behavioral research, as shown in Studies 3 and 4. These data are treated as continuous in the existing approach, which is suboptimal and can result in substantial bias without favorite conditions (e.g., normality). Accordingly, it would be valuable to extend the approach to address dichotomous and polytomous responses, possibly under the umbrella of item response theory.

Finally, Lu, Chow, and Loken (2016) proposed a Bayesian CFA with the spike-and-slab prior (SSP) and found that their approach outperformed the ridge regression prior and yielded more parsimonious loading structures. However, there is no research to compare the performance of Bayesian Lasso in estimating the loadings with that of the SSP under the CFA context. Future studies can further investigate if the SSP can be incorporated into the PCFA framework, and if it is possible, explain how that differs from the existing approach.

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(Appendices follow)

14

#### Appendix A

### Full Conditional Distribution for $\Lambda_i^{''}$ in the C-Step

Equation 1 can be rewritten as  $y_{ij} = \boldsymbol{\mu}_j + \boldsymbol{\Lambda}_j^T \boldsymbol{\omega}_i + \boldsymbol{\varepsilon}_{ij}$ , where  $\boldsymbol{\Lambda}_j^T = (\lambda_{j1}, \lambda_{j2}, \dots, \lambda_{jK})$  is the *j*th row of  $\boldsymbol{\Lambda}$ . After rearrange the latent factors and loading vector for item *j*, the factor vector is  $\boldsymbol{\omega}_i = \begin{pmatrix} \omega_i \\ \omega_i \end{pmatrix}$  where the  $K' \times 1$  vector  $\boldsymbol{\omega}_i'$  and  $K'' \times 1$  vector  $\boldsymbol{\omega}_i''$  are associated with the zero or specified loadings and unspecified loadings, respectively. Similarly, denote the rearranged factor matrix as  $\boldsymbol{\Omega} = \begin{pmatrix} \Omega' \\ \Omega'' \end{pmatrix} = \begin{pmatrix} \omega_{1}' \dots \omega_{N} \\ \omega_{1}' \dots \omega_{N}'' \end{pmatrix}$ . The conditional distribution for  $\boldsymbol{\Lambda}_i''$  can be expressed as:

$$p(\mathbf{\Lambda}_{j}^{"}|\mathbf{Y}, \mathbf{\Lambda}_{-j}, \mathbf{\Omega}, \boldsymbol{\mu}, \boldsymbol{\Phi}, \boldsymbol{\Psi}, \boldsymbol{\delta}_{lj}) \propto p(\mathbf{\Lambda}_{j}^{"}|\mathbf{y}_{j}, \mathbf{y}_{-j}, \mathbf{\Lambda}_{-j}, \mathbf{\Omega}, \boldsymbol{\Psi}, \boldsymbol{\delta}_{l})$$

$$\propto p(\mathbf{y}_{j}|\mathbf{y}_{-j}, \mathbf{\Lambda}_{j}^{"}, \mathbf{\Lambda}_{-j}, \mathbf{\Omega}, \boldsymbol{\Psi})p(\mathbf{\Lambda}_{j}^{"})$$

$$= \prod_{i=1}^{N} p(y_{ij}|y_{i(-j)}, \mathbf{\Lambda}_{j}^{"}, \mathbf{\Lambda}_{-j}, \boldsymbol{\omega}_{i}, \boldsymbol{\Psi})p(\mathbf{\Lambda}_{j}^{"})$$

$$\propto \prod_{i=1}^{N} \exp\left\{\frac{1}{2}\left[y_{ij} - \boldsymbol{\mu}_{j} - (\mathbf{\Lambda}_{j}^{"})^{T}\boldsymbol{\omega}_{i}^{"} - \boldsymbol{\psi}_{j}^{T}\boldsymbol{\Psi}_{-jj}^{-1}(\mathbf{y}_{i(-j)} - \boldsymbol{\mu}_{-j} - \mathbf{\Lambda}_{-j}^{T}\boldsymbol{\omega}_{i})\right]^{T} \times (\boldsymbol{\psi}_{jj} - \boldsymbol{\psi}_{j}^{T}\boldsymbol{\Psi}_{-jj}^{-1}(\mathbf{y}_{ij} - \boldsymbol{\mu}_{j} - (\mathbf{\Lambda}_{j}^{"})^{T}\boldsymbol{\omega}_{i}^{"} - \boldsymbol{\psi}_{j}^{T}\boldsymbol{\Psi}_{-jj}^{-1}(\mathbf{y}_{i(-j)} - \boldsymbol{\mu}_{-j} - \mathbf{\Lambda}_{-j}^{T}\boldsymbol{\omega}_{i})\right]\right\} \times p(\mathbf{\Lambda}_{j}^{"}).$$

where  $\psi_{jj}$  is the *j*th diagonal element of  $\Psi$ ,  $\psi_j = (\psi_{j1}, \ldots, \psi_{j,j-1}, \psi_{j,j+1}, \ldots, \psi_{j,j})^T$  is the vector of all off-diagonal elements of the *j*th column, and  $\Psi_{-jj}$  is the  $(J - 1) \times (J - 1)$  matrix resulting from deleting the *j*th row and *j*th column from  $\Psi$ .

Let  $\Psi_j^* = \psi_{jj} - \psi_j^T \Psi_{-jj}^{-1} \psi_j$ ,  $\mathbf{y}_{ij}^* = \mathbf{y}_{ij} - \mu_j - \psi_j^T \Psi_{-jj}^{-1} (\mathbf{y}_{i(-j)} - \mu_{-j} - \Lambda_{-j}^T \omega_i)$ , and  $\mathbf{Y}_j^* = (\mathbf{y}_{1j}^*, \mathbf{y}_{2j}^*, \dots, \mathbf{y}_{Nj}^*)^T$ . With Equation 5 as the prior for  $\Lambda_j^{"}$ , the right hand side becomes:

$$\prod_{i=1}^{n} \exp\left\{-\frac{1}{2}(\mathbf{y}_{ij}^{*}-(\boldsymbol{\Lambda}_{j}^{*})^{T}\boldsymbol{\omega}_{i}^{*})^{T}\boldsymbol{\Psi}_{j}^{*-1}(\mathbf{y}_{ij}^{*}-(\boldsymbol{\Lambda}_{j}^{*})^{T}\boldsymbol{\omega}_{i}^{*})\right\} \times p(\boldsymbol{\Lambda}_{j}^{*})$$

$$= \exp\left\{-\frac{1}{2}\boldsymbol{\Psi}_{j}^{*-1}(\mathbf{Y}_{j}^{*T}\mathbf{Y}_{j}^{*}-2(\boldsymbol{\Lambda}_{j}^{*})^{T}\boldsymbol{\Omega}^{*}\mathbf{Y}_{j}^{*}+(\boldsymbol{\Lambda}_{j}^{*})^{T}(\boldsymbol{\Omega}^{*}(\boldsymbol{\Omega}^{*})^{T})\boldsymbol{\Lambda}_{j}^{*})\right\} \times p(\boldsymbol{\Lambda}_{j}^{*}).$$

$$\Rightarrow p(\boldsymbol{\Lambda}_{j}^{*} \mid \mathbf{Y}, \boldsymbol{\Lambda}_{-j}, \boldsymbol{\Omega}, \boldsymbol{\mu}, \boldsymbol{\Phi}, \boldsymbol{\Psi}) \propto N(m, var), \text{ where }$$

$$m = (\boldsymbol{\Psi}_{j}^{*-1}(\boldsymbol{\Omega}^{*}(\boldsymbol{\Omega}^{*})^{T}) + \mathbf{D}_{\tau_{j}}^{-1})^{-1}(\boldsymbol{\Psi}_{j}^{*-1}\boldsymbol{\Omega}^{*}\mathbf{Y}_{j}^{*}) \text{ and } var =$$

$$(\boldsymbol{\Psi}_{j}^{*-1}(\boldsymbol{\Omega}^{*}(\boldsymbol{\Omega}^{*})^{T}) + \mathbf{D}_{\tau_{j}}^{-1})^{-1}.$$

The conditional distribution for  $\tau_{jk}$  can be expressed as:

$$p(\tau_{jk}^{2} | \mathbf{Y}, \mathbf{\Lambda}, \mathbf{\Omega}, \boldsymbol{\mu}, \boldsymbol{\Phi}, \boldsymbol{\Psi}) \propto p(\lambda_{jk} | \tau_{jk}^{2}) p(\tau_{jk}^{2})$$
$$\propto (\tau_{jk}^{2})^{-\frac{1}{2}} \exp\left\{-\frac{1}{2\tau_{jk}^{2}}\lambda_{jk}^{2}\right\} \exp\left\{-\frac{\delta_{l}^{2}}{2}\tau_{jk}^{2}\right\}.$$
$$\Rightarrow p\left(\frac{1}{\tau_{jk}^{2}} | \mathbf{Y}, \mathbf{\Lambda}, \mathbf{\Omega}, \boldsymbol{\mu}, \boldsymbol{\Phi}, \boldsymbol{\Psi}\right) \propto Inv-Gaussian\left(\sqrt{\frac{\delta_{l}^{2}}{\lambda_{jk}^{2}}}, \delta_{l}^{2}\right).$$

The conditional distribution for  $\delta_i$  can be expressed as:

$$p(\delta_l^2 | \mathbf{Y}, \mathbf{\Lambda}, \mathbf{\Omega}, \mathbf{\mu}, \mathbf{\Phi}, \mathbf{\Psi}) \propto p(\tau_{jk}^2 | \delta_l^2) p(\delta_l^2)$$
  
$$\propto \frac{\delta_l^2}{2} \exp\left\{-\frac{\delta_l^2}{2} \tau_{jk}^2\right\} \times (\delta_l^2)^{a_l - 1} \exp\left\{-b_l(\delta_l^2)\right\}$$
  
$$\propto (\delta_l^2)^{(a_l + 1) - 1} \exp\left\{-\left(\frac{\tau_{jk}^2}{2} + b_l\right)(\delta_l^2)\right\}$$
  
$$\propto Gamma\left(a_l + 1, \frac{\tau_{jk}^2}{2} + b_l\right).$$

(Appendices continue)

#### Appendix B

#### Full Conditional Distribution for $\Lambda_i^{''}$ in the E-Step

Different from the C-step,  $\Psi = diag(\psi_{jj})$  is modeled as a diagonal matrix. Let  $\mathbf{Y}_{j}^{*} = (\mathbf{y}_{1j}, \mathbf{y}_{2j}, \dots, \mathbf{y}_{Nj})^{T}$ . With notations similar above and Equation 6 as the prior, the conditional distribution for  $\Lambda_i^{"}$  can be expressed as:

$$p(\boldsymbol{\Lambda}_{j}^{"}|\mathbf{Y}, \boldsymbol{\Omega}, \boldsymbol{\mu}, \boldsymbol{\Phi}, \boldsymbol{\Psi}, \boldsymbol{\delta}_{lj}) \propto p(\mathbf{y}_{j} | \boldsymbol{\Lambda}_{j}^{"}, \boldsymbol{\Omega}, \boldsymbol{\Psi}) p(\boldsymbol{\Lambda}_{j}^{"} | \boldsymbol{\psi}_{jj})$$

$$\propto \prod_{i=1}^{N} \exp\left\{-\frac{1}{2}(y_{ij} - (\boldsymbol{\Lambda}_{j}^{"})^{T} \boldsymbol{\omega}_{i}^{"})^{T} \boldsymbol{\psi}_{jj}^{-1}(y_{ij} - (\boldsymbol{\Lambda}_{j}^{"T} \boldsymbol{\omega}_{i}^{"}) \times p(\boldsymbol{\Lambda}_{j}^{"} | \boldsymbol{\psi}_{jj})\right\}$$

$$\Rightarrow p(\boldsymbol{\Lambda}_{j}^{"} | \mathbf{Y}, \boldsymbol{\Omega}, \boldsymbol{\mu}, \boldsymbol{\Phi}, \boldsymbol{\Psi}, \boldsymbol{\delta}_{j}) \propto N(m, var), \text{ where }$$

$$m = (\boldsymbol{\Omega}^{"}(\boldsymbol{\Omega}^{"})^{T} + \mathbf{D}_{\tau_{j}}^{-1})^{-1}(\boldsymbol{\Omega}^{"}\mathbf{Y}_{j}^{*}) \text{ and } var = \psi_{jj}(\boldsymbol{\Omega}^{"}(\boldsymbol{\Omega}^{"})^{T} + \mathbf{D}_{\tau_{j}}^{-1})^{-1}.$$

The conditional distribution for  $\psi_i$  can be expressed as:

$$p(\boldsymbol{\psi}_{jj}^{-1} | \mathbf{Y}, \boldsymbol{\Lambda}_{j}^{''}, \boldsymbol{\Omega}, \boldsymbol{\mu}, \boldsymbol{\Phi}) \propto p(\mathbf{y}_{j} | \mathbf{y}_{-j}, \boldsymbol{\Lambda}_{j}^{''}, \boldsymbol{\Omega}, \boldsymbol{\Psi}) p(\boldsymbol{\Lambda}_{j}^{''} | \boldsymbol{\psi}_{jj}) p(\boldsymbol{\psi}_{jj}^{-1}).$$

which can be simplified as

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$$\propto Gamma \left( a_{0j} + \frac{N}{2} + \frac{K}{2} - 1, b_{0j} + \frac{1}{2} \left[ \sum_{i=1}^{N} (y_{ij} - (\Lambda_j^{"})^T \boldsymbol{\omega}_i^{"})^2 + (\Lambda_j^{"})^T \boldsymbol{D}_{\tau_j}^{-1} \Lambda_j^{"} \right] \right).$$

Similar to the C-step, the conditional distribution for  $\tau_{ik}$  and  $\delta_i$ can be expressed, respectively, as:

$$p(\frac{1}{\tau_{jk}^2}|\mathbf{Y}, \Lambda, \Omega, \mu, \Phi, \Psi) \propto Inv-Gaussian\left(\sqrt{\frac{\delta_l^2}{\lambda_{jk}^2}\psi_{jj}}, \delta_l^2\right),$$

and

$$p(\delta_l^2 | \mathbf{Y}, \boldsymbol{\Lambda}, \boldsymbol{\Omega}, \boldsymbol{\mu}, \boldsymbol{\Phi}, \boldsymbol{\Psi}) \propto Gamma(a_l + 1, \frac{\tau_{jk}^2}{2} + b_l)$$

#### Appendix C

#### Estimation of the Residual Covariance Matrix in the C-Step

The conditional distribution  $p(\Sigma | \mathbf{Y}, \boldsymbol{\mu}, \boldsymbol{\Omega}, \boldsymbol{\Lambda}, \boldsymbol{\Phi}, \boldsymbol{\tau}, \boldsymbol{\delta}_s)$  can be decomposed as follows:

$$p(\mathbf{\Sigma} | \mathbf{Y}, \mathbf{\mu}, \mathbf{\Omega}, \mathbf{\Lambda}, \mathbf{\Phi}, \mathbf{\tau}, \mathbf{\delta}_{s}) \propto p(\mathbf{Y} | \mathbf{\Sigma}, \mathbf{\mu}, \mathbf{\Omega}, \mathbf{\Lambda}, \mathbf{\Phi}) p(\mathbf{\tau}, \mathbf{\delta}_{s})$$
$$\propto |\mathbf{\Sigma}|^{N2} \exp\left[-\left(-\frac{1}{2}\mathbf{S}\mathbf{\Sigma}\right)\right] \prod_{i < j} \exp\left(-\frac{\sigma_{ij}^{2}}{2\tau_{ij}}\right) \times \prod_{j=1}^{J} \exp\left(-\frac{\delta_{s}\sigma_{jj}}{2}\right) \mathbf{I}(\mathbf{\Sigma})$$
$$> 0),$$

where  $\mathbf{\tau} = (\tau_{ij})_{i < j}$  is the vector of the latent scale parameters, and  $\mathbf{S} = \sum_{i=1}^{N} (\mathbf{y}_i - \boldsymbol{\mu} - \boldsymbol{\Lambda} \boldsymbol{\omega}_i) (\mathbf{y}_i - \boldsymbol{\mu} - \boldsymbol{\Lambda} \boldsymbol{\omega}_i).$ 

For k = 1, 2, ..., J, and without loss of generality, partition and rearrange the columns of  $\Sigma$  and S as follows:

$$\Sigma = \begin{pmatrix} \Sigma_{-kk} & \sigma_k \\ \sigma_k^T & \sigma_{kk} \end{pmatrix}, S = \begin{pmatrix} S_{-kk} & s_k \\ s_k^T & s_{kk} \end{pmatrix}.$$
  
Where  $\sigma_{kk}$  is the k-th diagonal element of  $\Sigma$ ,  $\sigma_k = (\sigma_{k1}, \ldots, \sigma_{kk})$ 

 $\sigma_{k,k-1}, \sigma_{k,k+1}, \ldots, \sigma_{k,j}$  is the vector of all off-diagonal elements of the k-th column, and  $\Sigma_{-kk}$  is the  $(p-1) \times (p-1)$  matrix resulting from deleting the k-th row and k-th column from  $\Sigma$ . Similar,  $s_{kk}$  is the k-th diagonal element of S,  $s_k$  is the vector of all-diagonal elements of the k-th column of S, and  $S_{-kk}$  is the matrix with the k-th row and k-th column of S deleted. Therefore we have:

$$p(\boldsymbol{\sigma}_{k},\boldsymbol{\sigma}_{kk}|\boldsymbol{\Sigma}_{-kk}, \mathbf{Y}, \boldsymbol{\Omega}, \boldsymbol{\mu}, \boldsymbol{\Lambda}, \boldsymbol{\Phi}, \boldsymbol{\tau}, \boldsymbol{\delta}_{s}) \propto (\boldsymbol{\sigma}_{kk} - \boldsymbol{\sigma}_{k}^{T}\boldsymbol{\Sigma}_{-kk}^{-1}\boldsymbol{\sigma}_{k})^{\frac{N}{2}}$$
$$\times \exp\left\{-\frac{1}{2}[\boldsymbol{\sigma}_{k}^{T}\boldsymbol{M}_{\tau}^{T}\boldsymbol{\sigma}_{k} + 2\boldsymbol{s}_{k}^{T}\boldsymbol{\sigma}_{k} + (\boldsymbol{s}_{kk} + \boldsymbol{\delta}_{s})\boldsymbol{\sigma}_{kk}]\right\}$$

where  $\mathbf{M}_{\tau}$  is the diagonal matrix with diagonal elements  $\tau_{k1}, \ldots,$  $\tau_{k,k-1}, \tau_{k,k+1}, \ldots, \tau_{kJ}$ 

Let  $\boldsymbol{\beta} = \boldsymbol{\sigma}_k$  and  $\boldsymbol{\gamma} = \boldsymbol{\sigma}_{kk} - \boldsymbol{\sigma}_k^T \boldsymbol{\Sigma}_{-kk}^{-1} \boldsymbol{\sigma}_k$ . It can be shown that:

$$p(\boldsymbol{\beta}|\boldsymbol{\Sigma}_{-kk}, \mathbf{Y}, \boldsymbol{\Omega}, \boldsymbol{\mu}, \boldsymbol{\Lambda}, \boldsymbol{\Phi}, \boldsymbol{\tau}, \boldsymbol{\delta}_s)$$

 $\propto N(-[(s_{kk} + \delta_s)\boldsymbol{\Sigma}_{-kk}^{-1} + \mathbf{M}_{\tau}^{-1}]^{-1}\boldsymbol{s}_k, [(s_{kk} + \delta_s)\boldsymbol{\Sigma}_{-kk}^{-1} + \mathbf{M}_{\tau}^{-1}]^{-1}),$  $p(\mathbf{\gamma} | \mathbf{\Sigma}_{-kk}, \mathbf{Y}, \mathbf{\Omega}, \mathbf{\mu}, \mathbf{\Lambda}, \mathbf{\Phi}, \mathbf{\tau}, \delta_s) \propto Gamma\left(\frac{N}{2} + 1, \frac{s_{kk} + \delta_s}{2}\right).$ After simulating observations from the above conditional distribution

tributions, we can obtain  $\sigma_k = \beta$ ,  $\sigma_k^T = \beta^T$  and  $\sigma_{kk} = \gamma$  +  $\sigma_k^T \Sigma_{-kk}^{-1} \sigma_k$ , then the k-th column and row of  $\Sigma$  are updated at a time. At the end,  $\boldsymbol{\psi} = \boldsymbol{\Sigma}^{-1}$  is computed.

The conditional distribution for  $\mathbf{\tau} = (\tau_{ij})_{i < j}$  can be expressed as:

$$\approx \prod_{i < j} \tau_{ij}^{-\frac{1}{2}} \exp\left\{-\frac{\sigma_{ij}^2 + \tau_{ij}^2 \delta_s^2}{2\tau_{ij}}\right\}.$$

It can be shown that for i < j,

$$p\left(\frac{1}{\tau_{ij}}|\mathbf{Y}, \mathbf{\Lambda}, \mathbf{\Omega}, \boldsymbol{\mu}, \boldsymbol{\Phi}, \boldsymbol{\Sigma}, \boldsymbol{\delta}_{s}\right) \propto Inv - Gaussian\left(\sqrt{\frac{\delta_{s}^{2}}{\sigma_{ij}^{2}}}, \boldsymbol{\delta}_{s}^{2}\right)$$

The conditional distribution for  $\delta_s$  can be expressed as:

$$p(\delta_{s} | \mathbf{Y}, \mathbf{\Lambda}, \mathbf{\Omega}, \mathbf{\mu}, \mathbf{\Phi}, \mathbf{\Sigma}, \mathbf{\tau}) \propto p(\mathbf{\tau} | \delta_{s}) p(\delta_{s})$$
$$\propto Gamma\left(a_{s} + \frac{J(J+1)}{2}, b_{s} + \frac{1}{2} \sum_{i=1}^{J} \sum_{j=1}^{J} |\sigma_{ij}|\right)$$

(Appendices continue)

#### Appendix D

#### **Additional Simulation Studies**

Additional simulation studies were conducted under more heterogeneous factor loadings, as shown in Table D1. The model settings were the same as in Study 2, except that the data were generated by a different factor loading matrix with more heterogenous loading values. In general, the results indicated that the performance under the condition with more heterogeneous loading was similar to that presented in Study 2.

We found that if at least one specified loading per item was imposed, parameter recovery was satisfactory in the C-step. To better understand the consequences of imposing this identifiability constraint, additional simulation studies were conducted. The settings were the same as those in Study 1 except that the data were generated with diagonal  $\Psi$  (i.e., the residuals in the true model were assumed to be independent). The C-step was fitted with three specified loadings per factor (that is,  $\lambda_{11}$ ,  $\lambda_{21}$  and  $\lambda_{31}$  are specified for F1 and  $\lambda_{82}$ ,  $\lambda_{92}$ ,  $\lambda_{10,2}$  are specified for F2). As shown in Table D2, the recoveries of most parameters were unacceptable. Accordingly, it is difficult to relax the constraint. Moreover, the constraint is easy to understand and to follow since there is usually a target factor for each item in scale development.

The consequence of correlated residuals is the existence of local dependence, which needs to be accounted for during modeling. Treating the correlated residuals as independent can distort the loading estimates, especially those associated with the correlated residuals, as shown in the E-step of Study 1 and 2. Additional simulation studies were conducted to further understand the consequence of correlated residuals. The settings were the same as those of Study 1 and 2 except that the data were generated with diagonal  $\Psi$ . The results of the E-step (Table D3 and D4) indicated that the estimates of all loading parameters were satisfactory. Accordingly, the biases of loading estimates in the E-step of Study 1 and 2 did come from the correlated residuals.

To investigate whether the model performs depending on the number of unspecified Lasso loadings or on the structure of the

 Table D1

 Simulation Studies With More Heterogeneous Loading Values

			E-st	ep			C-st	ep	
Par	True	BIAS	RMSE	SE	SIG%	BIAS	RMSE	SE	SIG%
λ <sub>11</sub>	0.8	0.034	0.062	0.096	1.000	-0.016	0.048	0.071	1.000
$\lambda_{21}$	0.8	0.024	0.066	0.091	1.000	0.001	0.054	0.076	1.000
$\lambda_{31}^{21}$	0.8	0.071	0.094	0.092	1.000	0.036	0.067	0.076	1.000
$\lambda_{41}$	0.3	0.124	0.156	0.087	1.000	0.071	0.107	0.109	0.960
λ <sub>51</sub>	0.5	0.008	0.077	0.089	1.000	0.022	0.069	0.090	1.000
$\lambda_{61}$	0.6	0.006	0.055	0.092	1.000	0.018	0.056	0.080	1.000
$\lambda_{71}$	0.3	-0.017	0.063	0.108	0.870	-0.060	0.084	0.087	0.890
$\lambda_{18,1}$	0.3	-0.004	0.046	0.132	0.620	-0.046	0.061	0.090	0.930
$\lambda_{62}$	0.3	-0.018	0.058	0.099	0.950	-0.068	0.087	0.089	0.900
$\lambda_{72}$	0.6	-0.019	0.058	0.091	0.990	0.017	0.056	0.089	1.000
$\lambda_{82}$	0.8	-0.002	0.069	0.097	0.990	-0.019	0.050	0.085	1.000
$\lambda_{02}$	0.3	0.191	0.211	0.091	0.990	0.107	0.129	0.130	0.980
$\lambda_{10,2}$	0.8	0.070	0.101	0.098	0.990	0.036	0.067	0.086	1.000
$\lambda_{11,2}$	0.5	-0.022	0.083	0.087	0.990	0.007	0.066	0.097	1.000
$\lambda_{12,2}$	0.6	-0.027	0.067	0.097	1.000	0.027	0.064	0.093	1.000
$\lambda_{13,2}$	0.3	-0.067	0.084	0.135	0.130	-0.058	0.071	0.090	0.900
$\lambda_{13}$	0.3	-0.086	0.100	0.137	0.070	-0.022	0.053	0.083	0.960
$\lambda_{123}$	0.3	0.017	0.057	0.116	0.950	-0.051	0.075	0.097	0.810
$\lambda_{13,3}$	0.8	0.019	0.062	0.117	0.990	0.011	0.055	0.079	1.000
$\lambda_{14,3}$	0.3	-0.024	0.092	0.088	0.870	0.004	0.070	0.099	0.940
$\lambda_{15,3}$	0.5	0.060	0.099	0.103	0.990	0.046	0.083	0.099	1.000
$\lambda_{16,3}$	0.6	0.051	0.094	0.107	0.990	0.042	0.084	0.093	1.000
$\lambda_{17,3}$	0.5	-0.032	0.084	0.097	0.990	0.018	0.063	0.097	1.000
$\lambda_{18,3}$	0.8	-0.024	0.065	0.120	0.990	0.003	0.051	0.080	1.000
$\lambda_0$	0	-0.024	0.072	0.110	0.012	-0.013	0.034	0.073	0.000
$\phi_{\nu\nu'}$	0.3	0.067	0.094	0.168	0.670	0.069	0.095	0.118	0.973
$\psi_{ii}$	0.53	-0.010	0.066	0.054	1.000	0.014	0.064	0.077	1.000
Ψ	0.3	_				-0.056	0.074	0.082	0.903
$\psi_h$	0.3	_				-0.038	0.058	0.063	1.000
ψ	0	_	_		_	-0.001	0.028	0.046	0.003

Table D2						
C-step Results	With	Three	Specified	Loadings	per Fac	tor

			N = 2	250			N = 1	500	
Par	True	BIAS	RMSE	SE	SIG%	BIAS	RMSE	SE	SIG%
λ <sub>11</sub>	0.7	-0.064	0.106	0.259	0.575	-0.052	0.121	0.239	0.650
$\lambda_{21}$	0.7	-0.064	0.107	0.260	0.600	-0.050	0.119	0.239	0.680
$\lambda_{31}^{21}$	0.7	-0.312	0.324	0.257	0.100	-0.287	0.310	0.243	0.290
λ <sub>41</sub>	0.7	-0.429	0.438	0.227	0.008	-0.389	0.407	0.220	0.110
$\lambda_{51}$	0.5	-0.300	0.309	0.184	0.000	-0.275	0.290	0.176	0.080
λ <sub>61</sub>	0.5	-0.330	0.336	0.183	0.000	-0.305	0.318	0.177	0.030
λ <sub>52</sub>	0.5	-0.307	0.319	0.220	0.000	-0.268	0.300	0.207	0.120
λ <sub>62</sub>	0.5	-0.245	0.263	0.264	0.000	-0.211	0.248	0.240	0.130
λ <sub>72</sub>	0.7	-0.360	0.376	0.305	0.000	-0.329	0.365	0.283	0.100
λ <sub>82</sub>	0.7	-0.376	0.387	0.306	0.008	-0.340	0.372	0.280	0.120
$\lambda_{92}$	0.7	-0.497	0.508	0.253	0.000	-0.449	0.480	0.240	0.080
$\lambda_{10,2}$	0.7	-0.496	0.507	0.254	0.000	-0.447	0.479	0.239	0.070
$\lambda_0$	0	0.009	0.046	0.157	0.000	0.007	0.051	0.155	0.000
φ <sub>12</sub>	0.3	-0.119	0.161	0.411	0.000	-0.073	0.162	0.386	0.080
ψ <sub>55</sub>	0.35	-0.023	0.092	0.161	0.283	-0.044	0.137	0.144	0.300
Ψ66	0.35	-0.025	0.084	0.165	0.275	-0.051	0.137	0.146	0.310
$\psi_{ii}$	0.51	-0.221	0.322	0.140	0.299	-0.237	0.338	0.126	0.351
Ψ̈́	0	0.326	0.338	0.165	0.316	0.306	0.328	0.147	0.364

*Note.*  $\psi_{jj}$  averaged across elements from j = 1 to 4 and 7 to 10;  $\lambda_0$  averaged across all zero loading estimates;  $\psi_0$  averaged across all zero residual covariance estimates.

loading matrix, additional simulation studies were conducted. The model settings were the same as Study 1 (10 items, two factors) except that the data were generated with diagonal  $\Psi$  (i.e., the residuals in the true model were assumed to be independent). In the E-step, the model was fitted with one specified loading per factor (that is,  $\lambda_{11}$  for Factor 1 and  $\lambda_{10,2}$  for Factor 2) or no specified loadings (that is, all loadings were unspeci-

fied). The results, found in Tables D5 and D6, can be compared with those in Table D3. The estimates in the case of no specified loading (Table D6) were poor. In contrast, the case of one specified loading per factor (Table D5) was acceptable, but slightly worse than the case of two specified loadings per factor (Table D3). This is consistent with the findings in Study 1 (with correlated residuals).

Table D3 E-Step Results With True Diagonal  $\psi$  and Two Specified Loadings per Factor (10 Items, Two Factors)

			N = 250			N = 500	
Par	True	BIAS	RMSE	SE	BIAS	RMSE	SE
λ <sub>11</sub>	0.7	0.021	0.078	0.134	0.004	0.068	0.118
$\lambda_{21}$	0.7	0.010	0.070	0.132	0.006	0.066	0.119
λ <sub>31</sub>	0.7	-0.001	0.071	0.130	0.001	0.068	0.118
$\lambda_{41}$	0.7	0.003	0.073	0.131	-0.013	0.065	0.114
λ <sub>51</sub>	0.5	-0.019	0.065	0.145	-0.017	0.051	0.130
λ <sub>61</sub>	0.5	-0.025	0.071	0.144	-0.021	0.053	0.132
λ <sub>52</sub>	0.5	-0.018	0.063	0.144	-0.025	0.053	0.130
λ <sub>62</sub>	0.5	-0.019	0.060	0.143	-0.012	0.051	0.130
λ <sub>72</sub>	0.7	0.004	0.058	0.132	0.005	0.066	0.118
λ <sub>82</sub>	0.7	0.000	0.071	0.131	0.003	0.071	0.117
λ <sub>92</sub>	0.7	0.015	0.069	0.133	0.009	0.067	0.118
$\lambda_{10,2}$	0.7	0.005	0.063	0.130	0.003	0.067	0.118
λο	0	-0.033	0.069	0.177	-0.027	0.066	0.168
φ <sub>12</sub>	0.3	0.075	0.094	0.219	0.075	0.087	0.211
$\psi_{55}$	0.35	0.001	0.040	0.041	0.006	0.029	0.029
$\psi_{66}$	0.35	0.012	0.038	0.042	0.006	0.028	0.029
$\psi_{jj}$	0.51	0.007	0.058	0.059	0.006	0.042	0.041

Note.  $\lambda_0$  averaged across all zero loading estimates.

Table D4									
E-Step Results	With True	Diagonal	and Two	o Specified	Loadings	per Factor	(18 Items,	Three	Factors)

			<i>N</i> = 250			N = 500	
Par	True	BIAS	RMSE	SE	BIAS	RMSE	SE
λ <sub>11</sub>	0.5	0.023	0.066	0.105	0.018	0.045	0.087
$\lambda_{21}$	0.7	0.023	0.065	0.094	0.024	0.062	0.077
λ <sub>31</sub>	0.7	0.014	0.068	0.093	0.027	0.061	0.077
$\lambda_{41}$	0.7	0.007	0.057	0.092	0.013	0.052	0.076
λ <sub>51</sub>	0.7	0.020	0.060	0.093	0.015	0.051	0.077
λ <sub>61</sub>	0.5	-0.001	0.057	0.108	-0.009	0.045	0.092
$\lambda_{71}$	0.5	-0.005	0.056	0.107	-0.005	0.048	0.092
λ <sub>18.1</sub>	0.5	0.000	0.062	0.105	0.013	0.042	0.086
$\lambda_{62}$	0.5	-0.007	0.055	0.109	-0.009	0.043	0.093
$\lambda_{72}$	0.5	-0.012	0.068	0.108	-0.010	0.043	0.093
$\lambda_{82}$	0.7	0.033	0.074	0.099	0.021	0.052	0.080
$\lambda_{92}$	0.7	0.014	0.065	0.097	0.016	0.049	0.080
λ <sub>10.2</sub>	0.7	0.016	0.069	0.097	0.006	0.045	0.079
$\lambda_{11,2}$	0.7	0.009	0.057	0.097	0.011	0.048	0.079
$\lambda_{12,2}$	0.5	-0.020	0.061	0.108	-0.005	0.041	0.091
λ <sub>13.2</sub>	0.5	-0.009	0.056	0.108	-0.003	0.047	0.091
λ <sub>13</sub>	0.5	-0.016	0.064	0.110	-0.019	0.044	0.092
λ <sub>12.3</sub>	0.5	-0.009	0.064	0.106	-0.012	0.047	0.090
$\lambda_{13,3}$	0.5	-0.010	0.061	0.106	-0.011	0.048	0.091
$\lambda_{14,3}$	0.7	0.006	0.063	0.095	0.003	0.062	0.080
$\lambda_{15,3}$	0.7	0.018	0.067	0.096	0.013	0.059	0.080
$\lambda_{16,3}$	0.7	0.022	0.068	0.097	0.014	0.068	0.080
λ <sub>17.3</sub>	0.7	-0.003	0.059	0.094	0.004	0.054	0.080
λ <sub>18.3</sub>	0.5	-0.016	0.063	0.108	-0.024	0.049	0.092
λ	0	-0.019	0.052	0.114	-0.018	0.045	0.101

Note.  $\lambda_0$  averaged across all zero loading estimates.

Additional simulation studies were also conducted for a more complex loading structure in Study 2 (18 items, three factors). The model settings were the same as those of Study 2 except that the data were generated with diagonal  $\Psi$ . In the E-step, the model was

fitted with one specified loading per factor (that is,  $\lambda_{11}$  for Factor 1,  $\lambda_{82}$  for Factor 2, and  $\lambda_{15,3}$  for Factor 3) or no specified loadings (that is, all loadings were unspecified). The results can be found in Table D7 and D8 and can be compared with those in Table D4.

Table D5E-Step Results With One Specified Loading per Factor (10 Items, Two Factors)

			N = 250			N = 500	
Par	True	BIAS	RMSE	SE	BIAS	RMSE	SE
λ <sub>11</sub>	0.7	-0.036	0.147	0.193	-0.023	0.114	0.149
$\lambda_{21}$	0.7	-0.057	0.143	0.187	-0.025	0.106	0.150
λ <sub>31</sub>	0.7	-0.056	0.148	0.187	-0.022	0.115	0.148
$\lambda_{41}$	0.7	-0.054	0.141	0.190	-0.038	0.111	0.146
λ <sub>51</sub>	0.5	-0.011	0.059	0.159	-0.017	0.053	0.145
λ <sub>61</sub>	0.5	-0.018	0.069	0.159	-0.021	0.053	0.147
λ <sub>52</sub>	0.5	-0.023	0.063	0.160	-0.025	0.053	0.146
λ <sub>62</sub>	0.5	-0.023	0.066	0.158	-0.012	0.051	0.146
$\lambda_{72}$	0.7	-0.056	0.150	0.188	-0.017	0.099	0.145
λ <sub>82</sub>	0.7	-0.061	0.155	0.187	-0.020	0.100	0.145
$\lambda_{92}$	0.7	-0.061	0.160	0.190	-0.020	0.096	0.144
$\lambda_{10,2}$	0.7	-0.058	0.151	0.188	-0.020	0.102	0.144
$\lambda_0^{10,2}$	0	0.029	0.142	0.226	-0.002	0.102	0.197

(Appendices continue)

Table D6									
E-Step Results	With All	Loadings	Being	Unspecified	(10	Items,	Two	Factors)	

			N = 250			N = 500	
Par	True	BIAS	RMSE	SE	BIAS	RMSE	SE
λ <sub>11</sub>	0.7	-0.385	0.474	0.289	-0.361	0.464	0.269
λ <sub>21</sub>	0.7	-0.378	0.462	0.280	-0.352	0.460	0.269
λ <sub>31</sub>	0.7	-0.380	0.464	0.281	-0.356	0.462	0.273
$\lambda_{41}$	0.7	-0.381	0.466	0.286	-0.356	0.454	0.264
λ <sub>51</sub>	0.5	-0.017	0.063	0.173	-0.034	0.069	0.166
λ <sub>61</sub>	0.5	-0.022	0.065	0.173	-0.030	0.061	0.168
λ <sub>52</sub>	0.5	-0.018	0.063	0.173	-0.012	0.058	0.163
λ <sub>62</sub>	0.5	-0.021	0.059	0.173	-0.007	0.057	0.165
$\lambda_{72}$	0.7	-0.384	0.470	0.290	-0.334	0.432	0.265
λ <sub>82</sub>	0.7	-0.380	0.467	0.285	-0.335	0.433	0.263
λ <sub>92</sub>	0.7	-0.384	0.469	0.290	-0.334	0.436	0.262
λ <sub>10.2</sub>	0.7	-0.391	0.473	0.283	-0.337	0.433	0.260
λ <sub>0</sub>	0	0.355	0.445	0.283	0.317	0.425	0.267

Note.  $\lambda_0$  averaged across all zero loading estimates.

Similar to the above, the estimates in the case of no specified loading (Table D8) were poor. In contrast, the case of one specified loading per factor (Table D7) was acceptable, but slightly worse than the case of two specified loadings per factor (Table D4) in general.

or more specified loadings per factor is preferred for more stable performance in the E-step.

The same simulated data sets were also used to investigate the same issue with the C-step. Table D2 above revealed that the C-step performed poorly up to three specified loadings per factor. In contrast, Table D9 confirmed that the constraint of one specified

In sum, the case of one specified loading per factor was acceptable in general and can serve as the minimum constraint, while two

Table D7								
E-Step Results	With One	Specified	Loading p	er Faci	tor (18 I	tems,	Three	Factors)

			N = 250			N = 500	
Par	True	BIAS	RMSE	SE	BIAS	RMSE	SE
λ <sub>11</sub>	0.5	0.015	0.086	0.133	0.012	0.077	0.102
$\lambda_{21}$	0.7	-0.072	0.219	0.130	-0.027	0.158	0.095
$\lambda_{31}^{21}$	0.7	-0.071	0.223	0.131	-0.018	0.159	0.097
$\lambda_{41}$	0.7	-0.081	0.224	0.132	-0.028	0.154	0.094
λ <sub>51</sub>	0.7	-0.068	0.218	0.135	-0.030	0.162	0.097
λ <sub>61</sub>	0.5	-0.049	0.153	0.127	-0.031	0.104	0.105
λ <sub>71</sub>	0.5	-0.052	0.153	0.127	-0.028	0.111	0.105
λ <sub>18.1</sub>	0.5	-0.009	0.087	0.131	0.008	0.071	0.100
$\lambda_{62}$	0.5	-0.019	0.082	0.125	-0.009	0.044	0.102
$\lambda_{72}$	0.5	-0.024	0.096	0.125	-0.011	0.048	0.103
$\lambda_{82}$	0.7	-0.018	0.180	0.126	0.002	0.120	0.092
$\lambda_{92}$	0.7	-0.047	0.175	0.122	-0.010	0.117	0.090
λ <sub>10.2</sub>	0.7	-0.032	0.183	0.124	-0.015	0.122	0.091
$\lambda_{11,2}$	0.7	-0.040	0.176	0.123	-0.008	0.119	0.090
$\lambda_{12,2}$	0.5	-0.040	0.117	0.126	-0.017	0.082	0.106
$\lambda_{13,2}$	0.5	-0.031	0.115	0.126	-0.014	0.086	0.107
λ <sub>13</sub>	0.5	-0.047	0.118	0.128	-0.029	0.074	0.103
λ <sub>12.3</sub>	0.5	-0.057	0.151	0.128	-0.035	0.118	0.104
λ <sub>13.3</sub>	0.5	-0.054	0.147	0.128	-0.034	0.113	0.105
λ <sub>14.3</sub>	0.7	-0.080	0.230	0.134	-0.030	0.159	0.093
λ <sub>15.3</sub>	0.7	-0.068	0.234	0.132	-0.028	0.173	0.095
λ <sub>16.3</sub>	0.7	-0.081	0.236	0.132	-0.028	0.173	0.093
λ <sub>17.3</sub>	0.7	-0.088	0.226	0.130	-0.033	0.171	0.093
λ <sub>18.3</sub>	0.5	-0.044	0.115	0.125	-0.032	0.072	0.102
λ <sub>0</sub>	0	0.020	0.142	0.138	0.000	0.101	0.115

Table D8								
E-Step Results	With All	Loadings	Being	Unspecified	(18	Items,	Three	Factors)

			N = 250			N = 500	
Par	True	BIAS	RMSE	SE	BIAS	RMSE	SE
λ <sub>11</sub>	0.5	-0.173	0.275	0.159	-0.169	0.293	0.116
$\lambda_{21}$	0.7	-0.452	0.557	0.165	-0.467	0.578	0.131
$\lambda_{31}^{21}$	0.7	-0.446	0.555	0.168	-0.462	0.575	0.132
$\lambda_{41}^{51}$	0.7	-0.461	0.565	0.167	-0.470	0.578	0.132
$\lambda_{51}$	0.7	-0.453	0.561	0.170	-0.474	0.579	0.131
λ <sub>61</sub>	0.5	-0.159	0.276	0.161	-0.191	0.304	0.121
$\lambda_{71}$	0.5	-0.179	0.290	0.159	-0.191	0.306	0.121
λ <sub>18.1</sub>	0.5	-0.172	0.273	0.158	-0.167	0.288	0.115
$\lambda_{62}$	0.5	-0.191	0.295	0.160	-0.178	0.301	0.123
$\lambda_{72}$	0.5	-0.179	0.287	0.158	-0.175	0.301	0.123
$\lambda_{82}$	0.7	-0.451	0.559	0.178	-0.484	0.589	0.131
$\lambda_{92}$	0.7	-0.449	0.543	0.176	-0.476	0.579	0.130
$\lambda_{10,2}$	0.7	-0.441	0.553	0.177	-0.481	0.584	0.129
$\lambda_{11,2}$	0.7	-0.439	0.546	0.176	-0.479	0.581	0.131
$\lambda_{12,2}$	0.5	-0.164	0.270	0.153	-0.178	0.297	0.121
$\lambda_{13,2}$	0.5	-0.161	0.268	0.152	-0.177	0.297	0.123
λ <sub>13</sub>	0.5	-0.164	0.277	0.158	-0.195	0.313	0.118
λ <sub>12.3</sub>	0.5	-0.196	0.300	0.154	-0.174	0.291	0.121
$\lambda_{13,3}$	0.5	-0.188	0.295	0.155	-0.172	0.294	0.122
$\lambda_{14,3}$	0.7	-0.463	0.567	0.172	-0.490	0.586	0.134
$\lambda_{15,3}$	0.7	-0.475	0.568	0.175	-0.487	0.586	0.133
$\lambda_{16,3}$	0.7	-0.452	0.552	0.177	-0.491	0.588	0.134
λ <sub>17.3</sub>	0.7	-0.467	0.561	0.175	-0.488	0.588	0.132
λ <sub>18.3</sub>	0.5	-0.179	0.288	0.156	-0.201	0.320	0.116
λ	0	0.230	0.374	0.174	0.245	0.405	0.127

Note.  $\lambda_0$  averaged across all zero loading estimates.

loading per item was still appropriate even when the residuals in the true model were assumed to be independent.

Additional simulation studies were conducted to compare the performance of the E-step with the BSEM approach using the ridge regression priors proposed by Muthén and Asparouhov (2012) in estimating the loadings. The BSEM approach relies on normal priors concentrated around zero with small variance. Mplus was used to analyze the same simulated data sets with diagonal  $\Psi$ . The results of loading estimates are presented in Tables D10 and D11. Compared with Tables D3 and D4, the E-step with the

Table D9C-Step Results With One Specified Loading per Item (10 Items, Two Factors)

			<i>N</i> = 250		N = 500				
Par	True	BIAS	RMSE	SE	BIAS	RMSE	SE		
λ <sub>11</sub>	0.7	-0.015	0.066	0.119	-0.020	0.047	0.096		
$\lambda_{21}$	0.7	-0.021	0.063	0.117	-0.018	0.046	0.095		
λ <sub>31</sub>	0.7	-0.022	0.065	0.118	-0.019	0.052	0.096		
$\lambda_{41}$	0.7	-0.018	0.067	0.117	-0.028	0.046	0.095		
$\lambda_{51}$	0.5	0.016	0.059	0.120	0.004	0.040	0.101		
$\lambda_{61}$	0.5	-0.089	0.110	0.137	-0.067	0.079	0.116		
$\lambda_{52}$	0.5	-0.083	0.102	0.137	-0.071	0.080	0.116		
$\lambda_{62}$	0.5	0.017	0.059	0.121	0.010	0.040	0.102		
$\lambda_{72}$	0.7	-0.022	0.055	0.118	-0.014	0.048	0.098		
$\lambda_{82}$	0.7	-0.021	0.064	0.119	-0.016	0.048	0.097		
$\lambda_{92}$	0.7	-0.020	0.062	0.119	-0.016	0.044	0.097		
$\lambda_{10,2}$	0.7	-0.029	0.063	0.116	-0.020	0.050	0.096		
λ	0	-0.012	0.034	0.105	-0.013	0.027	0.094		

Note.  $\lambda_0$  averaged across all zero loading estimates.

Loading	Loading Estimates for BSEM With the Ridge Regression Priors (10 Items, Two Factors)										
		N = 250					N = 500				
Par	True	BIAS	RMSE	SE	SIG%	BIAS	RMSE	SE			
λ <sub>11</sub>	0.7	-0.199	0.220	0.111	1.000	0.016	0.055	0.057			
λ <sub>21</sub>	0.7	-0.193	0.211	0.112	1.000	0.019	0.052	0.057			
λ <sub>31</sub>	0.7	-0.459	0.466	0.076	0.990	-0.119	0.130	0.047			
λ <sub>41</sub>	0.7	-0.456	0.464	0.076	1.000	-0.129	0.137	0.046			
λ <sub>51</sub>	0.5	-0.156	0.158	0.073	1.000	-0.1	0.105	0.050			
λ <sub>61</sub>	0.5	-0.161	0.162	0.072	1.000	-0.104	0.107	0.050			
λ <sub>52</sub>	0.5	-0.158	0.160	0.073	1.000	-0.113	0.116	0.049			
λ <sub>62</sub>	0.5	-0.160	0.162	0.073	1.000	-0.103	0.107	0.050			

0.076

0.076

0.113

0.111

0.087

1.000

1.000

1.000

1.000

0.446

-0.116

-0.12

0.02

0.013

-0.087

 Table D10

 Loading Estimates for BSEM With the Ridge Regression Priors (10 Items, Two Factors)

0.468

0.467

0.211

0.215

0.125

Bayesian Lasso was better in terms of BIAS and RMSE. More importantly, the Type I error rates for the BSEM approach were high, making it likely to misidentify truly zero loadings as significant. To better understand whether the bias in the E-step is a function

-0.460

-0.460

-0.192

-0.201

0.090

of the residual correlations, additional simulation studies were conducted. The settings were similar to those of Study 1, except that the data were generated with two additional level of residual correlations (0.3 and 0.7). Comparing the results (Tables D12 and D13) with the case of no correlated residual (Table D3), it seems that the bias was a function of the residual correlations to some extent, and the bias became larger when the residual correlation increased.

0.126

0.130

0.052

0.052

0.100

0.047

0.046

0.057

0.057

0.060

Table D11Loading Estimates for BSEM With the Ridge Regression Priors (18 Items, Three Factors)

			N = 250				N = 500			
Par	True	BIAS	RMSE	SE	SIG%	BIAS	RMSE	SE	SIG%	
λ <sub>11</sub>	0.5	0.156	0.187	0.120	1.000	-0.008	0.044	0.061	1.000	
$\lambda_{21}$	0.7	-0.133	0.188	0.117	1.000	0.006	0.048	0.056	1.000	
λ <sub>31</sub>	0.7	-0.412	0.439	0.076	0.729	-0.105	0.111	0.046	1.000	
λ <sub>41</sub>	0.7	-0.415	0.440	0.076	0.708	-0.117	0.122	0.046	1.000	
λ <sub>51</sub>	0.7	-0.406	0.434	0.076	0.708	-0.118	0.123	0.046	1.000	
λ <sub>61</sub>	0.5	-0.260	0.270	0.075	0.958	-0.110	0.114	0.050	1.000	
λ <sub>71</sub>	0.5	-0.266	0.275	0.075	0.917	-0.106	0.112	0.051	1.000	
$\lambda_{18,1}$	0.5	-0.181	0.200	0.076	1.000	-0.065	0.073	0.054	1.000	
λ <sub>62</sub>	0.5	-0.202	0.217	0.067	1.000	-0.070	0.076	0.050	1.000	
λ <sub>72</sub>	0.5	-0.206	0.224	0.067	0.979	-0.073	0.079	0.051	1.000	
$\lambda_{82}$	0.7	-0.056	0.156	0.103	1.000	0.045	0.063	0.058	1.000	
$\lambda_{92}$	0.7	-0.064	0.145	0.102	1.000	0.038	0.061	0.058	1.000	
$\lambda_{10,2}$	0.7	-0.314	0.356	0.072	0.708	-0.098	0.103	0.047	1.000	
λ <sub>11.2</sub>	0.7	-0.269	0.291	0.057	1.000	-0.122	0.124	0.038	1.000	
λ <sub>12.2</sub>	0.5	-0.069	0.131	0.057	1.000	0.078	0.082	0.038	1.000	
λ <sub>13.2</sub>	0.5	-0.160	0.187	0.067	1.000	-0.010	0.040	0.047	1.000	
λ <sub>13</sub>	0.5	-0.393	0.415	0.096	0.292	-0.137	0.140	0.055	1.000	
λ <sub>12,3</sub>	0.5	-0.312	0.323	0.073	0.604	-0.182	0.184	0.048	1.000	
λ <sub>13.3</sub>	0.5	-0.292	0.305	0.075	0.667	-0.158	0.161	0.048	1.000	
λ <sub>14.3</sub>	0.7	-0.432	0.456	0.074	0.667	-0.148	0.153	0.045	1.000	
λ <sub>15,3</sub>	0.7	-0.184	0.215	0.116	1.000	-0.041	0.057	0.053	1.000	
$\lambda_{16,3}$	0.7	-0.174	0.207	0.116	1.000	-0.041	0.071	0.053	1.000	
λ <sub>17,3</sub>	0.7	-0.434	0.456	0.074	0.667	-0.144	0.148	0.044	1.000	
λ <sub>18.3</sub>	0.5	-0.197	0.211	0.076	1.000	-0.131	0.135	0.053	1.000	
λο	0	0.026	0.105	0.079	0.181	-0.053	0.073	0.057	0.188	

Note.  $\lambda_0$  averaged across all zero loading estimates.

 $\lambda_{72}$ 

 $\lambda_{82}$ 

 $\lambda_{92}$ 

λ

 $\lambda_{10,2}$ 

0.7

0.7

0.7

0.7

0

SIG%

1.000

1.000

1.000

1.000

1.000

1.000

1.000

1.000

1.000

1.000

1.000

1.000

0.334

Table D12								
E-Step Results	With Two	Specified	Loadings	per	Factor	(Residual	Correlation	= .3)

			N = 250			N = 500	
Par	True	BIAS	RMSE	SE	BIAS	RMSE	SE
λ <sub>11</sub>	0.7	0.102	0.123	0.122	0.110	0.121	0.107
λ <sub>21</sub>	0.7	0.113	0.137	0.123	0.107	0.116	0.107
λ <sub>31</sub>	0.7	-0.035	0.069	0.110	-0.028	0.062	0.093
$\lambda_{41}$	0.7	-0.066	0.095	0.109	-0.054	0.073	0.091
λ <sub>51</sub>	0.5	-0.043	0.072	0.139	-0.050	0.068	0.125
λ <sub>61</sub>	0.5	-0.053	0.081	0.140	-0.052	0.069	0.124
λ <sub>52</sub>	0.5	-0.007	0.059	0.135	0.003	0.047	0.116
λ <sub>62</sub>	0.5	0.002	0.059	0.133	0.001	0.046	0.116
$\lambda_{72}$	0.7	-0.005	0.068	0.112	0.011	0.047	0.091
λ <sub>82</sub>	0.7	0.005	0.066	0.114	0.017	0.049	0.091
$\lambda_{92}$	0.7	0.026	0.079	0.116	0.026	0.056	0.093
$\lambda_{10,2}$	0.7	0.028	0.071	0.117	0.026	0.060	0.093
$\lambda_0$	0	-0.039	0.078	0.163	-0.039	0.070	0.150
φ <sub>12</sub>	0.3	0.083	0.101	0.209	0.090	0.100	0.197
$\psi_{55}$	0.35	0.017	0.045	0.042	0.016	0.034	0.029
$\psi_{66}$	0.35	0.014	0.044	0.042	0.016	0.034	0.029
$\psi_{ii}$	0.51	-0.007	0.080	0.057	-0.012	0.065	0.040

To better understand whether the bias in the E-step is a function of the sparsity of the residual covariance matrix, additional simulation studies were conducted. The settings were similar to those of Study 1 except that the data were generated with four residual correlations each with a magnitude of 0.3. Comparing the results (Table D14) with the case of two residual correlations each with the same magnitude (Table D12), it seems the difference was not significant. More work is needed to systematically investigate this issue.

To better understand whether the bias in the E-step diminishes when the cross-loadings are close to zero, additional simulation studies were conducted. The settings were similar to those of Study 1 except that the data were generated with residual correlations of 0.3 and no cross-loadings. Comparing the results in Table D15 with those in Table D12 above, the bias seemed to diminish a little bit but still existed when the cross-loadings were closed to zero.

To better understand how the Bayesian Lasso works when the aim is to estimate the residual covariance matrix, additional simulation studies were conducted. The settings were similar as those of Study 1 and 2 except that the loading structure was assumed to be known in the C-step (i.e., all the nonzero load-

 Table D13

 E-Step Results With Two Specified Loadings per Factor (Residual Correlation = .7)

			N = 250			N = 500	= 500	
Par	True	BIAS	RMSE	SE	BIAS	RMSE	SE	
λ <sub>11</sub>	0.7	0.249	0.255	0.114	0.244	0.252	0.108	
$\lambda_{21}$	0.7	0.247	0.253	0.113	0.245	0.254	0.108	
λ <sub>31</sub>	0.7	-0.196	0.207	0.091	-0.193	0.201	0.077	
$\lambda_{41}$	0.7	-0.233	0.248	0.095	-0.224	0.231	0.083	
λ <sub>51</sub>	0.5	-0.195	0.204	0.131	-0.193	0.198	0.124	
λ <sub>61</sub>	0.5	-0.187	0.195	0.130	-0.185	0.192	0.124	
λ <sub>52</sub>	0.5	0.100	0.112	0.098	0.107	0.116	0.087	
λ <sub>62</sub>	0.5	0.098	0.114	0.099	0.105	0.114	0.088	
λ <sub>72</sub>	0.7	0.019	0.070	0.089	0.033	0.069	0.079	
$\lambda_{82}$	0.7	-0.008	0.062	0.088	0.000	0.057	0.078	
λ <sub>92</sub>	0.7	0.002	0.061	0.089	-0.002	0.058	0.078	
λ <sub>10.2</sub>	0.7	0.014	0.069	0.090	0.006	0.056	0.079	
λ	0	-0.003	0.113	0.136	-0.001	0.117	0.131	
$\phi_{12}$	0.3	0.058	0.081	0.187	0.062	0.079	0.184	
ψ <sub>55</sub>	0.35	0.056	0.070	0.045	0.043	0.054	0.031	
$\psi_{66}$	0.35	0.052	0.069	0.045	0.045	0.054	0.031	
$\psi_{jj}$	0.51	-0.036	0.161	0.053	-0.042	0.149	0.037	

			N = 250			N = 500	
Par	True	BIAS	RMSE	SE	BIAS	RMSE	SE
λ <sub>11</sub>	0.7	0.095	0.121	0.125	0.094	0.109	0.109
$\lambda_{21}$	0.7	0.090	0.115	0.125	0.083	0.102	0.109
λ <sub>31</sub>	0.7	-0.024	0.079	0.115	-0.020	0.053	0.098
$\lambda_{41}$	0.7	-0.051	0.077	0.113	-0.048	0.074	0.096
λ <sub>51</sub>	0.5	-0.066	0.091	0.137	-0.075	0.090	0.125
$\lambda_{61}$	0.5	-0.055	0.080	0.135	-0.071	0.084	0.125
λ <sub>52</sub>	0.5	-0.003	0.061	0.129	0.008	0.050	0.114
$\lambda_{62}$	0.5	-0.015	0.068	0.129	0.008	0.048	0.114
$\lambda_{72}$	0.7	0.005	0.066	0.114	0.015	0.057	0.099
$\lambda_{82}$	0.7	0.035	0.082	0.120	0.048	0.073	0.103
$\lambda_{92}$	0.7	0.002	0.072	0.114	-0.006	0.054	0.096
$\lambda_{10,2}$	0.7	-0.010	0.071	0.113	-0.008	0.060	0.096
$\lambda_0$	0	-0.028	0.069	0.163	-0.030	0.063	0.151
φ <sub>12</sub>	0.3	0.107	0.121	0.205	0.104	0.114	0.195
$\psi_{55}$	0.35	0.031	0.053	0.043	0.028	0.041	0.030
ψ <sub>66</sub>	0.35	0.034	0.056	0.043	0.033	0.045	0.030
$\psi_{ii}$	0.51	-0.008	0.076	0.058	-0.010	0.065	0.041

 Table D14

 E-Step With Bayesian Lasso in Estimating the Loadings (Four Residual Correlations Each With a Magnitude of 0.3)

*Note.*  $\psi_{ij}$  averaged across elements from j = 1 to 4 and 7 to 10;  $\lambda_0$  averaged across all zero loading.

ings including the cross-loadings were assumed to be specified and all the zero loadings were fixed to zero). The results in Table D16 showed that, by assuming the loading structure as known, the power of detecting the within-factor nonzero residual covariance was still low, but higher than that with an unknown structure in Table 4 (0.020 vs. 0.005 in sample size 250; 0.200 vs. 0.040 in sample size 500). The power of detecting the between-factor nonzero residual covariance was satisfactory and higher than that in Table 4 (0.830 vs. 0.755 in sample size 250; 0.980 vs. 0.915 in sample size 500). It is interesting to observe that, similar to Table 4, the estimates of  $\lambda_{11}$  and  $\lambda_{21}$  were positively biased and the estimate of  $\psi_{21}$  was negatively biased. Moreover, the estimate of the between-factor residual covariance  $\psi_{74}$  was also more accurate and stable than that of the within-factor residual covariance  $\psi_{21}$ . Accordingly, the cases of known and unknown loading structures in the C-step exhibited similar patterns, and the power of detecting the nonzero off-diagonal elements of the covariance matrix was not only related to the Bayesian Lasso estimation, but also related to the position of the nonzero elements (that is, within factor or

Table D15E-Step With Bayesian Lasso Priors in Estimating the Loadings (No Cross-Loading)

			<i>N</i> = 250			N = 500	
Par	True	BIAS	RMSE	SE	BIAS	RMSE	SE
λ <sub>11</sub>	0.7	0.092	0.110	0.076	0.097	0.106	0.055
$\lambda_{21}$	0.7	0.091	0.107	0.076	0.098	0.108	0.055
λ <sub>31</sub>	0.7	-0.068	0.094	0.075	-0.045	0.067	0.054
$\lambda_{41}$	0.7	-0.067	0.094	0.077	-0.065	0.078	0.055
λ <sub>51</sub>	0.5	-0.042	0.089	0.075	-0.041	0.058	0.053
λ <sub>62</sub>	0.5	-0.023	0.057	0.075	-0.004	0.045	0.054
λ <sub>72</sub>	0.7	-0.007	0.070	0.074	-0.005	0.052	0.054
λ <sub>82</sub>	0.7	-0.019	0.066	0.074	-0.003	0.050	0.053
$\lambda_{92}$	0.7	-0.003	0.067	0.074	0.008	0.041	0.054
λ <sub>10.2</sub>	0.7	0.013	0.063	0.075	0.003	0.043	0.054
$\lambda_0$	0	0.011	0.058	0.104	0.008	0.042	0.090
$\phi_{12}$	0.3	-0.023	0.073	0.167	-0.022	0.059	0.152
ψ55	0.75	0.017	0.076	0.074	0.029	0.066	0.053
$\psi_{66}$	0.75	0.010	0.076	0.074	0.010	0.051	0.052
$\psi_{jj}$	0.51	-0.011	0.085	0.060	-0.014	0.073	0.042

*Note.*  $\psi_{ij}$  averaged across elements from j = 1 to 4 and 7 to 10;  $\lambda_0$  averaged across all zero loading estimates.

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Table D16				
C-Step With	Known Loading	Structure (10	Items, Two	Factors)

			N = 1	250			N = 1	500	
Par	True	BIAS	RMSE	SE	SIG%	BIAS	RMSE	SE	SIG%
λ	0.7	0.096	0.120	0.097	1.000	0.086	0.096	0.084	1.000
λ <sub>21</sub>	0.7	0.098	0.121	0.098	1.000	0.089	0.100	0.084	1.000
$\lambda_{21}^{21}$	0.7	-0.042	0.072	0.097	1.000	-0.029	0.055	0.080	1.000
$\lambda_{41}$	0.7	-0.085	0.102	0.083	1.000	-0.075	0.086	0.067	1.000
$\lambda_{51}$	0.5	-0.026	0.065	0.088	1.000	-0.036	0.055	0.069	1.000
$\lambda_{61}$	0.5	-0.035	0.064	0.087	1.000	-0.033	0.051	0.068	1.000
λ <sub>52</sub>	0.5	-0.010	0.048	0.087	1.000	-0.008	0.035	0.068	1.000
$\lambda_{\epsilon_2}$	0.5	-0.006	0.056	0.087	1.000	-0.011	0.039	0.068	1.000
λ <sub>72</sub>	0.7	-0.072	0.092	0.083	1.000	-0.055	0.069	0.065	1.000
$\lambda_{e2}$	0.7	-0.002	0.063	0.090	1.000	0.005	0.041	0.071	1.000
λ	0.7	-0.003	0.064	0.090	1.000	-0.003	0.042	0.071	1.000
$\lambda_{10,2}$	0.7	0.010	0.063	0.090	1.000	0.005	0.042	0.071	1.000
φ12	0.3	-0.029	0.084	0.101	0.830	-0.027	0.064	0.075	0.980
+12 Wai	0.3	-0.136	0.144	0.117	0.020	-0.126	0.131	0.110	0.200
+21 Wz4	0.3	0.007	0.049	0.069	1.000	0.007	0.037	0.052	1.000
+ /4 W55	0.35	0.085	0.096	0.093	1.000	0.074	0.080	0.081	1.000
1000	0.35	0.079	0.089	0.092	1.000	0.070	0.077	0.079	1.000
Ψ00 1/	0.51	0.011	0.097	0.109	1.000	0.003	0.078	0.094	1.000
$\psi_0^{\tau_D}$	0	0.022	0.041	0.060	0.003	0.019	0.035	0.050	0.003

*Note.*  $\psi_{ii}$  averaged across elements from j = 1 to 4 and 7 to 10;  $\psi_0$  averaged across all zero residual covariance estimates.

between factor). However, more work might be needed to systematically investigate this issue.

To study how the value of the correlation of the residuals of variables that load to the same factor influence the estimation of the true covariance structure using the Bayesian Lasso, additional simulation studies were conducted. The settings were similar to those of Study 1 except that the data were generated by  $\Psi$  with nonzero within-factor residual covariance  $\psi_{12}$  and all other off-diagonal elements equal to 0. Two levels of residual correlations (0.3 and 0.7) were considered, which led to  $\psi_{12} =$  $\psi_{21} = 0.152$  and  $\psi_{12} = \psi_{21} = 0.357$ , respectively. For ease of comparison, we only focused on the estimates of the elements in the residual covariance matrix. The results presented in Tables D17 and D18 indicate that the power of detecting the nonzero  $\psi_{12}$  was low whether the residual correlation was low (0.3) or high (0.7) and could be improved with a larger sample size. However, the accuracy and stability of estimates, in terms of BIAS and SE, became poorer when the correlation increased.

To study how the value of the correlation of the residuals of variables that load to different factors influences the estimation of the true covariance structure using the Bayesian Lasso, additional simulation studies were conducted. The settings were similar to those of Study 1 except that the data were generated by  $\Psi$  with nonzero between-factor residual covariance  $\psi_{74}$  and with all other off-diagonal elements equal to zero. Two levels of residual correlations (0.3 and 0.7) were considered, which led to  $\psi_{74} = \psi_{47} = 0.152$  and  $\psi_{74} = \psi_{47} = 0.357$ , respectively. The results presented in Tables D19 and D20 indicate that the power of detecting the nonzero  $\psi_{74}$  was low when the residual correlation was low (0.3), but became satisfactory (>0.800) when the residual correlation was high (0.7). Moreover, the power improved with a larger sample size. Similar to the above, the accuracy and stability of estimates, in terms of BIAS and SE, became poorer when the correlation increased. However, more work might be needed to systematically investigate this issue.

Table D17 C-Step With the Residuals of Variables That Load to the Same Factor (Residual Correlation = .3)

Par	True		N = 1	250	N = 500				
		BIAS	RMSE	SE	SIG%	BIAS	RMSE	SE	SIG%
ψ <sub>21</sub>	0.153	-0.023	0.038	0.105	0.000	-0.029	0.041	0.094	0.010
ψ55	0.35	0.055	0.066	0.088	1.000	0.052	0.060	0.071	1.000
ψ66	0.35	0.058	0.068	0.087	1.000	0.044	0.051	0.069	1.000
$\psi_{ii}$	0.51	0.042	0.072	0.114	1.000	0.030	0.056	0.096	1.000
ψ <sub>0</sub>	0	0.015	0.035	0.067	0.000	0.012	0.029	0.056	0.000

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C-Step With the Residuals of Variables That Load to the Same Factor (Residual Correlation = $.7$ )										
			N = 1	N = 250			N = 500			
Par	True	BIAS	RMSE	SE	SIG%	BIAS	RMSE	SE		
$\psi_{21}$	0.357	-0.155	0.162	0.165	0.010	-0.133	0.141	0.161		
$\psi_{55}$	0.35	0.102	0.112	0.120	1.000	0.083	0.091	0.100		

0.114

0.138

0.082

Table D18										
C-Step With the	Residuals of	Variables T	hat Load	to the	Same	Factor	(Residual	Correlation	= .	7)

0.103

0.109

0.045

Note.  $\psi_{jj}$  averaged across elements from j = 1 to 4 and 7 to 10;  $\lambda_0$  averaged across all zero loading estimates;  $\psi_0$  averaged across all zero residual covariance estimates.

1.000

1.000

0.000

0.076

0.020

0.021

0.082

0.093

0.039

0.097

0.124

0.072

#### Table D19

 $\psi_{66}$ 

 $\psi_{jj}$ 

 $\psi_0$ 

0.35

0.51

0

0.092

0.025

0.026

C-Step With the Residuals of Variables That Load to Different Factors (Residual Correlation = .3)

Par	True	N = 250					N = 500				
		BIAS	RMSE	SE	SIG%	BIAS	RMSE	SE	SIG%		
$\psi_{74}$	0.153	-0.070	0.077	0.070	0.020	-0.060	0.066	0.061	0.110		
$\psi_{55}$	0.35	0.053	0.064	0.086	1.000	0.048	0.055	0.070	1.000		
Ψ66	0.35	0.056	0.069	0.087	1.000	0.043	0.051	0.069	1.000		
$\psi_{ii}$	0.51	0.049	0.070	0.110	1.000	0.036	0.051	0.093	1.000		
ψ <sub>0</sub>	0	0.015	0.037	0.067	0.000	0.012	0.030	0.057	0.000		

#### Table D20

C-Step With the Residuals of Variables That Load to Different Factors (Residual Correlation = .7)

			N = 2	250	N = 500				
Par	True	BIAS	RMSE	SE	SIG%	BIAS	RMSE	SE	SIG%
$\psi_{74}$	0.357	-0.102	0.111	0.103	0.890	-0.088	0.092	0.091	0.980
ψ <sub>55</sub>	0.35	0.078	0.089	0.100	1.000	0.062	0.069	0.078	1.000
Ψ66	0.35	0.081	0.090	0.099	1.000	0.061	0.067	0.077	1.000
ψ,,	0.51	0.041	0.066	0.120	1.000	0.026	0.046	0.100	1.000
ψ	0	0.019	0.040	0.076	0.000	0.014	0.031	0.062	0.000

*Note.*  $\psi_{jj}$  averaged across elements from j = 1 to 4 and 7 to 10;  $\lambda_0$  averaged across all zero loading estimates;  $\psi_0$  averaged across all zero residual covariance estimates.

(Appendices continue)

SIG%

0.070

1.000

1.000

1.000

0.000

### Appendix E

## Humor Styles Questionnaire

Item	Content	Factor
1	I usually don't laugh or joke around much with other people	F1
2	If I am feeling depressed, I can usually cheer myself up with humor	F2
3	If someone makes a mistake, I will often tease them about it	F3
4	I let people laugh at me or make fun at my expense more than I should	F4
5	I don't have to work very hard at making other people laugh—I seem to be a naturally humorous person	F1
6	Even when I'm by myself, I'm often amused by the absurdities of life	F2
7	People are never offended or hurt by my sense of humor	F3
8	I will often get carried away in putting myself down if it makes my family or friends laugh	F4
9	I rarely make other people laugh by telling funny stories about myself	F1
10	If I am feeling upset or unhappy I usually try to think of something funny about the situation to make myself feel better	F2
11	When telling jokes or saying funny things, I am usually not very concerned about how other people are taking it	F3
12	I often try to make people like or accept me more by saying something funny about my own weaknesses, blunders, or faults	F4
13	I laugh and joke a lot with my closest friends	F1
14	My humorous outlook on life keeps me from getting overly upset or depressed about things	F2
15	I do not like it when people use humor as a way of criticizing or putting someone down	F3
16	I don't often say funny things to put myself down	F4
17	I usually don't like to tell jokes or amuse people	F1
18	If I'm by myself and I'm feeling unhappy, I make an effort to think of something funny to cheer myself up	F2
19	Sometimes I think of something that is so funny that I can't stop myself from saying it, even if it is not appropriate for the situation	F3
20	I often go overboard in putting myself down when I am making jokes or trying to be funny	F4
21	I enjoy making people laugh	F1
22	If I am feeling sad or upset. I usually lose my sense of humor	F2
23	I never participate in laughing at others even if all my friends are doing it	F3
24	When I am with friends or family, I often seem to be the one that other people make fun of or joke about	F4
25	I don't often joke around with my friends	F1
26	It is my experience that thinking about some amusing aspect of a situation is often a very effective way of coping with problems	F2
27	If I don't like someone. I often use humor or teasing to put them down	F3
28	If I am having problems or feeling unhappy, I often cover it up by joking around, so that even my closest friends don't know how I really feel	F4
29	I usually can't think of witty things to say when I'm with other people	F1
30	I don't need to be with other people to feel amused—I can usually find things to laugh about even when I'm by myself	F2
31	Even if something is really funny to me, I will not laugh or joke about it if someone will be offended	F3
32	Letting others laugh at me is my way of keeping my friends and family in good spirits	F4

Note. F1 = affiliative humor; F2 = self-enhancing humor; F3 = aggressive humor; F3 = self-defeating humor.

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