Bayesian linear mixed models using Stan: A tutorial for psychologists, linguists, and cognitive scientists

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Abstract

With the arrival of the R packages nlme and lme4, linear mixed models (LMMs) have come to be widely used in experimentally-driven areas like psychology, linguistics, and cognitive science. This tutorial provides a practical introduction to fitting LMMs in a Bayesian framework using the probabilistic programming language Stan. We choose Stan (rather than WinBUGS or JAGS) because it provides an elegant and scalable framework for fitting models in most of the standard applications of LMMs. We ease the reader into fitting increasingly complex LMMs, first using a two-condition repeated measures self-paced reading study, followed by a more complex 2×2 repeated measures factorial design that can be generalized to much more complex designs.

1 Introduction

Linear mixed models, or hierarchical/multilevel linear models, have become the main workhorse of experimental research in psychology, linguistics, and cognitive science, where repeated measures designs are the norm. Within the programming environment R [24], the nlme package [22] and its successor, lme4 [3] have revolutionized the use of linear mixed models (LMMs) due to their simplicity and speed: one can fit fairly complicated models relatively quickly, often with a single line of code. A great advantage of LMMs over traditional approaches such as repeated measures ANOVA and paired t-tests is that there is no need to aggregate over subjects and items to compute two sets of F-scores (or several t-scores) separately; a single model can take all sources of variance into account simultaneously. Furthermore, comparisons between conditions can easily be implemented in a single model through appropriate contrast coding.

Other important developments related to LMMs have been unfolding in computational statistics. Specifically, probabilistic programming languages like WinBUGS [18], JAGS [23] and Stan [27], among others, have made it possible to fit Bayesian LMMs quite easily. However, one prerequisite for using these programming languages is that some background statistical knowledge is needed before one can define the model. This difficulty is well-known; for example, Spiegelhalter and colleagues [26, 4] write: "Bayesian statistics has a (largely deserved) reputation for being mathematically challenging and difficult to put into practice...".

The purpose of this paper is to facilitate a first encounter with model specification in one of these programming languages, Stan. The tutorial is aimed primarily at psychologists, linguists, and cognitive scientists who have used 1me4 to fit models to their data, but may have only a basic knowledge of the underlying LMM machinery. A diagnostic test is that they may not be able to answer some or all of these questions: what is a design matrix; what is contrast coding; what is a random effects variance-covariance matrix in a linear mixed model? Our tutorial is not intended for statisticians or psychology researchers who could, for example, write their own Markov Chain Monte Carlo samplers in R or C++ or the like; for them, the Stan manual is the optimal starting point. The present tutorial attempts to ease the beginner into their first steps towards fitting Bayesian linear mixed models. More detailed presentations about linear mixed models are available in several textbooks; references are provided at the end of this tutorial.

We have chosen Stan as the programming language of choice (over JAGS and WinBUGS) because it is possible to fit arbitrarily complex models with Stan. For example, it is possible (if time consuming) to fit a model with 14 fixed effects predictors and two crossed random effects by subject and item, each involving a 14×14 variance-covariance matrix [2]; as far as we are aware, such models cannot, as far as we know, be fit in JAGS or WinBUGS.¹

In this tutorial, we take it as a given that the reader is interested in learning how to fit Bayesian linear mixed models. We do not try to explain the advantages this approach affords beyond the classical frequentist approach; for such justification, the reader is directed to the rich literature relating to a comparison between Bayesian versus frequentist statistics (such as the provocatively titled paper by Lavine [15]; and the highly accessible textbook by Kruschke [14]). We also assume that the reader is aware of Bayes' Theorem, which for our purposes amounts to the statement that the posterior distribution is proportional to the prior times the likelihood. For the purposes of this paper, the goal of a Bayesian analysis is simply to derive the posterior distribution of each parameter of interest, given some data and prior beliefs about the distributions of the parameters.

The tutorial is structured as follows. We begin by successively building up increasingly complex LMMs using the data-set reported by Gibson and Wu[8],

¹Whether it makes sense in general to fit such a complex model is a different issue; see [5], and [2] for recent discussion.

which has a simple two-condition design. At each step, we explain the structure of the model. The next section takes up inference for this two-condition design. Then we demonstrate how one can fit a somewhat more complex 2×2 factorial design.

This paper was written using a literate programming tool, knitr [29]; this integrates documentation for the accompanying code with the paper. The knitr file that generated this paper, as well as all the code and data used in this tutorial, can be downloaded from our website:

http://www.ling.uni-potsdam.de/~vasishth/statistics/BayesLMMs.html

We start with the two-condition repeated measures data-set [8] as a concrete running example. This simple example serves as a starter kit for fitting commonly used LMMs in the Bayesian setting. We assume that the reader has the relevant software installed; specifically, rstan in R. For detailed instructions, see

https://github.com/stan-dev/rstan/wiki/RStan-Getting-Started

2 Example 1: A two-condition repeated measures design

This section motivates the LMM with the self-paced reading data-set of Gibson and Wu [8]. We introduce the data-set, state our modeling goals here, and proceed to build up increasingly complex LMMs.

The scientific question Subject and object relative clauses have been widely used in reading studies to investigate sentence comprehension processes. A subject relative is a sentence like *The senator who interrogated the journalist resigned* where a noun (senator) is modified by a relative clause (who interrogated the journalist), and the modified noun is the grammatical subject of the relative clause. In an object relative, the noun modified by the relative clause is the grammatical object of the relative clause (e.g., *The senator who the journalist interrogated resigned*). In both cases, the noun that is modified (senator) is called the head noun.

A typical finding for English is that subject relatives are easier to process than object relatives [12]. Natural languages generally have relative clauses, and the subject relative advantage has until recently been considered to be true cross-linguistically. However, Chinese relative clauses apparently represent an interesting counter-example to this generalization; recent work by Hsiao and Gibson [10] has suggested that in Chinese, *object* relatives are easier to process than subject relatives at a particular point in the sentence (the head noun of the relative clause). We now present an analysis of a subsequently published data-set [8] that evaluates this claim.

The data The dependent variable of the experiment of Gibson and Wu [8] was the reading time rt of the head noun of the relative clause. This was recorded

in two conditions (subject relative and object relative), with 37 subjects and 15 items, presented in a standard Latin square design. There were originally 16 items, but one item was removed, resulting in $37 \times 15 = 555$ data points. However, eight data points from one subject (id 27) were missing. As a consequence, we have a total of 555 - 8 = 547 data points. The first few lines from the data frame are shown in Table 1; "o" refers to object relative and "s" to subject relative.

	1- :	:4		
row	subj	item	SO	rt
1	1	13	О	1561
2	1	6	\mathbf{S}	959
3	1	5	O	582
4	1	9	O	294
5	1	14	\mathbf{S}	438
6	1	4	\mathbf{S}	286
:	:	:	:	
	•	•	•	
547	9	11	O	350

Table 1: First six rows, and the last row, of the data-set of Gibson and Wu (2013), as they appear in the data frame.

We build up the Bayesian LMM from a fixed effects model to a varying intercepts model and finally to a varying intercepts, varying slopes model (the "maximal model" of Barr and colleagues [1]). The result is a probability model that expresses how the dependent variable, the reading time labeled rt, was generated in the experiment of Gibson and Wu [8].

As mentioned above, the goal of Bayesian modeling is to derive the posterior probability distribution of the model parameters from a prior probability distribution and a likelihood function. Stan makes it easy to compute this posterior distribution of each parameter of interest. The posterior distribution reflects what we should believe, given the data, regarding the value of that parameter.

2.1 Fixed Effects Model (Simple Linear Model)

We begin by making the working assumption that the dependent variable of reading time rt on the head noun is approximately log-normally distributed [25]. This assumes that the logarithm of rt is approximately normally distributed. The logarithm of the reading times, log rt, has some unknown grand mean β_0 . The mean of the log-normal distribution of rt is the sum of β_0 and an adjustment $\beta_1 \times so$ whose magnitude depends on the categorical predictor so, which has the value -1 when rt is from the subject relative condition, and +1 when rt is from the object relative condition. One way to write the model in terms of the logarithm of the reading times is as follows:

$$\log \mathsf{rt}_i = \beta_0 + \beta_1 \mathsf{so}_i + \varepsilon_i \tag{1}$$

The index i represents the i-th row in the data-frame (in this case, $i = 1, \ldots, 547$); the term ε_i represents the error in the i-th row. With the above ± 1 contrast coding, β_0 represents the grand mean of $\log rt$, regardless of relative clause type. It can be estimated by simply taking the grand mean of $\log rt$. The parameter β_1 is an adjustment to β_0 so that the mean of $\log rt$ is $\beta_0 + 1\beta_1$ when $\log rt$ is from the object relative condition, and $\beta_0 - 1\beta_1$ when $\log rt$ is from the subject relative condition. Notice that $2 \times \beta_1$ will be the difference in the means between the object and subject relative clause conditions. Together, β_0 and β_1 make up the part of the model which characterizes the effect of the experimental manipulation, relative clause type (so), on the dependent variable rt. We call this a fixed effects model because we estimate the β parameters, which are unvarying from subject to subject and from item to item. In R, this would correspond to fitting a simple linear model using the 1m function, with so as predictor and $\log rt$ as dependent variable.

The error ε is positive when $\log \mathtt{rt}_i$ is greater than the expected value $\mu_i = \beta_0 + \beta_1 \mathtt{so}_i$ and negative when $\log \mathtt{rt}_i$ is less than the expected value μ_i . Thus, the error is the amount by which the expected value differs from actually observed value. It is standardly assumed that the ε_i are independently and identically distributed as a normal distribution with mean zero and unknown standard deviation σ_e . Stan parameterizes the normal distribution by the mean and standard deviation, and we follow that convention here, by writing the distribution of ε as $N(0, \sigma_e)$ (the standard notation in statistics is in terms of mean and variance). A consequence of the assumption that the errors are identically distributed is that the distribution of ε should, at least approximately, have the same shape as the normal distribution. Independence implies that there should be no correlation between the errors—this is not the case in the data, since we have multiple measurements from each subject, and from each item.

Setting up the data We now fit the fixed effects Model 1. For the following discussion, please refer to the code in Listings 1 and 2 in the appendix. First, we read the Gibson and Wu [8] data into a data frame rDat in R, and then subset the critical region (Listing 1, lines 2 and 4). Next, we create a data list stanDat for Stan, which contains the data (Listing 1, line 7). Stan requires the data to be of type list; this is different from the 1m and 1mer functions, which assume that the data are in of type data-frame.

Defining the model The next step is to write the Stan model in a text file with extension .stan. A Stan model consists of several blocks. A block is a set of statements surrounded by brackets and preceded by the block name. We open up a file fixEf.stan in a text editor and write down the first block, the data block, which contains the declaration of the variables in the data object stanDat (Listing 2, lines 1-5). The strings real and int specify the data type for each variable. A real variable is a real number, and an int variable is an integer. For instance, N is the integer number of data points. The variables so and rt are arrays of length N whose entries are real. We constrain a variable

to take only a subset of the values allowed by its type (e.g. int or real) by specifying in brackets lower and upper bounds (e.g. <lower=-1,upper=1>).

Next, we turn to the parameters block, where the parameters are defined (Listing 2, lines 6-9). The fixed effects Model 1 has three parameters: the fixed intercept β_0 , the fixed slope β_1 , and the standard deviation σ_e of the error. The fixed effects β_0 and β_1 are in the vector beta of length two; note that although we called our parameters β_0 and β_1 in Model 1, in Stan, these are contained in a vector with indices 1 and 2, so β_0 is in beta[1] and β_1 in beta[2]. The third parameter, the standard deviation σ_e of the error (sigma_e), is also defined here, and is constrained to have lower bound 0 (Listing 2, line 8). Finally, the model block specifies the prior distribution and the likelihood (Listing 2, lines 10-15).

To understand the Stan syntax, compare the Stan code above to the specification of Model 1. The Stan code literally writes out this model. The block begins with a local variable declaration for mu, which is the mean of rt conditional on whether so is -1 for the subject relative condition or +1 for the object relative condition.

The prior distributions on the parameters beta and sigma_e would ordinarily be declared in the model block. If we don't declare any prior, it is assumed that they have a uniform prior distribution. Note that the distribution of sigma_e is truncated at zero because sigma_e is constrained to be positive (see the declaration real<lower=0> sigma_e; in the parameters block). So, this means that the error has a uniform prior with lower bound 0.

In the model block, the for-loop assigns to mu the mean for the log-normal distribution of rt[i], conditional on the value of the predictor so[i] for relative clause type. The statement rt[i] ~ lognormal(mu,sigma_e) means that the logarithm of rt is normally distributed with mean mu and standard deviation sigma_e. One could have equally well log-transformed the reading time and assumed a normal distribution instead of the lognormal.

Running the model We save the file fixEf.stan which we just wrote and fit the model in R with the function stan from the package rstan (Listing 1, lines 9 and 10). This call to the function stan will compile a C++ program which produces samples from the joint posterior distribution of the fixed intercept β_0 , the fixed slope β_1 , and the standard deviation σ_e of the error. Here, the function generates four *chains* of samples, each of which contains 2000 samples of each parameter. Samples 1 to 1000 are part of the *warmup*, where the chains settle into the posterior distribution. We analyze samples 1001 to 2000. The result is saved to an object fixEfFit of class StanFit.

Evaluating model convergence and summarizing results The first step after running the above function should be to look at the *trace plot* of each chain after warmup, using the command shown in Listing 1, lines 13 and 14. A trace plot has the chains plotted against the sample ID. In Figure 1, we see four different chains plotted against sample number going from 1001 to 2000. If the trace plot looks like a "fat, hairy caterpillar" [17] which does not bend, this

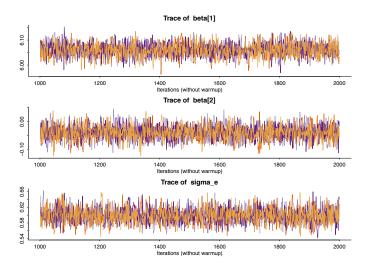


Figure 1: Trace plots of the fixed intercept β_0 , the fixed slope β_1 , and the standard deviation σ_e of the error for the fixed effects Model 1.

suggests that the chains have converged to the posterior distribution.

The second diagnostic which we use to assess whether the chains have converged to the posterior distribution is the statistic Rhat. Each parameter has the Rhat statistic associated with it [7]; this is essentially the ratio of between-chain variance to within-chain variance (analogous to ANOVA). The Rhat statistic should be approximately 1 ± 0.1 if the chain has converged. This is shown in the rightmost column of the model summary, see Table 2.

Having satisfied ourselves that the chains have converged, we turn to examine this posterior distribution. (If there is an indication that convergence has not happened, then, assuming that the model has no errors in it, increasing the number of samples usually resolves the issue.)

parameter	mean	2.5%	97.5%	Ŕ
\hat{eta}_0	6.06	0.03	6.11	1
\hat{eta}_1	-0.04	-0.09	0.01	1
$\hat{\sigma}_e$	0.60	0.56	0.64	1

Table 2: Examining the credible intervals and the R-hat statistic in the Gibson and Wu data.

The result of fitting the fixed effects Model 1 is the *joint posterior probability* distribution of the parameters β_0 , β_1 , and σ_e . The distribution is joint because each of the (4 chains × 1000 post-warmup iterations =)4000 posterior samples which the call to stan generates is a vector $\theta = (\beta_0, \beta_1, \sigma_e)^{\mathsf{T}}$ of three model parameters. Thus, the object fixEffit contains 4000 parameter vectors θ which

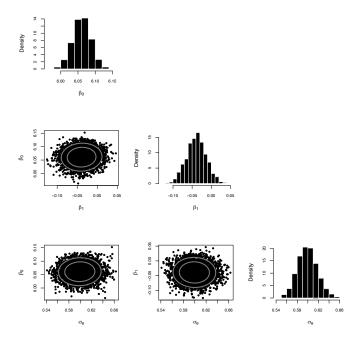


Figure 2: Bivariate joint posterior probability distribution of each element of θ with each other element (lower diagonal) and marginal posterior probability distribution of each element of θ separately (diagonal). All parameters are on the log scale, but note the difference in length scale between β_1 on the one hand and β_0 and σ_e on the other.

occupy a three dimensional space. Already in three dimensions, the posterior distribution becomes difficult to view in one graph. Figure 2 displays the joint posterior probability distribution of the elements of θ by projecting it down onto planes. In each of the three planes (lower triangular scattergrams) we see how one parameter varies with respect to the other. In the diagonal histograms, we visualize the marginal probability distribution of each parameter separately from the other parameters.

Of immediate interest is the marginal distribution of the slope β_1 . Figure 2 suggests that the posterior probability density of β_1 is mainly spread over the interval $(-\infty,0)$. One quantitative way to assess the posterior probability distribution is to examine its quantiles; see Table 2. Here, it is useful to define the concept of the *credible interval*. The $(1-\alpha)\%$ credible interval contains $(1-\alpha)\%$ of the posterior probability density. Unlike the $(1-\alpha)\%$ confidence interval from the frequentist setting, the $(1-\alpha)\%$ credible interval represents the range within which we are $(1-\alpha)\%$ certain that the true value of the parameter lies, given the prior and the data (see [21] for further discussion on CIs

vs credible intervals). A common convention is to use the interval ranging from the 2.5 to 97.5 percentiles. We follow this convention and 95% credible intervals in Table 2.

The samples of β_1 suggests that approximately 94% of the posterior probability density is below zero, suggesting that there is some evidence that object relatives are easier to process than subject relatives in Chinese, given the Gibson and Wu data. However, since the 95% credible interval includes 0, we may be reluctant to conclude that object relatives are easier to process. We will say more about the evaluation of research hypotheses further on.

2.2 Varying Intercepts Mixed Effects Model

The fixed effects Model 1 is inappropriate for the Gibson and Wu data because it does not take into account the fact that we have multiple measurements for each subject and item. As mentioned above, these multiple measurements lead to a violation of the independence of errors assumption. Moreover, the fixed effects coefficients β_0 and β_1 represent means over all subjects and items, ignoring the fact that some subjects will be faster and some slower than average; similarly, some items will be read faster than average, and some slower.

In linear mixed models, we take this by-subject and by-item variability into account by adding adjustment terms u_{0j} and w_{0k} , which adjust β_0 for subject j and item k. This partially decomposes ε_i into a sum of the terms u_{0j} and w_{0k} , which are adjustments to the intercept β_0 for the subject j and item k associated with \mathbf{rt}_i . If subject j is slower than the average of all the subjects, u_j would be some positive number, and if item k is read faster than the average reading time of all the items, then w_k would be some negative number. Each subject j has their own adjustment u_{0j} , and each item its own w_{0k} . These adjustments u_{0j} and w_{0k} are called random intercepts by Pinheiro and Bates [22] and varying intercepts by Gelman and Hill [6], and by adjusting β_0 by these we account for the variability between speakers, and between items.

It is standardly assumed that these adjustments are normally distributed around zero with unknown standard deviation: $u_0 \sim N(0, \sigma_u)$ and $w_0 \sim N(0, \sigma_w)$; the subject and item adjustments are also assumed to be mutually independent. We now have three sources of variance in this model: the standard deviation of the errors σ_e , the standard deviation of the by-subject random intercepts σ_u , and the standard deviation of the by-item varying intercepts σ_w . We will refer to these as variance components.

We now express the logarithm of reading time, which was produced by subjects $j=1,\ldots,37$ reading items $k=1,\ldots,15$, in conditions i=1,2 (1 refers to subject relatives, 2 to object relatives), as the following sum. Notice that we are now using a slightly different way to describe the model, compared to the fixed effects model. We are using indices for subject, item, and condition to identify unique rows. Also, instead of writing β_1 so, we index β_1 by the condition i. This follows the notation used in the textbook on linear mixed models, written by the authors of nlme [22], the precursor to lme4.

$$\log \mathsf{rt}_{ijk} = \beta_0 + \beta_{1i} + u_{0j} + w_{0k} + \varepsilon_{ijk} \tag{2}$$

Model 2 is an LMM, and more specifically a varying intercepts model. The coefficient β_{1i} is the one of primary interest; it will have some mean value $-\beta_1$ for subject relatives and $+\beta_1$ for object relatives due to the contrast coding. So, if our posterior mean for β_1 is negative, this would suggest that object relatives are read faster than subject relatives.

We fit the varying intercepts Model 2 in Stan in much the same way as the fixed effects Model 1. For the following discussion, please consult Listing 3 for the R code used to run the model, and Listing 4 for the Stan code.

Setting up the data The data which we prepare for passing on to the function stan now includes subject and item information (Listing 3, lines 2-8). The data block in the Stan code accordingly includes the number J, K of subjects and items, respectively; and the variable N records the number of rows in the data frame.

Defining the model Model 2, shown in Listing 4, still has the fixed intercept β_0 , the fixed slope β_1 , and the standard deviation σ_e of the error, and we specify these in the same way as we did for the fixed effects Model 1. In addition, the varying intercepts Model 2 has by-subject varying intercepts u_{0j} for $j = 1, \ldots, J$ and by-item varying intercepts w_{0k} for $k = 1, \ldots, K$. The standard deviation of u_0 is σ_u and the standard deviation of w_0 is σ_w . We again constrain the standard deviations to be positive.

The model block places normal distribution priors on the varying intercepts u_0 and w_0 . We implicitly place uniform priors on sigma_u, sigma_w, and sigma_e by omitting them from the model block. As pointed out earlier for sigma_e, these prior distributions have lower bound zero because of the constraint <lower=0> in the variable declarations.

The statement about how each row in the data is generated is shown in Listing 4, lines 26-29; here, both the fixed effects and the varying intercepts for subjects and items determine the expected value mu. The vector u has varying intercepts for subjects. Likewise, the vector w has varying intercepts for items. The for-loop in lines 26-29 now adds u[subj[i]] + w[item[i]] to the mean beta[1] of the distribution of rt[i]. These are subject- and item-specific adjustments to the fixed-effects intercept beta[1]. The term u[subj[i]] identifies the id of the subject for row i in the data-frame; thus, if i = 1, then subj[1]=1, and item[1]=13 (see Table 1).

Running the model We pass the list stanDat of data to stan, which compiles a C++ program to sample from the posterior distribution of Model 2. Stan samples from the posterior distribution of the model parameters, including the varying intercepts u_{0j} and w_{0k} for each subject j = 1, ..., J and item k = 1, ..., K.

It may be helpful to rewrite the model in mathematical form following the Stan syntax ([6] use a similar notation); the Stan statements are slightly different from the way that we expressed Model 2. Defining i as the row id in the data, i.e., $i = 1, \ldots, 547$, we can write:

Likelihood:
$$\mu_i = \beta_0 + u_{[subj[i]]} + w_{[item[i]]} + \beta_1 \times \mathfrak{so}_i$$

$$\mathsf{rt}_i \sim \mathsf{LogNormal}(\mu_i, \sigma_e)$$
 Priors:
$$u \sim \mathsf{Normal}(0, \sigma_u) \quad w \sim \mathsf{Normal}(0, \sigma_w)$$

$$\sigma_e, \sigma_u, \sigma_w \sim \mathsf{Uniform}(0, \infty)$$

$$\beta \sim \mathsf{Uniform}(-\infty, \infty)$$
 (3)

Here, notice that the *i*-th row in the statement for μ identifies the subject id ranging from $j = 1, \ldots, 37$, and the item id ranging from $k = 1, \ldots, 15$.

Summarizing the results The posterior distributions of each of the parameters is summarized in Table 3. The \hat{R} values suggest that model has converged. Note also that compared to Model 1, the estimate of σ_e is smaller; this is because the other two variance components are now being estimated as well. Note that the 95% credible interval for the estimate $\hat{\beta}_1$ includes 0; thus, there is some evidence that object relatives are easier than subject relatives, but we cannot exclude the possibility that there is no difference in the reading times between the two relative clause types.

parameter	mean	2.5%	97.5%	\hat{R}
\hat{eta}_0	6.06	5.92	6.20	1
$\hat{\beta}_1$	-0.04	-0.08	0.01	1
$\hat{\sigma}_e$	0.52	0.49	0.55	1
$\hat{\sigma}_u$	0.25	0.19	0.34	1
$\hat{\sigma}_w$	0.20	0.12	0.32	1

Table 3: The quantiles and the \hat{R} statistic in the Gibson and Wu data, the varying intercepts model.

2.3 Varying Intercepts, Varying Slopes Mixed Effects Model

Consider now that subjects who are faster than average (i.e., who have a negative varying intercept) may exhibit greater slowdowns when they read subject relatives compared to object relatives. Similarly, it is in principle possible that items which are read faster (i.e., which have a large negative varying intercept) may show a greater slowdown in subject relatives than object relatives. The

opposite situation could also hold: faster subjects may show smaller SR-OR effects, or items read faster may show smaller SR-OR effects. Although such individual-level variability was not of interest in the original paper by Gibson and Wu, it could be of theoretical interest (see, for example, [13]). Furthermore, as Barr and colleagues [1] point out, it is in principle desirable to include a fixed effect factor in the random effects as a varying slope if the experiment design is such that subjects see both levels of the factor (cf. [2]).

In order to express this structure in the LMM, we must make two changes in the varying intercepts Model 2.

Adding varying slopes The first change is to let the size of the effect for the predictor so vary by subject and by item. The goal here is to express that some subjects exhibit greater slowdowns in the object relative condition than others. We let effect size vary by subject and by item by including in the model by-subject and by-item varying slopes which adjust the fixed slope β_1 in the same way that the by-subject and by-item varying intercepts adjust the fixed intercept β_0 . This adjustment of the slope by subject and by item is expressed by adjusting β_1 by adding two terms u_{1j} and w_{1k} . These are random or varying slopes, and by adding them we account for how the effect of relative clause type varies by subject j and by item k. We now express the logarithm of reading time, which was produced by subject j reading item k, as the following sum. The subscript i indexes the conditions.

$$\log \mathtt{rt}_{ijk} = \underbrace{\beta_0 + u_{0j} + w_{0k}}_{\text{varying intercepts}} + \underbrace{\beta_1 + u_{1ij} + w_{1ik}}_{\text{varying slopes}} + \varepsilon_{ijk} \tag{4}$$

Defining a variance-covariance matrix for the random effects The second change which we make to Model 2 is to define a covariance relationship between by-subject varying intercepts and slopes, and between by-items intercepts and slopes. This amounts to adding an assumption that the by-subject slopes u_1 could in principle have some correlation with the by-subject intercepts u_0 ; and by-item slopes w_1 with by-item intercept w_0 . We explain this in detail below.

Let us assume that the adjustments u_0 and u_1 are normally distributed with mean zero and some variances σ_{u0}^2 and σ_{u1}^2 , respectively; also assume that u_0 and u_1 have correlation ρ_u . It is standard to express this situation by defining a variance-covariance matrix Σ_u (sometime this is simply called a variance matrix). This matrix has the variances of u_0 and u_1 respectively along the diagonals, and the covariances on the off-diagonals. (The covariance between two variables X, Y, Cov(X,Y) is defined as the product of their correlation ρ and their standard deviations σ_X and σ_Y : $Cov(X,Y) = \rho \sigma_X \sigma_Y$.)

$$\Sigma_u = \begin{pmatrix} \sigma_{u0}^2 & \rho_u \sigma_{u0} \sigma_{u1} \\ \rho_u \sigma_{u0} \sigma_{u1} & \sigma_{u1}^2 \end{pmatrix}$$
 (5)

Similarly, we can define a variance-covariance matrix Σ_w for items, using the standard deviations σ_{w0} , σ_{w1} , and the correlation ρ_w .

$$\Sigma_w = \begin{pmatrix} \sigma_{w0}^2 & \rho_w \sigma_{w0} \sigma_{w1} \\ \rho_w \sigma_{w0} \sigma_{w1} & \sigma_{w1}^2 \end{pmatrix} \tag{6}$$

The standard way to express this relationship between the subject intercepts u_0 and slopes u_1 , and the item intercepts w_0 and slopes w_1 , is to define a bivariate normal distribution as follows:

$$\begin{pmatrix} u_0 \\ u_1 \end{pmatrix} \sim N \begin{pmatrix} \begin{pmatrix} 0 \\ 0 \end{pmatrix}, \Sigma_u \end{pmatrix} \quad \begin{pmatrix} w_0 \\ w_1 \end{pmatrix} \sim N \begin{pmatrix} \begin{pmatrix} 0 \\ 0 \end{pmatrix}, \Sigma_w \end{pmatrix}$$
 (7)

An important point to notice here is that any $n \times n$ variance-covariance matrix has associated with it an $n \times n$ correlation matrix. In the subject variance-covariance matrix Σ_u , the correlation matrix is

$$\begin{pmatrix} 1 & \rho_{01} \\ \rho_{01} & 1 \end{pmatrix} \tag{8}$$

In a correlation matrix, the diagonal elements will always be 1, because a variable always has a correlation of 1 with itself. The off-diagonals will have the correlations between the variables. Note also that, given the variances σ_{u0}^2 and σ_{u1}^2 , we can always recover the variance-covariance matrix, if we know the correlation matrix. This is because of the above-mentioned definition of covariance.

A correlation matrix can be decomposed into a square root of the matrix, using the Cholesky decomposition. Thus, given a correlation matrix C, we can obtain its square root L; an obvious consequence is that we can square L to get the correlation matrix C back. This is easy to illustrate with a simple example. Suppose we have a correlation matrix:

$$C = \begin{pmatrix} 1 & -0.5 \\ -0.5 & 1 \end{pmatrix} \tag{9}$$

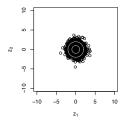
We can use the Cholesky decomposition function in R, chol, to derive the lower triangular square root L of this matrix. This gives us:

$$L = \begin{pmatrix} 1 & 0 \\ -0.5 & 0.8660254 \end{pmatrix} \tag{10}$$

We can confirm that this is a square root by multiplying L with itself to get the correlation matrix back (squaring a matrix is done by multiplying the matrix by its transpose):

$$LL^{\mathsf{T}} = \begin{pmatrix} 1 & 0 \\ -0.5 & 0.8660254 \end{pmatrix} \begin{pmatrix} 1 & -0.5 \\ 0 & 0.8660254 \end{pmatrix} = \begin{pmatrix} 1 & -0.5 \\ -0.5 & 1 \end{pmatrix} \tag{11}$$

The reason that we bring up the Cholesky decomposition here is that we will use it to generate the by-subject and by-item adjustments to the intercept and slope fixed-effects parameters.



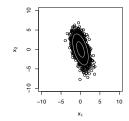


Figure 3: Uncorrelated random variables $\mathbf{z} = (z_1, z_2)^{\mathsf{T}}$ (left) and correlated random variables $\mathbf{x} = (x_1, x_2)^{\mathsf{T}}$ (right).

Generating correlated random variables using the Cholesky decomposition The by-subject and by-item adjustments are generated using the following standard procedure for generating correlated random variables $\mathbf{x} = (x_1, x_2)$:

1. Given a vector of standard deviances (e.g., σ_{u0}, σ_{u1}), create a diagonal matrix:

$$\tau = \begin{pmatrix} \sigma_{u0} & 0\\ 0 & \sigma_{u0} \end{pmatrix} \tag{12}$$

- 2. Premultiply the diagonalized matrix τ with the Cholesky decomposition L of the correlation matrix C to get a matrix Λ .
- 3. Generate values from a random variable $\mathbf{z} = (z_1, z_2)^{\intercal}$, where z_1 and z_2 each have independent N(0, 1) distributions (left panel of Figure 3).
- 4. Multiply Λ with \mathbf{z} ; this generates the correlated random variables \mathbf{x} (right panel of Figure 3).

This digression into Cholesky decomposition and the generation of correlated random variables is important to understand for building the Stan model. We will define a vague prior distribution on L, and a vague prior on the standard deviances. This allows us to generate the by-subject and by-item adjustments to the fixed effects intercepts and slopes.

Defining the model With this background, implementing the varying intercepts, varying slope Model 4 is straightforward; see Listing 6 for the code. The data block is the same as before. The parameters block contains several new parameters. This time, we have vectors $sigma_u$ and $sigma_w$ which are $(\sigma_{u0}, \sigma_{u1})^{\mathsf{T}}$ and $(\sigma_{w0}, \sigma_{w1})^{\mathsf{T}}$, instead of scalar values as in Model 2. The variables L_u, L_w, z_u, and z_w, which have been declared in the parameters block, play a role in the transformed parameters block, a block which we did not use in the earlier models. The transformed parameters block generates the by-subject and by-item varying intercepts and slopes using the parameters $sigma_u$, $sigma_w$, L_u, L_w, z_u, and z_w. The J pairs of by-subject varying intercepts and slopes

are in the rows of the $J \times 2$ matrix \mathbf{u} , and the K pairs of by-item varying intercepts and slopes are in the rows of the $K \times 2$ matrix \mathbf{w} .

These varying intercepts and slopes are obtained through the statements diag_pre_multiply(sigma_u, L_u) * z_u and diag_pre_multiply(sigma_w, L_w) * z_w. This statement generates varying intercepts and slopes from the joint probability distribution of equation 7. The parameters L_u, L_w are the Cholesky decompositions of the subject and item correlation matrices, respectively, and z_u, and z_w are N(0,1) random variables.

It is helpful to walk through steps 1 to 4 involved in generating the varying intercepts and slopes using the procedure described above for generating correlated random variables. The statement diag_pre_multiply(sigma_u,L_u) * z_u computes the transpose matrix product (steps 1 and 2). The right multiplication of this product by z_u, a matrix of normally distributed random variables (step 3), yields the varying intercepts and slopes (step 4).

$$\begin{pmatrix}
u_{01} & u_{11} \\
u_{02} & u_{12} \\
\vdots & \vdots \\
u_{0J} & u_{1J}
\end{pmatrix} = \left(\operatorname{diag}(\sigma_{u0}, \sigma_{u1}) \mathbf{L}_{u} \mathbf{z}_{u}\right)^{\mathsf{T}}$$

$$= \left(\begin{pmatrix} \sigma_{u0} & 0 \\ 0 & \sigma_{01} \end{pmatrix} \begin{pmatrix} \ell_{11} & 0 \\ \ell_{21} & \ell_{22} \end{pmatrix} \begin{pmatrix} z_{11} & z_{12} & \dots & z_{1J} \\ z_{21} & z_{22} & \dots & z_{2J} \end{pmatrix}\right)^{\mathsf{T}}$$
(13)

Turning to the model block, here, we place priors on the parameters declared in the parameters block, and define how these parameters generate log rt (Listing 6, lines 30-42). The definition of the prior L_u \sim lkj_corr_cholesky(2.0) implicitly places a so-called lkj prior with shape parameter $\eta=2.0$ on the correlation matrices

$$\begin{pmatrix} 1 & \rho_u \\ \rho_u & 1 \end{pmatrix}$$
 and $\begin{pmatrix} 1 & \rho_w \\ \rho_w & 1 \end{pmatrix}$ (14)

where ρ_u is the correlation between the by-subject varying intercept σ_{u0} and slope σ_{u1} (cf. the covariance matrix of Equation 5) and ρ_w is the correlation between the by-item varying intercept σ_{w0} and slope σ_{w1} . The lkj distribution with shape parameter $\eta=1.0$ is a uniform prior over all 2×2 correlation matrices; it scales up to larger correlation matrices. The parameter η has an effect on the shape of the lkj distribution. Our choice of $\eta=2.0$ implies that the correlations in the off-diagonals are near zero, reflecting the prior belief that there is no correlation between intercepts and slopes.

The statement $to_vector(z_u) \sim normal(0,1)$ places a normal distribution with mean zero and standard deviation one on z_u . The same goes for z_w . The for-loop assigns to mu the mean of the log-normal distribution from which we draw rt[i], conditional on the value of the predictor so[i] for relative clause type and the subject and item identity.

We can now fit the varying intercepts, varying slopes Model 4; see Listing 7 for the code. We see in the model summary in Table 4 that the model has converged, and that the credible intervals of the parameter of interest, β_1 , still include 0. In fact, the posterior probability of the parameter being less than 0 is now 90% (this information can be extracted as shown in Listing 5, lines 6-8).

parameter	mean	2.5%	97.5%	\hat{R}
\hat{eta}_0	6.06	5.92	6.21	1
\hat{eta}_1	-0.04	-0.09	0.02	1
$\hat{\sigma}_e$	0.51	0.48	0.55	1
$\hat{\sigma}_{u0}$	0.25	0.19	0.34	1
$\hat{\sigma}_{u1}$	0.07	0.01	0.14	1
$\hat{\sigma}_{w0}$	0.20	0.13	0.32	1
$\hat{\sigma}_{w1}$	0.04	0.0	0.10	1

Table 4: The quantiles and the \hat{R} statistic in the Gibson and Wu data, the varying intercepts, varying slopes model.

Figure 4 plots the varying slope's posterior distribution against the varying intercept's posterior distribution for each subject. The correlation between u_0 and u_1 is negative, as captured by the marginal posterior distributions of the correlation ρ_u between u_0 and u_1 . Thus, Figure 4 suggests that the slower a subject's reading time is on average, the slower they read object relatives. In contrast, Figure 4 shows no clear pattern for the by-item varying intercepts and slopes. We briefly discuss inference next.

3 Inference

Having fit a varying intercepts, varying slopes Model 4, we now explain one way to carry out statistical inference, using credible intervals. We have used this approach to draw inferences from data in previously published work (e.g., [4], [9]). There are of course other approaches possible for carrying out inference. Bayes Factors are an example; see Lee and Wagenmakers [16]. Another is to define a Region of Practical Equivalence [14]. The reader can choose the approach they find the most appealing.

3.1 Inference using credible intervals

The result of fitting the varying intercepts, varying slopes Model 4 is the posterior distribution of the model parameters. As mentioned above in connection with Models 1-3, direct inference from the posterior distributions is possible. For instance, we can find the posterior probability with which the fixed intercept β_1 or the correlation ρ_u between by-subject varying intercepts and slopes take on any given value by consulting the marginal posterior distributions whose

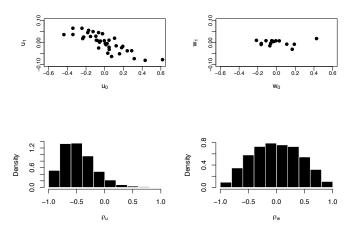


Figure 4: The top row shows the relationship between the varying slopes (y-axis) and intercepts (x-axis) for each subject (left panel) and item (right panel). The bottom row shows the posterior distribution of the parameter of correlation between the varying slopes and intercepts for each subject (left panel) and item (right panel).

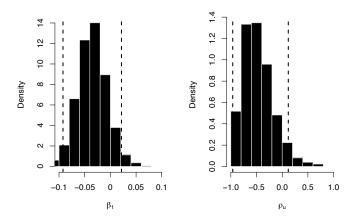


Figure 5: Upper and lower bounds on the credible intervals (dashed lines) plotted over the marginal posterior distribution of the fixed slope β_1 (left) and of the correlation ρ_u between the by-subject varying intercepts and varying slopes (right).

histograms are shown in Figure 5. The information conveyed by such graphs can be sharpened by using the 95% credible interval, mentioned earlier. Approximately 95% of the posterior density of β_1 lies between the 2.5 percentile -0.092 and the 97.5 percentile 0.023. This leads us to conclude that the slope β_1 for relative clause type so is less than zero with probability 90% (see line 8 in Listing 5). Since 0 is included in the credible interval, it is difficult to draw the inference that object relative clauses are read faster than subject relative clauses. However, one could perhaps still make a weak claim to that effect, especially if a lot of evidence has accumulated in other experiments that supports such a conclusion (see [28] for a more detailed discussion).

What about the correlations between varying intercepts and varying slopes for subject and for item? What can we infer from the analysis about these relationships? The 95% credible interval for ρ_u is (-1,0.1). Our belief that ρ_u is less than zero is rather uncertain, although we can conclude that ρ_u is less than zero with probability 90%. There is only weak evidence that subjects who read faster than average exhibit greater slowdowns at the head noun of object relative clauses than subjects who read slower than average. For the by-item varying intercepts and slopes, it is pretty clear that we do not have enough data (15 items) to draw any conclusions. For these data, it probably makes sense to fit a simpler model [2], with only varying intercepts and slopes for subject, and only varying intercepts for items; although there is no harm done here if we fit a model with a full variance-covariance matrix for both subjects and items.

In sum, regarding our main research question, our conclusion here is that we cannot say that object relatives are harder to process than subject relatives, because the credible interval for β_1 includes 0. However, one could argue that there is *some* weak evidence in favor of the hypothesis, since the posterior probability of the parameter being negative is approximately 90%.

4 Example 2: Generalizing the linear mixed model to factorial designs

The Gibson and Wu [8] data-set has a two-condition design. This section presents a varying intercepts, varying slopes model for a 2×2 factorial design. Because of the more general matrix formulation we use here, the Stan code can be deployed with minimal changes for much more complex designs, including correlational studies.

Our example is the 2×2 repeated measures factorial design of Husain et al [11] (Experiment 1), also a self-paced reading study on relative clauses. The dependent variable was the reading time rt of the relative clause verb. The factors were relative clause type, which we code with the predictor so (so = +1 for object relatives and so = -1 for subject relatives) and distance between the head noun and the relative clause verb, which we code with the predictor dist (dist = +1 for far and dist = -1 for near). Their interaction is the product of the dist and so contrast vectors, and labeled as the predictor int. The 60

subjects were speakers of Hindi, an Indo-Aryan language spoken primarily in India. The 24 items were presented in a standard, fully balanced Latin square design. This resulted in a total of 1440 data points ($60 \times 24 = 1440$). The first few lines from the data frame are shown below.

row	subj	item	so	dist	$_{ m rt}$
1	1	14	S	n	1561
2	1	16	O	\mathbf{n}	959
3	1	15	O	f	582
4	1	18	\mathbf{S}	\mathbf{n}	294
5	1	4	O	\mathbf{n}	438
6	1	17	\mathbf{S}	f	286
:	:	:	:	:	:
1440	9	13	\mathbf{s}	f	516

Table 5: The first six rows, and the last row, of the data-set of Husain et al. (2014, Experiment 1), as they appear in the data frame.

The theoretical interest is in determining whether relative clause type and distance influence reading time, and whether there is an interaction between these two factors. We use Stan to determine the posterior probability distribution of the fixed effect β_1 for relative clause type, the fixed effect β_2 for distance, and their interaction β_3 .

We fit a varying intercepts, varying slopes model to this data-set. This is an extension of Model 4. The grand mean β_0 of log rt is adjusted by subject and by item through the varying intercepts u_0 and w_0 , which are unique values for each subject and item respectively. Likewise, the three fixed effects β_1 , β_2 , and β_3 which are associated with the predictors so, dist, and int, respectively, are adjusted by the by-subject varying slopes u_1 , u_2 , and u_3 and by-item varying slopes w_1 , w_2 , and w_3 .

It is more convenient to represent this model in matrix form. We build up the model specification by first noting that, for each subject, the by-subject varying intercept u_0 and slopes u_1 , u_2 , and u_3 have a multivariate normal prior distribution with mean zero and covariance matrix Σ_u . Similarly, for each item, the by-item varying intercept w_0 and slopes w_1 , w_2 , and w_3 have a multivariate normal prior distribution with mean zero and covariance matrix Σ_w . We can write this as follows:

$$\begin{pmatrix} u_0 \\ u_1 \\ u_2 \\ u_3 \end{pmatrix} \sim \mathcal{N} \begin{pmatrix} \begin{pmatrix} 0 \\ 0 \\ 0 \\ 0 \end{pmatrix}, \Sigma_u \end{pmatrix} \quad \begin{pmatrix} w_0 \\ w_1 \\ w_2 \\ w_3 \end{pmatrix} \sim \mathcal{N} \begin{pmatrix} \begin{pmatrix} 0 \\ 0 \\ 0 \\ 0 \end{pmatrix}, \Sigma_w \end{pmatrix}$$
(15)

The error ε is assumed to have a normal distribution with mean zero and standard deviation σ_e . Thus, the varying intercepts, varying slopes here will be the same as Model 4, just with two additional predictors dist, int along with

their associated fixed-effect slopes β_2 , β_3 and random-effect slopes u_2 , u_3 , w_2 , and w_3 .

We proceed to implement the model in Stan. First we read in the data-set (see Listing 7). Instead of passing the predictors so, dist, and their interaction int to stan as vectors, as we did with so earlier, we make so, dist, and int into a design matrix X using the function model.matrix available in $R.^2$ The first column of the design matrix X consists of all ones. The second column is the predictor so which codes the factor for relative clause type. The third column the predictor dist which codes the factor for distance. The fourth column is the predictor int which codes the interaction between relative clause type and distance. The model matrix thus consists of a fully factorial 2×2 design, with blocks of this design repeated for each subject. For the full data-set, we could write it very compactly in matrix form as follows:

$$\log(\mathbf{rt}) = \mathbf{X}\beta + \mathbf{Z}_{u}\mathbf{u} + \mathbf{Z}_{w}\mathbf{w} + \varepsilon \tag{16}$$

Here, **X** is the $N \times P$ model matrix (with N = 1440, since we have 1440 data points; and P = 4 since we have the intercept plus three other fixed effects), β is a $P \times 1$ vector of fixed effects parameters, \mathbf{Z}_u and \mathbf{Z}_w are the subject and item model matrices $(N \times P)$, and u and w are the by-subject and by-item adjustments to the fixed effects estimates; these are identical to the design matrix \mathbf{X} in the model with varying intercepts and varying slopes included. For more examples of similar model specifications in Stan, see the R package RePsychLing on github (https://github.com/dmbates/RePsychLing).

Having defined the model, we proceed to assemble the list stanDat of data, relying on the above matrix formulation; please refer to Listing 7. The number N of observations, the number J of subjects and K of items, the reading times rt, and the subject and item indicator variables subj and item are familiar from the previous models presented. The integer P is the number of fixed effects (four including the intercept). Model 16 includes a varying intercept u_0 and varying slopes u_1, u_2, u_3 for each subject, and so the number $\mathbf{n}_{-}\mathbf{u}$ of by-subject random effects equals P. Likewise, Model 16 includes a varying intercept w_0 and varying slopes w_1, w_2, w_3 for each item, and so the number n_w of by-item random effects also equals P. The data block contains the corresponding variables. We declare the fixed effects design matrix X as an array of N row vectors whose components are the predictors associated with the N reading times. Likewise for the subject and item random effects design matrices Z_u and Z_w , which correspond to Z_u and \mathbf{Z}_w respectively in Model 16. The vector beta contains the fixed effects $\beta_0, \beta_1, \beta_2, \text{ and } \beta_3.$ The matrices L_u, L_w and the arrays z_u, z_w of vectors (not to be confused with the design matrices Z_u and Z_w) will generate the varying intercepts and slopes u_0, \ldots, u_3 and w_0, \ldots, w_3 , using the procedure described for Model 4. The vector sigma_u contains the standard deviations of the by-subject varying intercepts and slopes u_0, \ldots, u_3 , and the vector sigma_w contains the standard deviations of the by-item varying intercepts and slopes

²Here, we would like to acknowledge the contribution of Douglas Bates in specifying the model in this general matrix form.

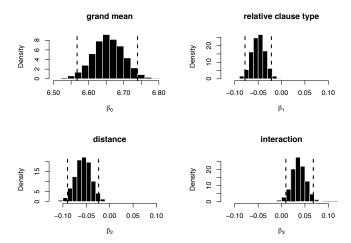


Figure 6: Marginal posterior distribution and HPD intervals of the fixed effects grand mean β_0 , slope β_1 for relative clause type, slope β_2 for distance, and interaction β_3 . All fixed effects are on the log-scale.

 w_0, \ldots, w_3 . The variable sigmale is the standard deviation σ_e of the error ε . The transformed parameters block generates the by-subject intercepts and slopes u_0, \ldots, u_3 and the by-item intercepts and slopes w_0, \ldots, w_3 .

We place lkj priors on the random effects correlation matrices through the lkj_corr_cholesky(2.0) priors on their Cholesky factors L_u and L_w. We implicitly place uniform priors on the fixed effects β_0, \ldots, β_3 , the random effects standard deviations $\sigma_{u0}, \ldots, \sigma_{u3}$, and $\sigma_{w0}, \ldots, \sigma_{w3}$ and the error standard deviation σ_e by omitting any prior specifications for them in the model block. We specify the likelihood with the probability statement that rt[i] is distributed log-normally with mean X[i] * beta + Z_u[i] * u[subj[i]] + Z_w[i] * w[item[i]] and standard deviation sigma_e. The next step towards model-fitting is to pass the list stanDat to stan, which compiles a C++ program to sample from the posterior distribution of the model parameters.

Figure 6 plots histograms of the marginal posterior distribution of the fixed effects. The HPD interval of the fixed effect $\hat{\beta}_1$ for relative clause type is entirely below zero. This is evidence that object relatives are read faster than subject relatives. The HPD interval of the fixed effect $\hat{\beta}_2$ for distance is also entirely below zero. This is evidence of a slowdown when the verb (where reading time was measured) is closer to the head noun of the relative clause. The HPD of the interaction $\hat{\beta}_3$ between relative clause type and distance is greater than zero, which is evidence for a greater slowdown on subject relatives when the distance between the verb and head noun is short.

A major advantage of the above matrix formulation is that we do not need to write a new Stan model for a future repeated measures factorial design. All we have to do now is define the design matrix X appropriately, and include it

(along with appropriately defined Z_u and Z_w for the subjects and items random effects) as part of the data specification that is passed to Stan.

5 Concluding remarks, and further reading

We hope that this tutorial has given the reader a flavor of what it would be like to fit Bayesian linear mixed models. There is of course much more to say on the topic, and we hope that the interested reader will take a look at some of the excellent books that have recently come out. We suggest below a sequence of reading that we found helpful. A good first general textbook is by Gelman and Hill [6]; it begins with the frequentist approach and only later transitions to Bayesian models. The forthcoming book by Mcelreath [20] is also excellent. For those looking for a psychology-specific introduction, the books by Kruschke [14] and Lee and Wagenmakers [16] are to be recommended, although for the latter the going might be easier if the reader has already looked at Gelman and Hill [6]. As a second book, [17] is recommended; it provides many interesting and useful examples using the BUGS language, which are discussed in exceptionally clear language. Many of these books use the BUGS syntax [18], which the probabilistic programming language JAGS [23] also adopts; however, Stan code for these books is slowly becoming available on the Stan home page (https://github.com/stan-dev/example-models/wiki). For those with introductory calculus, a slightly more technical introduction to Bayesian methods by Lynch [19] is an excellent choice. Finally, the textbook by Gelman and colleagues [5] is the definitive modern guide, and provides a more advanced treatment.

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```
## read in data:
1
    rDat <- read.table ( "gibsonwu2012data.txt" , header=TRUE )</pre>
2
    ## subset critical region:
3
    rDat <- subset( rDat , region == "headnoun" )</pre>
    ## create data as list for Stan, and fit model:
    stanDat <- list( rt = rDat$rt, so = rDat$type, N = nrow(rDat) )</pre>
    library(stan)
8
    fixEfFit <- stan ( file = "fixEf.stan" , data = stanDat ,</pre>
9
                    iter = 2000 , chains = 4 )
10
11
    ## plot traceplot, excluding warm-up:
12
    13
14
15
    ## examine quantiles of posterior distributions:
    print( fixEfFit , pars = c("beta","sigma_e") ,
17
           probs = c(0.025, 0.5, 0.975)
18
19
    ## examine quantiles of parameter of interest:
20
    beta1 <- extract ( fixEfFit , pars=c("beta[2]"))$beta</pre>
    print ( signif ( quantile ( beta1, probs = c(0.025, 0.5, 0.975))
22
             , 2))
```

Listing 1: Code for the fixed effects Model 1.

```
data {
1
       int<lower=1> N;
                                         //number of data points
2
                                         //reading time
3
       real rt[N];
       real<lower=-1,upper=1> so[N];
4
                                       //predictor
5
    parameters {
6
       vector[2] beta;
                                    //intercept and slope
7
       real<lower=0> sigma_e;
                                    //error sd
8
9
    model {
10
11
       real mu;
                                            // likelihood
       for (i in 1:N){
12
13
         mu <- beta[1] + beta[2]*so[i];</pre>
         rt[i] ~ lognormal(mu,sigma_e);
14
15
16
```

Listing 2: Stan code for the fixed effects Model 1.

```
## format data for Stan:
1
    stanDat<-list(subj=as.integer(factor(rDat$subj))),</pre>
2
3
                   item=as.integer(factor(rDat$item)),
                   rt=rDat$rt,
                   so=rDat$so,
5
                   N=nrow(rDat),
6
                   J=length(unique(rDat$subj)),
7
                   K=length(unique(rDat$item)))
8
9
    ## Sample from posterior distribution:
10
11
    ranIntFit <- stan(file="ranInt.stan", data=stanDat,</pre>
                       iter=2000, chains=4)
12
     ## Summarize results:
13
    print(ranIntFit,pars=c("beta","sigma_e","sigma_u","sigma_w"),
14
           probs=c(0.025,0.5,0.975))
15
    beta1 <- extract(ranIntFit,pars=c("beta[2]"))$beta</pre>
17
18
    print(signif(quantile(beta1,probs=c(0.025,0.5,0.975)),2))
19
     ## Posterior probability of betal being less than 0:
20
    mean(beta1<0)
```

Listing 3: Code for running Model 2, the varying intercepts model.

```
data {
 1
       int<lower=1> N;
                                          //number of data points
2
3
       real rt[N];
                                          //reading time
       real<lower=-1,upper=1> so[N];
                                          //predictor
 4
       int<lower=1> J;
                                          //number of subjects
5
       int<lower=1> K;
                                          //number of items
 6
       int<lower=1, upper=J> subj[N];
                                          //subject id
7
       int<lower=1, upper=K> item[N];
                                          //item id
8
9
10
11
     parameters {
       vector[2] beta;
                                   //fixed intercept and slope
12
13
       vector[J] u;
                                    //subject intercepts
                                   //item intercepts
       vector[K] w;
14
15
       real<lower=0> sigma e;
                                   //error sd
       real<lower=0> sigma_u;
                                   //subj sd
16
       real<lower=0> sigma_w;
                                   //item sd
17
18
19
     model {
20
      real mu;
21
       //priors
22
                                  //subj random effects
23
          normal(0,sigma_u);
       w ~ normal(0,sigma_w);
                                  //item random effects
24
       // likelihood
       for (i in 1:N){
26
         mu <- beta[1] + u[subj[i]] + w[item[i]] + beta[2]*so[i];</pre>
27
         rt[i] ~ lognormal(mu,sigma_e);
28
29
       }
    }
```

Listing 4: Stan code for running Model 2, the varying intercepts model.

Listing 5: Code for running Model 3, the varying intercepts, varying slopes model.

```
data {
1
2
       int<lower=1> N;
                                           //number of data points
                                           //reading time
       real rt[N];
3
       real<lower=-1,upper=1> so[N];
                                           //predictor
4
                                           //number of subjects
 5
       int<lower=1> J;
       int<lower=1> K;
                                           //number of items
 6
                                           //subject id
       int<lower=1, upper=J> subj[N];
                                           //item id
       int<lower=1, upper=K> item[N];
 8
9
10
11
     parameters {
12
       vector[2] beta;
                                           //intercept and slope
       real<lower=0> sigma_e;
                                           //error sd
13
14
       vector<lower=0>[2] sigma_u;
                                           //subj sd
       vector<lower=0>[2] sigma_w;
                                           //item sd
15
       cholesky_factor_corr[2] L_u;
16
17
       cholesky_factor_corr[2] L_w;
       matrix[2,J] z_u;
18
19
       matrix[2,K] z_w;
20
21
     transformed parameters{
22
       matrix[2,J] u;
23
24
       matrix[2,K] w;
25
       u <- diag pre multiply(sigma u,L u) * z u; //subj random effects
26
         <- diag_pre_multiply(sigma_w,L_w) * z_w; //item random effects
27
28
29
     model {
30
       real mu;
31
32
       //priors
            lkj_corr_cholesky(2.0);
       Lu
33
       L_w ~ lkj_corr_cholesky(2.0);
34
       to_vector(z_u) ~ normal(0,1);
to_vector(z_w) ~ normal(0,1);
35
36
       //likelihood
37
38
       for (i in 1:N){
         mu <- beta[1] + u[1,subj[i]] + w[1,item[i]]</pre>
39
                + (beta[2] + u[2,subj[i]] + w[2,item[i]])*so[i];
40
         rt[i] ~ lognormal(mu,sigma_e);
41
42
       }
```

Listing 6: The Stan code for Model 3, the varying intercepts, varying slopes model.

```
rDat<-read.table("HusainEtAlexpt1data.txt",header=TRUE)
1
     rDat$subj <- with(rDat,factor(subj))</pre>
2
     rDat$item <- with(rDat,factor(item))
3
     X <- unname(model.matrix(~1+so+dist+int, rDat))</pre>
5
6
     stanDat <- within(list(),</pre>
8
9
       N<-nrow(X)
      P <- n_u <- n_w <- ncol(X)
10
11
       X <- X
       z_u < - x
12
       Z_w <- X
13
       J <- length(levels(rDat$subj))</pre>
14
       K <- length(levels(rDat$item))</pre>
15
       rt <- rDat$rt
16
       subj <- as.integer(rDat$subj)</pre>
17
       item <- as.integer(rDat$item)</pre>
18
19
20
     factorialFit <- stan(file="factorialModel.stan",</pre>
21
                             data=stanDat,
22
                             iter=2000, chains=4)
23
```

Listing 7: Preparation of data for analyzing the Husain et al. data-set, and running the model.

```
data {
1
       int<lower=0> N;
                                       //no trials
2
       int<lower=1> P;
                                       //no fixefs
       int<lower=0> J;
                                       //no subjects
4
       int<lower=1> n u;
                                       //no subj ranefs
5
                                       //no items
       int<lower=0> K;
6
       int<lower=1> n w;
                                       //no item ranefs
       int<lower=1,upper=J> subj[N]; //subject indicator
       int<lower=1,upper=K> item[N]; //item indicator
9
       row vector[P] X[N];
                                       //fixef design matrix
10
                                       //subj ranef design matrix
11
       row_vector[n_u] Z_u[N];
       row_vector[n_w] Z_w[N];
                                       //item ranef design matrix
12
       vector[N] rt;
                                       //reading time
13
14
15
     parameters {
                                       //fixef coefs
       vector[P] beta;
16
17
       cholesky factor corr[n u] L u; //cholesky factor of subj ranef corr matrix
       cholesky_factor_corr[n_w] L_w; //cholesky factor of item ranef corr matrix
18
       vector<lower=0>[n_u] sigma_u; //subj ranef std
19
20
       vector<lower=0>[n_w] sigma_w; //item ranef std
       real<lower=0> sigma_e;
                                       //residual std
21
       vector[n_u] z_u[J];
                                       //subj ranef
22
       vector[n_w] z_w[K];
                                       //item ranef
23
24
25
     transformed parameters {
                                       //subj ranefs
       vector[n_u] u[J];
26
       vector[n_w] w[K];
                                       //item ranefs
27
28
         matrix[n_u,n_u] Sigma_u;
                                       //subj ranef cov matrix
29
         matrix[n_w,n_w] Sigma_w;
                                       //item ranef cov matrix
30
         Sigma u <- diag pre multiply(sigma u,L u);</pre>
31
32
         Sigma_w <- diag_pre_multiply(sigma_w,L_w);</pre>
         for(j in 1:J)
33
34
           u[j] <- Sigma_u * z_u[j];
         for(k in 1:K)
35
           w[k] \leftarrow Sigma_w * z_w[k];
36
37
38
     model {
39
       //priors
40
       L u ~ lkj corr cholesky(2.0);
41
      L_w ~ lkj_corr_cholesky(2.0);
42
       for (j in 1:J)
z_u[j] ~ normal(0,1);
43
44
       for (k in 1:K)
45
         z_w[k] normal(0,1);
46
       //likelihood
47
       for (i in 1:N)
48
         rt[i] ~ lognormal(X[i] * beta +
49
                            Z_u[i] * u[subj[i]] +
50
                            Z_w[i] * w[item[i]],
51
52
                            sigma_e);
53
```

Listing 8: Stan code for Husain et al data.