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Accepted for publication in the Journal of Forecasting.

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**Running title:** A Bayesian model to forecast elections from partial information

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**Data availability statement:**

Two datasets are used in this paper. They have been downloaded from <https://eci.gov.in/files/file/12787-bihar-legislative-election-2020/> and <https://dataverse.harvard.edu/file.xhtml?fileId=4788675&version=8.0>.

**Declaration of interest:**

The authors declare no conflict of interest.

# Forecasting Elections from Partial Information Using a Bayesian Model for a Multinomial Sequence of Data

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February 7, 2024

## Abstract

Predicting the winner of an election is of importance to multiple stakeholders. To formulate the problem, we consider an independent sequence of categorical data with a finite number of possible outcomes in each. The data is assumed to be observed in batches, each of which is based on a large number of such trials and can be modelled via multinomial distributions. We postulate that the multinomial probabilities of the categories vary randomly depending on batches. The challenge is to predict accurately on cumulative data based on data up to a few batches as early as possible. On the theoretical front, we first derive sufficient conditions of asymptotic normality of the estimates of the multinomial cell probabilities and present corresponding suitable transformations. Then, in a Bayesian framework, we consider hierarchical priors using multivariate normal and inverse Wishart distributions and establish the posterior convergence. The desired inference is arrived at using these results and ensuing Gibbs sampling. The methodology is demonstrated with election data from two different settings — one from India and the other from the United States of America. Additional insights of the effectiveness of the proposed methodology are attained through a simulation study.

*Keywords: Election data, Gibbs sampling, Hierarchical priors, Posterior convergence, Forecasting from partial information.*

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# 1 Introduction

Voter models have been of interest for a long time. With the strengthening of democracy and the role of media, statistical models for the prediction of elections have been on the rise around the world. Elections are studied at different stages – starting with pre-poll predictions, going on to post-poll predictions, and also calling the election before the final result is declared as the counting of ballots is in progress. The current work focuses on the last context, where the counting of results is publicly disclosed in batches (or rounds). Stemmed from public interest, this often leads to competition among the media houses “to call the election”, i.e. to predict the outcome correctly and as early as possible.

Using the exit poll data has been a popular approach in this context. It however has its own pitfalls, as has been shown in many papers (see, e.g. [Stiers and Dassonneville \[2018\]](#)). Because of the inadequacy of the exit poll data, Associated Press (AP) designed its methodology for both survey and forecasting ([Slodysko \[2020\]](#)). Interestingly, they still do not call a closely contested race unless the trailing candidate cannot mathematically secure a victory. Another common approach is to extract relevant information from social media, such as Twitter or Google Trends, and to use that to predict the winner of an election. [O’Connor et al. \[2010\]](#) and [Tumasjan et al. \[2010\]](#) are two notable works in this regard. A major criticism behind this approach is that intentional bias in the social media posts often lead to error-prone conclusions in the forecasting problems ([Anuta et al. \[2017\]](#)). Further, [Haq et al. \[2020\]](#) provides a great review of the existing literature on how to call an election and one can see that most of the methods are either based on data from social media or are ad-hoc and devoid of sound statistical principles. To that end, it would be of paramount importance to develop a good forecasting methodology based only on the available information of votes secured by the candidates in the completed rounds of counting. This can be treated as a collection of multinomial data where the number of categories is equal to the number of candidates and the random fluctuations in cell probabilities point to the randomness of the vote shares in different rounds.

Multinomial distribution is the most common way to model categorical data. In this work, we study the behaviour of a collection of multinomial distributions with random fluctuations in cell (or ‘category’, as we interchangeably use the two terminology) probabilities. We consider situations where this multinomial data is observed in batches, with the number of trials in each batch being possibly different. The cell probabilities for every multinomial random variable in the data are assumed to be additive combinations of a fixed component which is constant across batches, and a random perturbation. These perturbations for different batches are taken to be independent. Before delving deeper into the model and related results, we present an interesting application where the proposed setup can be utilized effectively.

## 1.1 Motivating Example

This study is largely motivated by the forecasting results in political elections during the process of counting of votes; in particular, we look at one such instance in India – the electoral data from the legislative assembly election held in Bihar (a state in India), during October-November of 2020. There were two major alliances – the National Democratic Alliance (NDA), and the *Mahagathbandhan* (MGB). The counting of votes was declared in different rounds, and the contest being very close, there were widespread speculated forecasts in various media for nearly 20 hours since the counting began, until the final results were eventually known.

To illustrate the challenges of this forecasting exercise, let us consider Hilsa and Parbatta, two of the 243 constituencies in Bihar where the contests were among the closest. In both of them, like

in most other constituencies in the state, the fight was primarily confined between NDA and MGB. For both Hilsa and Parbatta, the counting of votes was completed in 33 rounds. The number of votes counted in the different rounds varied greatly (between two to seven thousand, except the last one or two rounds which typically had less votes). In Figure 1, we depict the number of votes by which the candidates led (roundwise numbers are given on the left panel, cumulative numbers are given in the right panel). For instance, we observe that among the first 10 rounds of counting for Hilsa, the NDA candidate led in the sixth, eighth and ninth rounds, while the MGB candidate led in counting of votes in the remaining seven rounds. Considering cumulative vote counts in Hilsa at the end of successive rounds, lead changed once; that was in the final round with NDA winning by meager 12 votes. In contrast, the lead changed as many as five times in Parbatta, viz. in the 16th, 20th, 23rd, 26th and 31st rounds, before NDA claimed the victory with a margin of 951 votes.

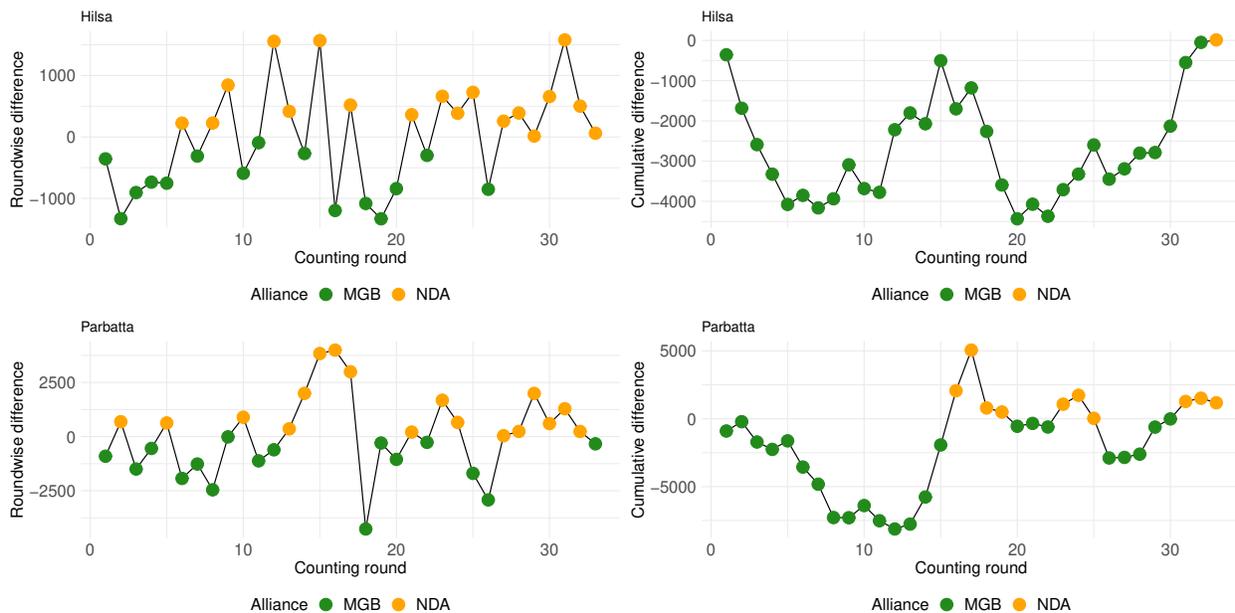


Figure 1: Round-wise margin of differences in votes (left) and cumulative margins of votes (right) for the two main alliances in Hilsa and Matihani, two most closely contested constituencies.

The above clearly demonstrates the difficulty in forecasting the winner (candidate who eventually gets the highest number of votes from all rounds of counting) of individual constituencies, as early (after as few rounds of counting) as possible. Our goal is to construct a method that relies on sound statistical principles to predict the winner for each seat based on trends, thereby making a call about which alliance is going to secure a majority.

We note that the methodology can be adapted to other political applications as well. For example, in the context of the presidential election of the United States of America (USA), one may consider the data from a few counties to project the winner of the respective state. We illustrate it later in this paper, although in a limited capacity because of not having complete information on the actual timing when the county-wise results were progressively declared.

## 1.2 Related literature and our contribution

Literature on a non-identical sequence of categorical data is somewhat limited. In the binomial setting (i.e. with two categories of outcome), [Le Cam \[1960\]](#) introduced the concept of independent

but non-identical random indicators, which was later discussed and applied in a few other studies (see, e.g. Volkova [1996], Hong et al. [2009], Fernández and Williams [2010]). However, in these papers, the cell probabilities are considered to be different but non-random. Introducing the randomness in the probabilities makes it more attractive and suitable for many real life applications. To that end, Cargnoni et al. [1997] proposed a conditionally Gaussian dynamic model for multinomial time series data and used it to forecast the flux of students in the Italian educational system. This model has later been adopted and modified in different capacities, see Landim and Gamerman [2000], Terui et al. [2010], Quinn et al. [2010], Terui and Ban [2014] for example.

In this paper, we develop an appropriate framework for the above-mentioned setup under mild assumptions on the behaviour of the random perturbations. The model is implemented through Bayesian techniques in a setting similar to Cargnoni et al. [1997]. We complement their work by providing relevant theoretical results of the proposed method and describe how the model can be used to forecast and infer about different features of the entire data. Next, in line with the motivations discussed earlier, in addition to simulation, we demonstrate with two contextual applications of our proposed methodology. First one is related to election prediction where we use the data from Bihar legislative assembly election held in 2020. The second is for US election; although in absence of availability of sequencing of the actual data, we consider different hypothetical possibilities. To the best of our knowledge, in the context of these applications, the present work is the first attempt to utilize a randomly varying probability structure for a sequence of non-identical multinomial distributions. We emphasize that the goal of this study is to forecast aspects of the final outcome rather than inference on the multinomial cell probabilities, although the latter can also be easily done through the proposed framework.

### 1.3 Organization

The rest of the paper is arranged as follows. In Section 2.1, we introduce the preliminary notations, and in Section 2.2, we describe the model in a formal statistical framework, and state the main results. The proofs of the theorems are provided in Section 7. In Section 2.3, we discuss the estimation of the posterior distribution of the model parameters via Gibbs sampling. The prediction aspects of the model are described in Section 2.4. Section 3 discusses simulation results which demonstrate the suitability of the model in the broad context. Real data application in an Indian context is presented in Section 4. We also present a hypothetical case study from American politics in Section 5. Finally, we conclude with a summary, some related comments and ways to extend the current work in Section 6.

## 2 Methods

### 2.1 Preliminaries

In this section, we describe the notations and assumptions, and then for better understanding of the reader, draw necessary connections to the previously discussed examples.

At the core of the framework, we have a sequence of independent trials, each of which results in one among the possible  $C$  categories of outcomes. The outcomes of the trials are recorded in aggregated form, which we refer to as batches (or rounds). We assume that the cell probabilities remain the same within a batch, but vary randomly across different batches. Hence, the observed data from an individual batch can be modelled by the standard multinomial distribution; and collectively we have a sequence of independent but non-identical multinomial random variables.

We adopt the following notations. Altogether, the multinomial data is available in  $K$  batches. For each batch, the number of trials is given by  $n_j$ ,  $1 \leq j \leq K$ , and each  $n_j$  is assumed to be large. Let  $N_j = \sum_{i=1}^j n_i$  denote the cumulative number of trials up to the  $j^{\text{th}}$  batch. For simplicity, let  $N = N_K$ , the total number of trials in all the batches. We use  $\mathbf{X}_j = (X_{1j}, X_{2j}, \dots, X_{Cj})^T$  to denote the observed counts for the  $C$  different categories in case of the  $j^{\text{th}}$  multinomial variable in the data. Also, let  $\mathbf{Y}_j = (Y_{1j}, Y_{2j}, \dots, Y_{Cj})^T$  be the cumulative counts till the  $j^{\text{th}}$  batch. Clearly,

$$\sum_{i=1}^j X_{ci} = Y_{cj}, \quad \text{for } c \in \{1, 2, \dots, C\}. \quad (2.1)$$

Let  $p_{1j}, p_{2j}, \dots, p_{Cj}$  denote the probabilities for the  $C$  categories for each of the independent trials in the  $j^{\text{th}}$  batch. We focus on  $\{p_{1j}, p_{2j}, \dots, p_{(C-1)j}\}$ , since the probability of the last category is then automatically defined. Our objective is to estimate the probability distribution, and subsequently various properties, of  $(h(\mathbf{Y}_l) \mid \mathcal{F}_j)$  for  $l > j$ , where  $h$  is some suitable measurable function and  $\mathcal{F}_j$  is the sigma-field generated by the data up to the  $j^{\text{th}}$  batch.

If we focus on the aforementioned voting context of Bihar (or, the USA),  $C$  stands for the number of candidates contesting in a constituency (or, a state), and each voter casts a vote in favour of exactly one of these candidates. Here,  $n_j$  represents the number of voters whose votes are counted in the  $j^{\text{th}}$  round (or, the  $j^{\text{th}}$  county), while  $N_j$  denotes the corresponding cumulative number of votes. Similarly,  $\mathbf{X}_j$  shows the number of votes received by the candidates in the  $j^{\text{th}}$  round (or, the  $j^{\text{th}}$  county), and  $\mathbf{Y}_j$  represents the corresponding vector of cumulative votes. We can use our method to find the probability of winning for the  $i^{\text{th}}$  candidate, given the counted votes from the first  $j$  number of rounds (or, counties). In this case, the measurable function of interest is  $h_{1,i}(\mathbf{Y}_K) = \mathbb{I}(Y_{iK} > \max_{l \neq i} Y_{lK})$ , where  $\mathbb{I}(\cdot)$  denotes the indicator function. Similarly, using the function  $h_2(\mathbf{Y}_K) = \mathbf{Y}_K^{(1)} - \mathbf{Y}_K^{(2)}$ , where  $\mathbf{Y}_K^{(1)}$  and  $\mathbf{Y}_K^{(2)}$  respectively denote the first and second order statistics of  $\mathbf{Y}_K$ , we can predict the margin of victory as well.

## 2.2 Model framework and main results

We propose to consider the multinomial cell probabilities as random variables, having a fixed part and a randomly varying part. Thus, for  $c = 1, 2, \dots, C$ ,  $p_{cj}$  is written as  $p_c + \varepsilon_{cj}$ , where  $p_c$  is the constant preference component that remains the same across the batches for category  $c$ , and  $\varepsilon_{cj}$ 's, for  $c \in \{1, 2, \dots, C\}$ , are zero-mean random perturbations that model possible fluctuations in the preference probabilities for the categories across the batches. We enforce  $\varepsilon_{Cj} = -\sum_{c=1}^{C-1} \varepsilon_{cj}$ . In line with the earlier notations, we use  $\mathbf{p} = (p_1, p_2, \dots, p_C)^T$ ,  $\tilde{\mathbf{p}} = (p_1, p_2, \dots, p_{C-1})^T$ ,  $\mathbf{p}_j = (p_{1j}, p_{2j}, \dots, p_{Cj})^T$  and  $\boldsymbol{\varepsilon}_j = (\varepsilon_{1j}, \varepsilon_{2j}, \dots, \varepsilon_{(C-1)j})^T$ , for convenience. Note that the covariances between the components of  $\boldsymbol{\varepsilon}_j$ 's are likely to be negative due to the structure of the multinomial distribution. The randomness of  $\boldsymbol{\varepsilon}_j$ 's need to be also carefully modelled so as to ensure that the category probabilities  $p_{ij}$ 's lie in the interval  $[0, 1]$ . We shall use  $\mathcal{S}$  to denote the possible set of values for  $\boldsymbol{\varepsilon}_j$ .

Let  $\text{Cov}(\cdot, \cdot)$  denote the variance-covariance matrix of two random variables and  $\mathcal{N}_r(\boldsymbol{\theta}, \Psi)$  denote a  $r$ -variate normal distribution with mean  $\boldsymbol{\theta}$  and dispersion matrix  $\Psi$ . Throughout this article,  $\mathbf{0}$  denotes a vector of all zeroes and  $\mathbf{I}$  denotes an identity matrix of appropriate order. Following is a critical assumption that we use throughout the paper.

**Assumption 1.** *The random variables  $(\boldsymbol{\varepsilon}_j)_{1 \leq j \leq K}$ , as defined above, are independent of each other and are distributed on  $\mathcal{S}$ , with  $\mathbb{E}(\boldsymbol{\varepsilon}_{cj}) = 0$  for  $1 \leq c \leq C-1$ . Further, the density of  $\sqrt{n_j} \boldsymbol{\varepsilon}_j$  converges uniformly to the density of  $\mathcal{N}_{C-1}(\mathbf{0}, \Xi_\varepsilon)$ , for some positive definite matrix  $\Xi_\varepsilon$ , typically unknown.*

Then, the proposed model can be written as

$$\mathbf{X}_j | \boldsymbol{\varepsilon}_j \sim \text{Multinomial} \left( n_j, p_1 + \varepsilon_{1j}, p_2 + \varepsilon_{2j}, \dots, p_{C-1} + \varepsilon_{(C-1)j}, p_C - \sum_{c=1}^{C-1} \varepsilon_{cj} \right), \quad (2.2)$$

where  $(\mathbf{X}_j)_{1 \leq j \leq K}$  are independent of each other and  $(\boldsymbol{\varepsilon}_j)_{1 \leq j \leq K}$  satisfy Assumption 1.

We shall adopt a Bayesian framework to implement the above model. Before that, it is imperative to present a couple of frequentist results on the distribution of the votes, as they help us in setting up the problem, and serve as the basis of the framework and the results stated in Section 2.3. Below, Proposition 1 shows the prior distribution of the voting percentages, while Lemma 1 helps us in getting the distribution of the transformed prior that we actually use in our Bayesian hierarchical models.

**Proposition 1.** *For the proposed model (eq. (2.2)) the following results are true if Assumption 1 is satisfied.*

- (a) *Unconditional first and second order moments of  $X_{cj}$ ,  $X_{c'j}$  ( $c, c' = 1, 2, \dots, C$ ,  $c \neq c'$ ) are given by*

$$\mathbb{E}(X_{cj}) = n_j p_c, \quad \text{Cov}(X_{cj}, X_{c'j}) = n_j \begin{bmatrix} p_c(1-p_c) & -p_c p_{c'} \\ -p_c p_{c'} & p_{c'}(1-p_{c'}) \end{bmatrix} + n_j(n_j - 1) \text{Cov}(\varepsilon_{cj}, \varepsilon_{c'j}). \quad (2.3)$$

- (b) *Unconditional first and second order moments of  $Y_{cj}$ ,  $Y_{c'j}$  ( $c, c' = 1, 2, \dots, C$ ,  $c \neq c'$ ) are given by*

$$\mathbb{E}(Y_{cj}) = N_j p_c, \quad \text{Cov}(Y_{cj}, Y_{c'j}) = N_j \begin{bmatrix} p_c(1-p_c) & -p_c p_{c'} \\ -p_c p_{c'} & p_{c'}(1-p_{c'}) \end{bmatrix} + \sum_{i=1}^j n_i(n_i - 1) \text{Cov}(\varepsilon_{ci}, \varepsilon_{c'i}). \quad (2.4)$$

- (c) *As  $n_j \rightarrow \infty$ ,  $\hat{\mathbf{p}}_j = (\hat{p}_{1j}, \hat{p}_{2j}, \dots, \hat{p}_{(C-1)j})^T = (X_{1j}/n_j, X_{2j}/n_j, \dots, X_{(C-1)j}/n_j)^T$  satisfies the following:*

$$\sqrt{n_j}(\hat{\mathbf{p}}_j - \tilde{\mathbf{p}}) \xrightarrow{\mathcal{L}} \mathcal{N}_{C-1}(\mathbf{0}, \Xi), \quad (2.5)$$

where  $\xrightarrow{\mathcal{L}}$  denotes convergence in law, and  $\Xi = \Xi_p + \Xi_\varepsilon$ . Here,  $\Xi_\varepsilon$  is positive definite with finite norm and

$$\Xi_p = \begin{bmatrix} p_1(1-p_1) & -p_1 p_2 & \dots & -p_1 p_{C-1} \\ -p_2 p_1 & p_2(1-p_2) & \dots & -p_2 p_{C-1} \\ \vdots & \vdots & \ddots & \vdots \\ -p_{C-1} p_1 & -p_{C-1} p_2 & \dots & p_{C-1}(1-p_{C-1}) \end{bmatrix}. \quad (2.6)$$

**Remark 1.** *In part (c) of Proposition 1, if the variances and covariance of all components of  $\boldsymbol{\varepsilon}_j$  are  $o(1/n_j)$  then as  $n_j \rightarrow \infty$ ,  $\sqrt{n_j}(\hat{\mathbf{p}}_j - \tilde{\mathbf{p}}) \xrightarrow{\mathcal{L}} \mathcal{N}_{C-1}(\mathbf{0}, \Xi_p)$ .*

The above remark includes the scenario of independent and identically distributed multinomials and is therefore of special interest. Detailed proof of Proposition 1 is deferred to Section 7. Note that the asymptotic variance is a function of  $\mathbf{p}$ , thereby motivating us to adapt an appropriate adjustment in the form of using a suitable variance stabilizing transformation. While the standard

variance stabilizing transformation is given by the sine-inverse function (Anscombe [1948]), there are some better variance stabilizing transformations, as discussed in Yu [2009]. Based on that work, we define the following modified transformation, having superior relative errors, skewness and kurtosis:

$$\mathbf{L}_j = \begin{pmatrix} L_{1j} \\ L_{2j} \\ \vdots \\ L_{(C-1)j} \end{pmatrix} = \begin{pmatrix} \sin^{-1} \frac{2\hat{p}_{1j} - 1}{1 + 2a/n_j} \\ \sin^{-1} \frac{2\hat{p}_{2j} - 1}{1 + 2a/n_j} \\ \vdots \\ \sin^{-1} \frac{2\hat{p}_{(C-1)j} - 1}{1 + 2a/n_j} \end{pmatrix}, \quad (2.7)$$

where  $a$  is a positive constant. Throughout this paper, we consider  $a = 3/8$ , one of the most popular choices in this regard, cf. Anscombe [1948] and Yu [2009].

**Lemma 1.** *For every  $1 \leq j \leq K$  in the framework of the proposed model (eq. (2.2)), under Assumption 1*

$$\sqrt{n_j + 0.5}(\mathbf{L}_j - \boldsymbol{\mu}) \xrightarrow{\mathcal{L}} \mathcal{N}_{C-1}(\mathbf{0}, \Sigma), \quad \text{as } n_j \rightarrow \infty, \quad (2.8)$$

where  $\boldsymbol{\mu} = (\sin^{-1}(2p_1 - 1), \sin^{-1}(2p_2 - 1), \dots, \sin^{-1}(2p_{C-1} - 1))$ , and  $\Sigma$  is a variance-covariance matrix with  $\|\Sigma\| < \infty$ .

In particular, if the variances and covariances of  $\varepsilon_{1j}, \dots, \varepsilon_{(C-1)j}$  are  $o(1/n_j)$  then  $\Sigma$  is a variance-covariance matrix with diagonal entries as 1, and off-diagonal entries given by

$$\Sigma_{cc'} = -\sqrt{\frac{p_c p_{c'}}{(1-p_c)(1-p_{c'})}} \text{ for } c \neq c' \in \{1, 2, \dots, C-1\}.$$

**Proof.** For a function  $g : \mathbb{R}^{C-1} \rightarrow \mathbb{R}^{C-1}$  of the form:

$$g(x_1, x_2, \dots, x_{C-1}) = \begin{pmatrix} g_1(x_1) \\ g_2(x_2) \\ \vdots \\ g_{C-1}(x_{C-1}) \end{pmatrix},$$

using eq. (2.5) and the multivariate delta theorem (see Cox [2005] and Ver Hoef [2012] for the theorem and related discussions) we get

$$\sqrt{n_j} (g(\hat{p}_{1j}, \hat{p}_{2j}, \dots, \hat{p}_{(C-1)j}) - g(p_1, p_2, \dots, p_{C-1})) \xrightarrow{\mathcal{L}} \mathcal{N}_{C-1}(\mathbf{0}, \Omega), \quad (2.9)$$

where the  $(c, c')^{th}$  entry of  $\Omega$  is

$$\Omega_{c,c'} = \begin{cases} (g'_c(p_c))^2 \Xi_{c,c} & \text{for } c = c', \\ -g'_c(p_c)g'_{c'}(p_{c'}) \Xi_{c,c'} & \text{for } c \neq c'. \end{cases}$$

Using  $g_1(t) = g_2(t) = \dots = g_{C-1}(t) = \sin^{-1}((2t-1)/(1+2a/n_j))$  in the above, the required result follows. For the special case in the second part, the exact form of  $\Xi_p$  from eq. (2.6) can be used.  $\square$

The above lemma serves as a key component of our method elaborated in the following subsec-

tion. Note that  $g_1(t)$  is a monotonic function on  $[0, 1]$  and therefore, we can easily use the above result to make inference about  $\tilde{\boldsymbol{p}}$ , and subsequently about  $(h(\mathbf{Y}_l) \mid \mathcal{F}_j)$  for  $l > j$ .

### 2.3 Bayesian estimation

In this paper, for implementation of the model, we adopt a Bayesian framework which is advantageous from multiple perspectives. First, we find this to yield more realistic results in several real-life data context. Second, it helps in reducing the computational complexity, especially for a large dataset. Third, this approach has the flexibility to naturally extend to similar problems in presence of covariates.

Following eq. (2.8) and the notations discussed above, the proposed model (see eq. (2.2)) is equivalent to the following in an asymptotic sense.

$$\mathbf{L}_j \sim \mathcal{N}_{C-1} \left( \boldsymbol{\mu}, \frac{\Sigma}{n_j + 0.5} \right), \quad 1 \leq j \leq K. \quad (2.10)$$

Observe that both  $\boldsymbol{\mu}$  and  $\Sigma$  are unknown parameters and are fixed across different batches. In the Bayesian framework, appropriate priors need to be assigned to these parameters to ensure “good” behaviour of the posterior distributions. To that end, we consider the following hierarchical structure of the prior distributions.

$$\boldsymbol{\mu} \sim \mathcal{N}_{C-1}(\boldsymbol{\alpha}, \Sigma_p), \quad \Sigma \sim \text{Inverse Wishart}(\Psi, \nu), \quad \Sigma_p \sim \text{Inverse Wishart}(\Psi_p, \nu_p). \quad (2.11)$$

The inverse Wishart prior is the most natural conjugate prior for the covariance matrix, cf. Chen [1979], Haff [1980], Barnard et al. [2000], Champion [2003]. The conjugacy property facilitates amalgamation into Markov chain Monte Carlo (MCMC) methods based on Gibbs sampling. This leads to easier simulations from the posterior distribution as well as posterior predictive distribution after each round. On this note, the inverse Wishart distribution, if imposed only on  $\Sigma$  with  $\Sigma_p$  known, restricts the flexibility in terms of modeling prior knowledge (Hsu et al. [2012]). To address this issue, we apply the above hierarchical modeling as is most commonly used in Bayesian computations. This allows greater flexibility and stability than diffuse priors (Gelman et al. [2006]). Studies by Kass et al. [2006], Gelman and Hill [2006], Bouriga and Féron [2013] discuss in depth how the hierarchical inverse Wishart priors can be used effectively for modeling variance-covariance matrices.

Before outlining the implementation procedure of the above model, we present an important result which establishes the large sample property of the posterior means of  $\boldsymbol{\mu}$  and  $\Sigma$ .

**Theorem 1.** *Let  $\boldsymbol{\mu}_0$  and  $\Sigma_0$  be the true underlying mean and true covariance matrix, respectively. Let  $\boldsymbol{\mu}_{PM}$  and  $\Sigma_{PM}$  be the posterior means. Then, for the hierarchical prior distributions given by eq. (2.11), the following are true:*

$$\lim_{K \rightarrow \infty} \mathbb{E}[\Pi_K(\|\boldsymbol{\mu}_{PM} - \boldsymbol{\mu}_0\| > \epsilon \mid \mathbf{L}_1, \mathbf{L}_2, \dots, \mathbf{L}_K)] = 0, \quad (2.12)$$

$$\lim_{K \rightarrow \infty} \mathbb{E}[\Pi_K(\|\Sigma_{PM} - \Sigma_0\| > \epsilon \mid \mathbf{L}_1, \mathbf{L}_2, \dots, \mathbf{L}_K)] = 0. \quad (2.13)$$

In terms of implementation, it is usually a complicated procedure to find out a closed form for the joint posterior distribution of the parameters. A common approach is to adopt Gibbs sampling which reduces the computational burden significantly. It is an MCMC method to obtain a sequence of realizations from a joint probability distribution. Here, every parameter is updated in an iterative

manner using the conditional posterior distributions given other parameters. We refer to [Geman and Geman \[1984\]](#) and [Durbin and Koopman \[2002\]](#) for more in-depth readings on Gibbs sampling.

Recall that  $\mathcal{F}_j$  denotes the sigma-field generated by the data up to the  $j^{\text{th}}$  instance. In addition, below,  $\Gamma_k$  stands for a  $k$ -variate gamma function and  $\text{tr}(\cdot)$  is the trace of a matrix.

For  $j = 1$ , we have  $\mathbf{L}_1 | \boldsymbol{\mu}, \Sigma \sim \mathcal{N}_{C-1}(\boldsymbol{\mu}, \Sigma / (n_1 + 0.5))$ . The joint posterior likelihood is therefore given by

$$\begin{aligned}
f(\boldsymbol{\mu}, \Sigma, \Sigma_p | \mathcal{F}_1) &\propto |\Sigma|^{-1/2} \exp \left[ -\frac{n_1 + 0.5}{2} (\mathbf{L}_1 - \boldsymbol{\mu})^T \Sigma^{-1} (\mathbf{L}_1 - \boldsymbol{\mu}) \right] \\
&\times |\Sigma_p|^{-1/2} \exp \left[ -\frac{1}{2} (\boldsymbol{\mu} - \boldsymbol{\alpha})^T \Sigma_p^{-1} (\boldsymbol{\mu} - \boldsymbol{\alpha}) \right] \\
&\times \frac{|\Psi|^{\nu/2}}{2^\nu \Gamma_2(\nu/2)} |\Sigma|^{-(\nu+C)/2} \exp \left[ -\frac{1}{2} \text{tr}(\Psi \Sigma^{-1}) \right] \\
&\times \frac{|\Psi_p|^{\nu_p/2}}{2^{\nu_p} \Gamma_2(\nu_p/2)} |\Sigma_p|^{-(\nu_p+C)/2} \exp \left[ -\frac{1}{2} \text{tr}(\Psi_p \Sigma_p^{-1}) \right]. \tag{2.14}
\end{aligned}$$

In an exact similar way, the posterior likelihood based on the data up to the  $j^{\text{th}}$  instance can be written as

$$\begin{aligned}
f(\boldsymbol{\mu}, \Sigma, \Sigma_p | \mathcal{F}_j) &\propto |\Sigma|^{-j/2} \exp \left[ -\sum_{i=1}^j \frac{n_i + 0.5}{2} (\mathbf{L}_i - \boldsymbol{\mu})^T \Sigma^{-1} (\mathbf{L}_i - \boldsymbol{\mu}) \right] \\
&\times |\Sigma_p|^{-1/2} \exp \left[ -\frac{1}{2} (\boldsymbol{\mu} - \boldsymbol{\alpha})^T \Sigma_p^{-1} (\boldsymbol{\mu} - \boldsymbol{\alpha}) \right] \\
&\times \frac{|\Psi|^{\nu/2}}{2^\nu \Gamma_2(\nu/2)} |\Sigma|^{-(\nu+C)/2} \exp \left[ -\frac{1}{2} \text{tr}(\Psi \Sigma^{-1}) \right] \\
&\times \frac{|\Psi_p|^{\nu_p/2}}{2^{\nu_p} \Gamma_2(\nu_p/2)} |\Sigma_p|^{-(\nu_p+C)/2} \exp \left[ -\frac{1}{2} \text{tr}(\Psi_p \Sigma_p^{-1}) \right]. \tag{2.15}
\end{aligned}$$

In the Gibbs sampler, we need to use the conditional posterior distributions of  $\boldsymbol{\mu}$ ,  $\Sigma$  and  $\Sigma_p$ . It is easy to note that

$$f(\Sigma | \boldsymbol{\mu}, \Sigma_p, \mathcal{F}_j) \propto |\Sigma|^{-(\nu+j+C)/2} \exp \left[ -\frac{1}{2} \text{tr} \left( \Psi \Sigma^{-1} + \sum_{i=1}^j (n_i + 0.5) (\mathbf{L}_i - \boldsymbol{\mu})(\mathbf{L}_i - \boldsymbol{\mu})^T \Sigma^{-1} \right) \right]. \tag{2.16}$$

Thus, we can write the following.

$$\Sigma | \boldsymbol{\mu}, \Sigma_p, \mathcal{F}_j \sim \text{Inverse Wishart} \left( \Psi + \sum_{i=1}^j (n_i + 0.5) (\mathbf{L}_i - \boldsymbol{\mu})(\mathbf{L}_i - \boldsymbol{\mu})^T, \nu + j \right). \tag{2.17}$$

In an identical fashion, it is possible to show that

$$\Sigma_p | \boldsymbol{\mu}, \Sigma, \mathcal{F}_j \sim \text{Inverse Wishart} (\Psi_p + (\boldsymbol{\mu} - \boldsymbol{\alpha})(\boldsymbol{\mu} - \boldsymbol{\alpha})^T, \nu_p + 1). \tag{2.18}$$

For the conditional posterior distribution of  $\boldsymbol{\mu}$ , observe that

$$\begin{aligned}
& f(\boldsymbol{\mu}|\Sigma, \Sigma_p, \mathcal{F}_j) \\
& \propto \exp \left[ -\frac{1}{2} \left( \sum_{i=1}^j (n_i + 0.5) (\mathbf{L}_i - \boldsymbol{\mu})^T \Sigma^{-1} (\mathbf{L}_i - \boldsymbol{\mu}) + (\boldsymbol{\mu} - \boldsymbol{\alpha})^T \Sigma_p^{-1} (\boldsymbol{\mu} - \boldsymbol{\alpha}) \right) \right] \\
& \propto \exp \left[ -\frac{1}{2} \left( \boldsymbol{\mu}^T [\Sigma_p^{-1} + (N_j + j/2)\Sigma^{-1}] \boldsymbol{\mu} - \boldsymbol{\mu}^T \left[ \Sigma_p^{-1} \boldsymbol{\alpha} + \Sigma^{-1} \left( \sum_{i=1}^j (n_i + 0.5) \mathbf{L}_i \right) \right] \right) \right]. \quad (2.19)
\end{aligned}$$

Let  $V_j = \Sigma_p^{-1} + (N_j + j/2)\Sigma^{-1}$ . Comparing the above expression with the density of multivariate normal distribution, and readjusting the terms as necessary, we can show that

$$\boldsymbol{\mu}|\Sigma, \Sigma_p, \mathcal{F}_j \sim \mathcal{N}_{C-1} \left( V_j^{-1} \left[ \Sigma_p^{-1} \boldsymbol{\alpha} + \Sigma^{-1} \left( \sum_{i=1}^j (n_i + 0.5) \mathbf{L}_i \right) \right], V_j^{-1} \right). \quad (2.20)$$

Existing theory on Bayesian inference ensures that if the above conditional posterior distributions are iterated many times, it would converge to the true posterior distributions for the parameters. It is naturally of prime importance to make sure that convergence is achieved when we implement the algorithm and collect a posterior sample. In that regard, we use the Gelman-Rubin statistic, cf. [Gelman et al. \[1992\]](#). This is an efficient way to monitor the convergence of the Markov chains. Here, multiple parallel chains are initiated from different starting values. Following the standard theory of convergence, all of these chains eventually converge to the true posterior distributions and hence, after enough iterations, it should be impossible to distinguish between different chains. Leveraging this idea and applying an ANOVA-like technique, the Gelman-Rubin statistic compares the variation between the chains to the variation within the chains. Ideally, this statistic converges to 1, and a value close to 1 indicates convergence. In all our applications, we start multiple chains by randomly generating initial values for the parameters and start collecting samples from the posterior distributions only when the value of the Gelman-Rubin statistic is below 1.1. Also, we take samples from iterations sufficiently apart from each other so as to ensure independence of the realizations.

## 2.4 Prediction

In this section, we discuss the prediction procedure for the proposed model. Note that, because of the conjugacy we observed above, the posterior predictive distribution for the  $l^{th}$  instance based on the data up to the  $j^{th}$  instance ( $l > j$ ) is given by a normal distribution. The mean parameter of this distribution can be computed by the following equation:

$$\mathbb{E}(\mathbf{L}_l|\mathcal{F}_j) = \mathbb{E}(\mathbb{E}(\mathbf{L}_l|\boldsymbol{\mu}, \mathcal{F}_j)) = \mathbb{E}(\boldsymbol{\mu}|\mathcal{F}_j). \quad (2.21)$$

Similarly, using the property of conditional variance, the dispersion parameter of the posterior predictive distribution is

$$\text{Cov}(\mathbf{L}_l|\mathcal{F}_j) = \mathbb{E}[\text{Cov}(\mathbf{L}_l|\boldsymbol{\mu}, \Sigma, \mathcal{F}_j)] + \text{Cov}(\mathbb{E}(\mathbf{L}_l|\boldsymbol{\mu}, \Sigma, \mathcal{F}_j)) = \frac{\mathbb{E}(\Sigma|\mathcal{F}_j)}{n_l + 0.5} + \text{Cov}(\boldsymbol{\mu}|\mathcal{F}_j). \quad (2.22)$$

Once again, it is complicated to get closed form expressions for the above expectation and variance. Therefore, we make use of the Gibbs sampler to simulate realizations from the posterior predictive distribution. Based on the data up to the  $j^{th}$  instance, we generate  $M$  samples (for

large  $M$ ) for the parameters from the posterior distributions. Let us call them  $\Sigma_{j,s}, \boldsymbol{\mu}_{j,s}$  for  $s = 1, 2, \dots, M$ . Then, we can approximate the mean of the posterior predictive distribution in eq. (2.21) by  $\sum_{s=1}^M \boldsymbol{\mu}_{j,s}/M$ . Similarly, the first term in the right hand side of eq. (2.22) can be approximated by  $\hat{\Sigma}_j/(n_l + 0.5)$  where  $\hat{\Sigma}_j$  is the sample mean of  $\Sigma_{j,s}$  for  $s = 1, 2, \dots, M$ , while the second term in that equation can be approximated by the sample dispersion matrix of  $\boldsymbol{\mu}_{j,s}$  for  $s = 1, 2, \dots, M$ . Next, we generate many samples (once again, say  $M$ ) from the posterior predictive distribution with these estimated parameters.

This sample from the posterior predictive distribution is then used to infer on various properties of  $(h(\mathbf{Y}_l) \mid \mathcal{F}_j)$ , as discussed in Section 2.1.

### 3 Simulation Study

In this section, we consider a few toy election scenarios and evaluate the effectiveness of our proposed methodology across various scenarios of elections. We generate  $K$  batches of election data as multinomial random draws following eq. (2.2) for three different simulated election outcomes (SEO). Each election data is assumed to have  $C$  candidates. The expected values of  $n_j$  (the number of votes cast in the  $j^{\text{th}}$  batch) are considered to be the same for all  $j$ . Let  $n$  denote the common expected value for different batches. We consider different value of  $C$  (3 or 5),  $K$  (25 or 50),  $n$  (100 or 1000 or 5000 or 50000) to understand the robustness of the method.

The SEOs we consider differ in the way the  $\boldsymbol{\varepsilon}_j$ 's are generated. First, we take the simplest situation of  $\varepsilon_{cj} = 0$  for all  $c, j$ , which corresponds to an iid collection of multinomial random variables. Second,  $\boldsymbol{\varepsilon}_j$  is simulated from a truncated multivariate normal distribution with mean  $\mathbf{0}$  and dispersion matrix  $n_j^{-0.5}\mathbf{I}$ . Third, we impose dependence across different coordinates of  $\boldsymbol{\varepsilon}_j$ , and simulate it from a truncated multivariate normal distribution with mean  $\mathbf{0}$  and dispersion matrix  $n_j^{-0.5}A$ , where  $A$  is a randomly generated positive definite matrix with finite norm. Observe that all the three SEOs satisfy Assumption 1. For every combination of  $K, C, n$ , each experiment is repeated many times and we compute the average performance over these repetitions.

In this simulation study, we particularly explore two aspects in detail. First, we find out the accuracy of our methodology in terms of predicting the candidate with maximum votes at the end. Here, a decision is made when the winner is predicted with at least 99.5% probability and when the predicted difference in votes for the top two candidates is more than 5% of the remaining votes to be counted. Second, we focus on two specific cases which are similar to our real-life applications, and examine how well we can predict various properties of  $h(\mathbf{Y}_l)$ , for different choices of  $h(\cdot)$ . One of these cases has  $C = 5, K = 50$  and we look for different options for  $n$ . In the other case we have  $C = 3, K = 25, n = 5000$ .

In case of the first problem, for every iteration, the fixed values of  $\boldsymbol{p}$  are obtained from a Dirichlet distribution with equal parameters to ensure a more general view of the performance of the proposed method. Detailed results are reported in Appendix A, in Tables A1, A2, and A3. We note that the results are not too sensitive with respect to the choices of  $C$  and  $K$ , but the accuracy increases with  $n$ . Below, in Table 1, we look at the results for  $C = 5, K = 50$  for the three different SEOs. It is evident that under the iid assumption (SEO1), the method makes a correct call almost all the times. When the components of  $\boldsymbol{\varepsilon}_j$  are independent random variables (SEO2), a correct call is made more than 90% of the times for large enough samples ( $n > 5000$ ). For the third SEO, where the components of  $\boldsymbol{\varepsilon}_j$  are simulated from a multivariate distribution with a general positive definite matrix as the covariance matrix, the accuracy of making a correct call drops to 70% for  $n = 50000$ . Another interesting observation is that the accuracy depends heavily on the average final margin between the top two categories. In other words, if the underlying probabilities of the

Table 1: Accuracy (in %) of predicting the category with maximum votes (corresponding final margins, averaged over all repetitions, are given in parentheses) for different SEOs. All results correspond to the case of  $C = 5$ ,  $K = 50$ .

SEO	$n$	Correct (avg margin)	Incorrect (avg margin)	No call (avg margin)
SEO1	100	94.5% (1584)	1.5% (73)	4% (168)
	1000	98.25% (14388)	0.5% (516)	1.25% (378)
	5000	99.5% (71562)	0.25% (1461)	0.25% (858)
	50000	100% (730428)		
SEO2	100	70% (605)	8.25% (205)	21.75% (314)
	1000	84.75% (10986)	4.75% (1286)	10.5% (2891)
	5000	91.75% (67476)	3.25% (6624)	5% (12241)
	50000	95.5% (720905)	2.25% (71467)	2.25% (77728)
SEO3	100	42.4% (154)	22% (106)	35.6% (110)
	1000	45% (2439)	23.2% (1423)	31.8% (1536)
	5000	55.8% (18972)	14.2% (9671)	30% (10822)
	50000	70% (305700)	8.8% (79016)	21.2% (184868)

top two candidates are close (which directly relates to a lower margin of difference between the final counts), then the method achieves lower accuracy, and vice-versa. We also observe that in those cases, the method records *no call* (i.e. a decision cannot be made with the prescribed rules) more often than incorrect calls.

Moving on to the particular choice of  $C = 3, K = 25, n = 5000$ , we aim to find out the effectiveness of the proposed approach in identifying the leading candidates for varying degrees of difference between the true vote probabilities of the topmost two candidates. To that end, for  $\mathbf{p} = (p_1, p_2, p_3)$ , without loss of generality, we assume  $p_1 > p_2 > p_3$ . Now, data are generated for the above three simulated elections by fixing  $p_1 - p_2 = \delta$  where  $\delta \in \{0.01, 0.05, 0.1, 0.25\}$ . For every choice of  $\delta$  and for every SEO, we repeat the experiments and find out the mean accuracy in predicting the top candidate. Additionally, we also find out the average number of observations used by the method for making the prediction. These results are displayed in Table 2.

It is evident that for SEO1, the method provides great prediction. Even when there is only 1% difference in the probabilities of the top two candidates, the method predicts the leading candidate correctly 96% of the times. For larger than 1% difference, it never fails to predict the leading candidate correctly. For SEO2, our method can predict the topmost candidate with more than 75% accuracy whenever the probabilities of the top two candidates differ by at least 5%. This accuracy reaches the value of nearly 90% for  $\delta = 0.1$  and is 100% if the difference in those two cell probabilities is 0.25. For the third SEO, the prediction accuracy is about 70% for  $\delta = 0.1$ . The accuracy improves steadily as  $\delta$  increases. One can say that the method will be able to predict the winner with high level of accuracy whenever there is a considerable difference between the  $p_i$  values for the top two candidates. If that difference is minute, which corresponds to a very closely contested election, the accuracy will drop. Additionally, we observe that in the most extreme cases, about 20% of the times, the method never declares a winner with desired certainty, which is in line with what we should expect.

We now focus on the second specific case of  $C = 5, K = 50, n = 50000$ . For different SEOs, the values of  $\mathbf{p}$  are generated randomly from a Dirichlet distribution, and we evaluate how well our method can predict the cumulative proportions of counts for the five categories. Experiments are repeated many times and we compute the overall root mean squared error (RMSE) for every SEO,

Table 2: Accuracy (in %) of predicting the candidate with maximum vote count for different SEOs and for different values of  $\delta = p_1 - p_2$ . All results correspond to the case of  $n = 5000$ ,  $C = 3$ ,  $K = 25$ . Numbers inside the parentheses indicate what percentage of data are used on an average before making a call.

SEO	$\delta$	Correct (data used)	Incorrect (data used)	No call
DGP1	0.01	96% (14%)	4% (13%)	
	0.05	100% (12%)		
	0.10	100% (12%)		
	0.25	100% (12%)		
SEO2	0.01	64.5% (22%)	30.5% (14%)	5%
	0.05	76% (22%)	22.5% (14%)	1.5%
	0.10	89.5% (16%)	10% (14%)	0.5%
	0.25	100% (12%)		
SEO3	0.01	48% (26%)	34% (14%)	18%
	0.05	63.5% (24%)	21% (13%)	15.5%
	0.10	68% (22%)	16.5% (15%)	15.5%
	0.25	74.5% (16%)	8.5% (13%)	17%

Table 3: Root mean squared error (RMSE, in %) in estimating the overall cumulative proportions of counts for all categories.

Data used	SEO1	SEO2	SEO3
5 rounds	0.06%	1.95%	4.38%
15 rounds	< 0.01%	0.95%	2.21%
25 rounds	< 0.01%	0.63%	1.46%
35 rounds	< 0.01%	0.41%	0.92%
45 rounds	< 0.01%	0.22%	0.51%

based on the predictions made at different stages. Refer to Table 3 for these results. Once again, for SEO1, the method achieves great accuracy very early. With only 15 rounds of data (approximately 30% of the total observations), the predictions are precise. At a similar situation, for SEO2, the RMSE is less than 1%, and it decreases steadily to fall below 0.5% by round 35. Finally, for the third SEO, RMSE of less than 1% is recorded after 35 rounds as well. These results show that our method can accurately predict the overall cumulative proportions of counts with about 35 rounds of data. Translating this to the sales forecasting problems, we hypothesize that the proposed method will be able to predict the annual market shares of different categories correctly at least 3 to 4 months before the year-end.

## 4 Application on the data from Bihar election

We examine the effectiveness of our proposed methodology in the election calling context, with the real data from the Bihar Legislative Assembly Election held in 2020, across 243 constituencies (seats). Indian elections have multi-party system, with a few coalition of parties formed before (sometimes after) the election, playing the major role in most contests. While these coalitions are at times temporary in nature and there is no legal binding or legitimacy, typically the alliance winning the majority (at least 50%) of the constituencies comes to power. In the Bihar 2020 election, the

contests in most constituencies were largely limited to two such alliances – the National Democratic Alliance (NDA), and the *Mahagathbandhan* (MGB). Although there were a couple of other alliances, they did not impact the election outcomes and for the purpose of this analysis, we consider them in “others” category. For each constituency, the counting of votes took place in several rounds, and we want to examine how these round-wise numbers can be useful in predicting the final winner ahead of time.

The result of the election is downloaded from the official website of [Election Commission of India \[2020\]](#). A brief summary of the data is provided in Table 4. All summary statistics in this table are calculated over the 243 constituencies in Bihar (our analysis ignores the postal votes, which are few in numbers and do not affect the results). Note that the number of votes cast in the constituencies vary between 119159 and 225767, with the average being 172480. The votes are counted in 25 to 51 rounds across the state, and the mode is found out to be 32. While in most rounds about little over 5000 votes are counted, in some cases, especially the last few rounds, show fairly less numbers. The last three rows in the table provide the minimum, maximum and average number of votes counted per round for all the constituencies. We also point out that the final margin between winner and runner-up parties in these constituencies show a wide range. The lowest is recorded in Hilsa (13 votes) whereas the maximum is recorded in Balrampur (53078 votes).

Table 4: Summary of the Bihar election data. All summary statistics are calculated over the 243 constituencies.

	Minimum	Maximum	Average	Median
Total votes cast	119159	225767	172480	172322
Number of rounds for counting	25	51	32.47	32
Final margin	13	53078	16787	13913
Minimum votes counted per round	114	5173	1618	1029
Maximum votes counted per round	4426	8756	6745	6728
Average votes counted per round	3070	6505	5345	5381

As mentioned above, in the following analysis, we focus on the two dominant categories (essentially the two main alliances in every constituency) and the rest of the alliances or the parties are collated into the third category. As our main objective is to predict the winner and the margin of victory, it is sensible to consider this three-category-setup. We also make the logical assumption that each vote is cast independently, thereby ensuring that in each round of counting, we have an independent multinomial distribution. However, there is no information on how the counting happens sequentially in each round, and therefore we assume the individual probabilities of the three different categories in different counting rounds to be random variables satisfying Assumption 1. Consequently, it is an exciting application of the proposed modeling framework.

Recall the hierarchical prior distributions for the parameters in the model (eq. (2.11)). For  $\alpha$ , we use the past election’s data and use the proportion of votes received by the corresponding alliances (scaled, if needed, to have  $\sum \alpha_i = 1$ ). For  $\Psi$  and  $\Psi_p$ , we use identity matrices of appropriate order. Both  $\nu$  and  $\nu_p$  are taken to be 5. We point out that the results are not too sensitive to these choices. Next, using these priors, we implement our method and estimate the probability of winning for every alliance after the counting of votes in each round. Corresponding margin of victory is also estimated in the process. Based on these values, we propose the following rules to call a race. First, we allow at least 50% votes to be counted in order to make a prediction. It is otherwise considered to be *too early to call*. On the other hand, akin to the procedure laid out by [Lapinski et al. \[2020\]](#), our decision is to call the race in a particular constituency when we are at least 99.5% confident of

the winner and when the predicted winning margin is at least 5% of the remaining votes. Unless these conditions are met, it is termed as *too close to call*.

We start with a summary of the results obtained after fitting the model to the election data. Complete results are provided in Table A4 in Appendix A. In 227 out of the 243 constituencies (approximately 93.4%), the proposed method calls the race correctly. Our method considers that it was always *too close to call* for Bakhri and Barbigaha where the eventual win margin was 439 and 238 respectively. For 14 other constituencies (approximately 5.7%), the method calls the race incorrectly, as the winning candidate was significantly behind at the time of calling in those constituencies.

We next take a detailed look at the prediction patterns for all constituencies, in different aspects, through Figure 2. In both panels of the Figure we color the correct predictions by green, the incorrect predictions by red, and no calls by blue. In the left panel of the Figure we plot the remaining vote percentages at the time of calling along the color gradient. In the right panel we plot the final margins along the color gradient. It is interesting to observe that in the right panel all red points, which correspond to incorrect calls, appear with lighter shades. Thus, our method performs perfectly for all constituencies but one where the final margins of victories are above 10000. From the left panel it is also evident that the proposed method calls the races very early for all such constituencies. Only exception to this is the constituency Baisi, which is discussed in more detail below. We can also observe that seats painted in lighter shades in left panel are also in lighter shades on right panel, implying that the closely contested constituencies require more votes to be counted before the race can be called.

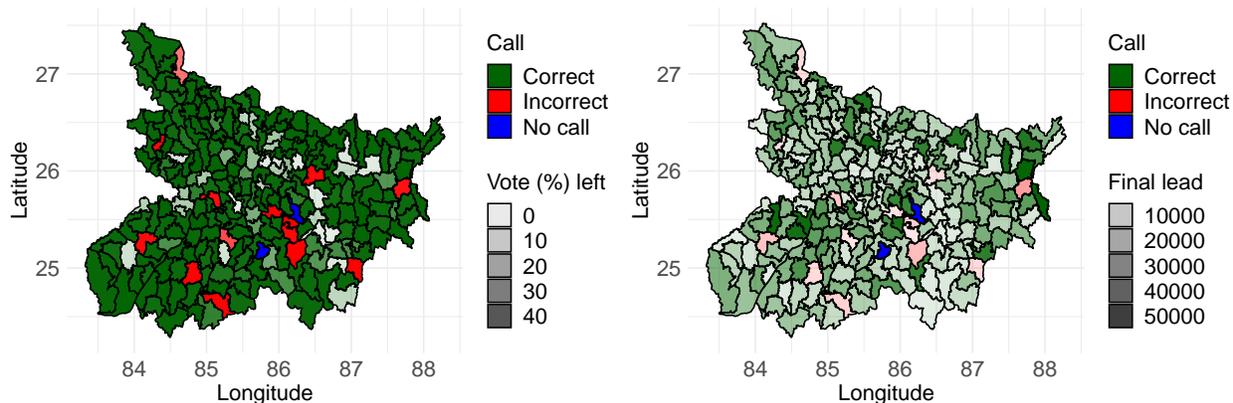


Figure 2: Forecasting accuracy for different constituencies in the Bihar election data. Different colours indicate whether it is a correct call or not. Left panel shows how quickly a call is made, as a darker color indicates that the call is made with higher proportion of votes remaining to be counted. In the right panel, a lighter shade indicates smaller final margin, i.e. a closely contested election.

To further explore the above point, the accuracy and the average percentage of votes counted before making a call, corresponding to the final margins of victories, are displayed in Table 5. There are 52 constituencies which observed very closely fought election (final margin is less than 5000 votes). In 41 out of them, our method correctly predicts the winner and in 2 of them, it cannot make a call. On average, around 70% votes are counted for all these constituencies before we can make a call. This, interestingly, drops drastically for the other constituencies. On average, only about 60% votes are counted before we make a call in the constituencies where eventual margin is between 5000 and 10000. In 4 out of 34 such constituencies though, our method predicts inaccurately. For

the constituencies with higher eventual margins (greater than 10000), the prediction turns out to be correct in 156 out of 157 constituencies. In these cases, around only about 50 to 55 percentage of votes are counted on average before making a call. Finally, the last row of the table corroborates the earlier observation that the size of the constituency does not have considerable effect on the prediction accuracy.

Table 5: Summary of the prediction accuracy in terms of calling the race in different constituencies, according to the final margin of victory.

Final margin	< 2000	2000–5000	5000–10000	10000–20000	> 20000	Total
Constituencies	23	29	34	74	83	243
Correct call	16	25	30	73	83	227
Incorrect call	5	4	4	1	0	14
Too close to call	2	0	0	0	0	2
Average counting	68.0%	74.1%	60.5%	55.2%	51.9%	58.3%
Average votes	172136	170946	171107	171636	174427	172480

Not only the winner of an election, but the proposed approach also predicts the final margin of victory, along with a prediction interval, for the winner. In Figure 3, we present the true margin and the predicted win margin for all the 227 constituencies where correct calls are made. Corresponding prediction intervals are also shown in the same plot. It can be observed that the predicted margins closely resemble the true margins.

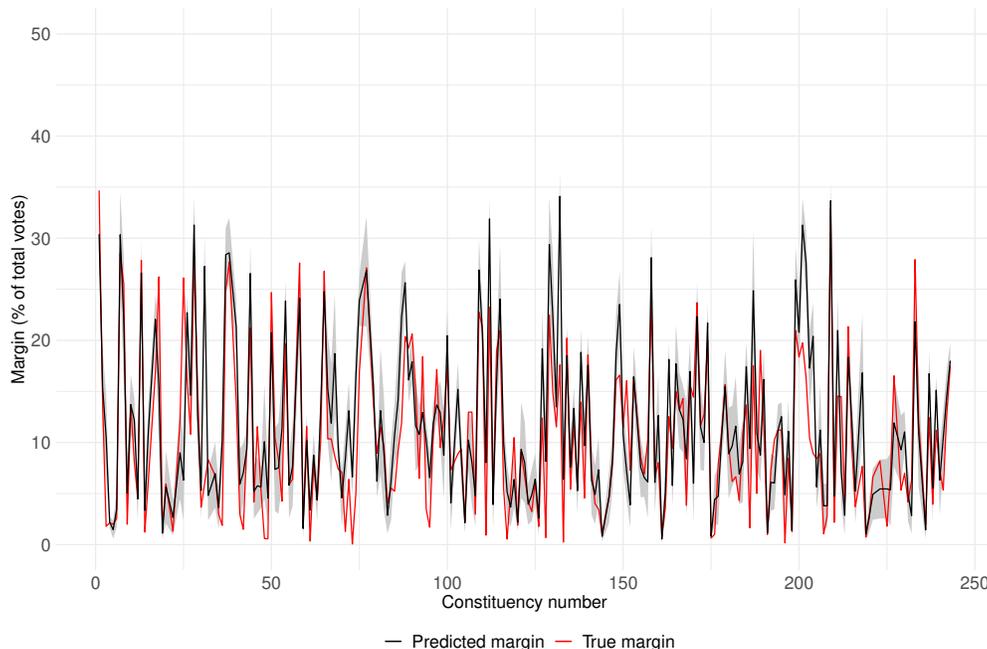


Figure 3: True margin of victory and the predicted margin of victory (both in % of total votes) for all constituencies where the method made a correct call. Prediction interval is displayed in grey.

Overall, it is evident that the method performs really well in terms of calling an election race based on only the counting of votes in different rounds. It is imperative to point out that better accuracy can be achieved with more detailed information about different constituencies. Especially,

information about the sequence in which the counting is carried out, ethnic composition or socio-economic status of people in different areas are necessary to build a more accurate procedure. Our proposed method in this paper works with only the multinomial assumption and achieves more than 93% accuracy in calling the race with the requirement of at least 50% votes being counted. To that end, in order to gain more insight about the accuracy of the method, we rerun the experiment with different values for the minimum requirement of counting. The number of constituencies with correct call, incorrect call and no call are displayed against the minimum requirement in Table 6. We see that the method achieves 90% accuracy (215 out of 243) even when only 30% votes are counted. For minimum counting of 60% or more votes, the results do not improve substantially, except for the fact that no call can be made in more number of constituencies.

Table 6: Prediction accuracy across all 243 constituencies for the proposed method corresponding to minimum % of votes required to be counted before calling a race.

Minimum requirement	Correct call	Incorrect call	Too close to call
10%	182	61	0
20%	199	44	0
30%	215	28	0
40%	221	21	1
50%	227	14	2
60%	234	7	2
70%	235	4	4
80%	235	3	5
90%	235	1	7

Next, we look into the robustness of the method, by focusing on four particular constituencies in more detail. These are Hilsa (minimum final margin), Balrampur (maximum final margin), Baisi (final margin more than 10000, but our proposed methodology makes a wrong call with nearly 55% votes counted) and Barbigha (final margin of 238, the proposed method never calls the race in anyone’s favour). Now, for each constituency, we create synthetic data by randomly switching the order of the individual rounds of counting and implement our method on the resulting data. This exercise helps us to understand whether the predictions would have been different if the data were available in a different order. In particular, it points towards objectively judging to what extent the success or failure of the proposed methodology in this real data can be attributed to luck. Note that the proposed modeling framework assumes independence across different rounds, and hence the method should perform similarly if the votes are actually counted in random order. In Table 7 below, the results of these experiments are presented. We see that for Balrampur, even with randomly permuted rounds, the method calls the race correctly every time. For Baisi, the method yields correct forecast 99.97% of times, although in the original order of the data it makes an incorrect call. Meanwhile, Hilsa and Barbigha observed very closely contested elections. For these two constituencies, in the experiment with permuted data, the method records *too close to call* decisions more commonly, 72.55% and 70.53% respectively. These results suggest that quite possibly the sequence of roundwise counting of votes is not random. Thus, if more information is available about individual rounds from which votes are counted, it would be possible to modify the current method to achieve greater accuracy.

Table 7: Summary of the prediction accuracy in four constituencies, for synthetic data generated through many permutations of individual rounds.

Constituency	Final margin	Original call	Results in permuted data		
			Correct	Incorrect	No call
Hilsa	13	Incorrect	12.57%	14.87%	72.55%
Balrampur	53078	Correct	100%	0%	0%
Baisi	16312	Incorrect	99.97%	0.03%	0%
Barbiga	238	No call	17.19%	12.28%	70.53%

## 5 Application to the USA election with hypothetical sequences

In this section, in an attempt to impress upon the reader that the proposed methodology can be extended to other election forecasting contexts as well, we focus on the statewide results from the 2020 Presidential Election of the USA (data source: [MIT Election Data and Science Lab \[2020\]](#)). It is well known that in American politics, every election relies primarily on a few states, popularly termed as the ‘swing states’, where the verdict can reasonably go to either way. For the other states, the results are guessable with absolute certainty, and are of less interest in any such research. Thus, in this application, our focus is on the eleven swing states of the 2020 Presidential Election – Arizona, Florida, Georgia, Michigan, Minnesota, Nevada, New Hampshire, North Carolina, Pennsylvania, Texas and Wisconsin. Since the final outcome of the election in these cases are extremely likely to be either democratic or republican, with a minimal chance of a third-party-victory, we can consider a three-category-setup. Then, akin to Section 4, it can be easily argued that the setup follows the proposed modeling framework. We also adopt the same setting for prior distributions and the procedure to call the race.

Our objective is to demonstrate that partial data of the counties can provide accurate forecast of the final result in these turbulent states. However, we note that the information of the particular order in which results were declared for different counties are unavailable. That is why we call it a hypothetical application, and we circumvent the issue by taking a permutation-based approach to implement the proposed method. For every state, we randomly assign the orders to the counties and assume that the data are available in that sequence. This permutation procedure is repeated 100 times, and we evaluate the performance of the proposed method at a summary level across these permutations.

Let us start with a brief overview of the number of counties and the number of voters in the swing states (see Table 8). All summary statistics are calculated based on the eleven states. We can see that the number of counties as well as the number of voters vary substantially; although the final margin of victory remains within 7.4% of the total votes. It is in fact less than 1% in three states (Arizona, Georgia and Wisconsin), making them the most closely contested states in this election.

Turn attention to the results of the forecasting procedure on the permuted datasets. In the last three columns of Table 9, we display the proportion of times our method correctly calls the race, finds it too close to call, or incorrectly calls the race based on partial county-level data,

We observe that in New Hampshire, Florida, Nevada and Texas, partial data provide the correct prediction in almost all cases. One may connect this to the fact that the final margin in these states were on the higher side, especially if we look at the values in terms of proportion to the total number of votes. Arizona, on the other hand, was the most closely contested state with a big voter pool. Our approach is still able to predict the correct winner 60% of the times whereas it decides the race

Table 8: Summary of the 2020 US Presidential Election data. All summary statistics are calculated based on the 11 swing states.

	Minimum	Maximum	Average	Median
Number of counties	10	254	84.64	72
Total votes cast	803,833	11,315,056	5,229,904	4,998,482
Final margin (of total votes)	0.25%	7.37%	2.94%	2.40%
Minimum votes per county	66	16515	3136	1817
Maximum votes per county	4426	2,068,144	914,028	755,969
Average votes per county	31427	225,686	85322	66739

Table 9: Performance summary for the application on US Presidential Election data. Total number of votes is given in millions, final margin is given as percentage of total votes, and the last three columns indicate percentage of times the algorithm made a correct call, could not call or made a wrong call.

State	Counties	Total votes	Final margin	Correct	Too close	Incorrect
Arizona	15	3,385,294	10,457	60%	39%	1%
Florida	67	11,067,456	371,686	99%	0%	1%
Georgia	159	4,998,482	12,670	59%	1%	40%
Michigan	83	5,539,302	154,188	72%	0%	28%
Minnesota	87	3,277,171	233,012	82%	0%	18%
Nevada	17	1,404,911	33,706	97%	3%	0%
New Hampshire	10	803,833	59,277	100%	0%	0%
North Carolina	100	5,524,802	74,481	74%	1%	25%
Pennsylvania	67	6,915,283	80,555	54%	4%	42%
Texas	254	11,315,056	631,221	96%	0%	4%
Wisconsin	72	3,297,352	20,608	54%	1%	45%

to be too close to call on 39% occasions. Georgia and Wisconsin are the two other states with the lowest final margins. There, our algorithm makes wrong judgements in more than 40% situations.

In the other four states, namely Michigan, Minnesota, North Carolina and Pennsylvania, the results are not good. Albeit the final margins are considerably large, the final conclusions are wrong in at least 18% of the permuted datasets. This phenomena brings forward an important limitation of the methodology. In our procedure, the key assumption is that the perturbations, albeit random and different in different counties, have the same asymptotic distribution. In such political applications, this assumption may be violated on certain occasions. To demonstrate it empirically, let us look at the margins of differences across the counties in these states (Figure 4).

From the figure, it is evident that in all these states (most prominently in Michigan, Minnesota, Georgia and Wisconsin), the final verdict relied heavily on only a couple of counties. Quite naturally, a partial dataset with the information of those particular counties is extremely likely to provide contradictory results to a partial dataset without those counties. In North Carolina, on the other hand, several counties observed high margin of difference in favor of the democratic party, while the final ruling went the opposite way. These situations point to a potential deviation from the aforementioned assumption, and we hypothesize that in such circumstances, the method can be updated to include more information about individual counties. It calls for an interesting future research direction and we discuss it in more detail in the next section.

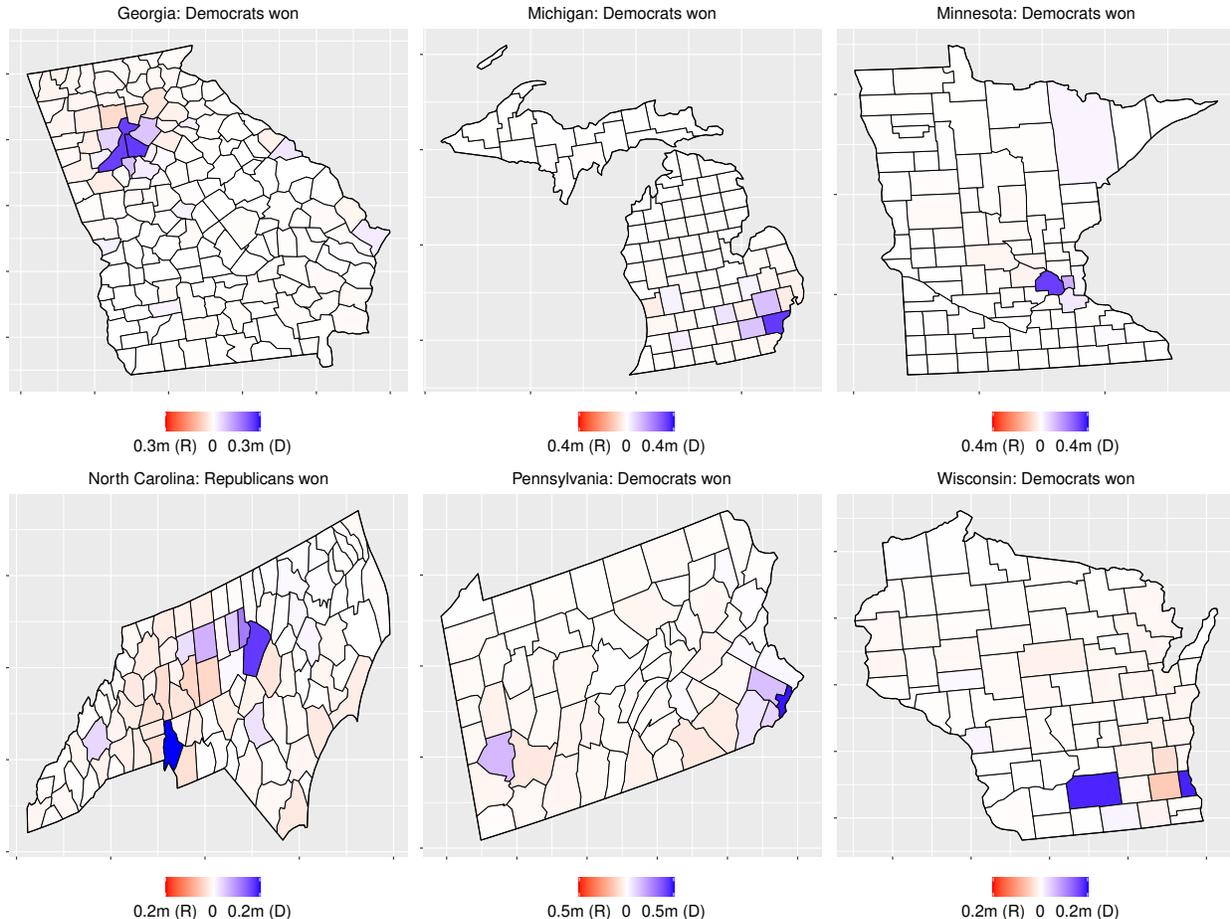


Figure 4: County-wise margin of difference in votes for the two main parties (D and R stand for democratic and republican party respectively) in the six states where the method did not perform well. The unit “m” indicates millions.

## 6 Conclusions

In this work, we consider a hierarchical Bayes model for batches of multinomial data with cell probabilities changing randomly across batches. The goal is to predict various properties for aggregate data correctly and as early as possible. We illustrate the effectiveness of the methodology in poll prediction problems which motivated this study. The performance of the methodology is good, especially considering the limited information. We plan to implement the method in future elections. A potential future direction in this specific context would be to improve the method when more information on the constituencies or the rounds of counting are available. It is likely that additional information on relevant covariates would improve forecast accuracy. It is possible to modify the proposed methodology to incorporate the covariates, but that is a more challenging and interesting problem and requires full treatment. We defer it for a future work.

The proposed model accommodates randomness in the multinomial cell probabilities, depending on the batch. This is pragmatic given the actual data pattern in most practical situations, as otherwise early calls are made which end up being often wrong. This is also intuitively justifiable as, for example, different rounds of votes can have differential probabilities for the candidates, potentially owing to different locations or the changes in people’s behaviour. This paper aims to

tackle these situations appropriately.

It is worth mention that a common practice in Bayesian analysis of sequences of multinomial data is to apply Dirichlet priors for the cell probabilities. This approach requires the iid assumption and hence do not work under the modeling framework of this study. In fact, if we assume iid behaviour for the multinomial data across the batches and use that model for the examples discussed above, then the performances are much worse than what we achieve. Further, we want to point out that even without the variance stabilizing transformation, the proposed Bayesian method in conjunction with the results from Proposition 1 can be used in similar problems. The prediction accuracy of that model is comparable to our approach. We employ the transformation for superior prediction performance, albeit marginally in some cases, across all scenarios.

Finally, note that in some contexts, it may be unrealistic to assume that values of the future  $n_j$ 's are known. In that case, the decision can be taken by anticipating them to be at the level of the average of  $n_j$ 's observed so far. It is intuitively clear from Section 2.4, as one can argue that the effectiveness of the methodology will not alter to any appreciable extent as long as the sample sizes are large. It can also be backed up by relevant simulation studies.

The methodology proposed in this work can be applied in other domain as well, most notably in sales forecasting. Retail giants and companies track the sales for different products in regular basis. This is useful from various perspectives. One of the key aspects is to project or assess the overall (annual) sales pattern from the data collated at monthly, weekly or even daily level. It is of great value to a company to infer about its eventual market share for the period (year) as early as possible, hence to validate if its (possibly new) marketing plans are as effective as targeted. The retailer would find these inferences useful from the perspective of managing their supply chain, as well as in their own marketing. Other possible applications of the proposed methodology (possibly with necessary adjustments) include analyzing hospital-visits data, in-game sports prediction etc. Some of these applications can make use of additional covariate information; we propose to take that up in a follow-up article.

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## 7 Proofs

**Proposition 1.** Following eq. (2.2),  $\mathbb{E}(X_{lj}|\varepsilon_{lj}) = n_j(p_l + \varepsilon_{lj})$ . Then, using Assumption 1 and applying the relationship between total and conditional expectation,

$$\mathbb{E}(X_{lj}) = \mathbb{E}[\mathbb{E}(X_{lj}|\varepsilon_{lj})] = n_j p_l. \quad (7.1)$$

We also know that  $\text{Var}(X_{lj}|\varepsilon_{lj}) = n_j(p_l + \varepsilon_{lj})(1 - p_l - \varepsilon_{lj})$ . Then, we can use the relationship  $\text{Var}(X_{lj}) = \text{Var}(\mathbb{E}(X_{lj}|\varepsilon_j)) + \mathbb{E}(\text{Var}(X_{lj}|\varepsilon_j))$  to get the following:

$$\text{Var}(X_{lj}) = \text{Var}(n_j(p_l + \varepsilon_{lj})) + \mathbb{E}[n_j(p_l + \varepsilon_{lj})(1 - p_l - \varepsilon_{lj})] = n_j p_l(1 - p_l) + n_j(n_j - 1)\text{Var}(\varepsilon_{lj}). \quad (7.2)$$

Finally, note that the covariance of  $X_{1j}$  and  $X_{2j}$ , conditional on  $\varepsilon_j$ , is equal to  $-n_j(p_1 + \varepsilon_{1j})(p_2 + \varepsilon_{2j})$ . One can use this and apply similar technique as before to complete the proof of part (a).

Part (b) follows from (a) using the fact that the multinomial data in different batches are independent of each other.

To prove part (c), we use the multinomial central limit theorem which states that if  $(D_1, D_2, \dots, D_C) \sim \text{Multinomial}(n, \pi_1, \pi_2, \dots, \pi_C)$ , then

$$\sqrt{n} \left( \begin{pmatrix} D_1/n \\ D_2/n \\ \vdots \\ D_{C-1}/n \end{pmatrix} - \begin{pmatrix} \pi_1 \\ \pi_2 \\ \vdots \\ \pi_{C-1} \end{pmatrix} \right) \xrightarrow{\mathcal{L}} \mathcal{N}_2 \left( 0, \begin{bmatrix} \pi_1(1 - \pi_1) & -\pi_1\pi_2 & \dots & -\pi_1\pi_{C-1} \\ -\pi_1\pi_2 & \pi_2(1 - \pi_2) & \dots & -\pi_2\pi_{C-1} \\ \vdots & \vdots & \ddots & \vdots \\ -\pi_1\pi_{C-1} & -\pi_2\pi_{C-1} & \dots & \pi_{C-1}(1 - \pi_{C-1}) \end{bmatrix} \right). \quad (7.3)$$

Let us define  $V_{\varepsilon_j}$  to be the covariance matrix of  $(\hat{p}_{1j}, \hat{p}_{2j}, \dots, \hat{p}_{(C-1)j})$ , conditional on  $\varepsilon_j$ . Then, for the  $(c, c')$ <sup>th</sup> element of  $V_{\varepsilon_j}$ , we have

$$(V_{\varepsilon_j})_{c,c'} = \begin{cases} (p_c + \varepsilon_{cj})(1 - p_c - \varepsilon_{cj}) & \text{for } c = c', \\ -(p_c + \varepsilon_{cj})(p_{c'} + \varepsilon_{c'j}) & \text{for } c \neq c'. \end{cases}$$

Now, under Assumption 1 and from eq. (7.3), as  $n_j \rightarrow \infty$ ,

$$\sqrt{n_j} V_{\varepsilon_j}^{-1/2} (\hat{\mathbf{p}}_j - \tilde{\mathbf{p}} - \varepsilon_j) | \varepsilon_j \xrightarrow{\mathcal{L}} \mathcal{N}_{C-1}(\mathbf{0}, \mathbf{I}). \quad (7.4)$$

Using the assumptions in part (c) of Proposition 1, as  $n_j \rightarrow \infty$ ,  $V_{\varepsilon_j}$  converges in probability to  $\Xi_p$ . Hence, by an application of Slutsky's Theorem, eq. (7.4) is equivalent to:

$$\sqrt{n_j} \Xi_p^{-1/2} (\hat{\mathbf{p}}_j - \tilde{\mathbf{p}} - \varepsilon_j) | \varepsilon_j \xrightarrow{\mathcal{L}} \mathcal{N}_{C-1}(\mathbf{0}, \mathbf{I}). \quad (7.5)$$

Let  $f_{\varepsilon}(\cdot)$  denote the density function of  $\sqrt{n_j} \varepsilon_j$ ,  $\phi(\cdot, \Xi_{\varepsilon})$  be the density function of  $\mathcal{N}_{C-1}(\mathbf{0}, \Xi_{\varepsilon})$ , and  $\Phi(\cdot, \Xi_{\varepsilon})$  be the distribution function of  $\mathcal{N}_{C-1}(\mathbf{0}, \Xi_{\varepsilon})$ . Note that the previous convergence in eq. (7.5) is conditional on  $\varepsilon_j$ , as  $n_j \rightarrow \infty$ . To find the unconditional distribution of  $\hat{\mathbf{p}}_j - \tilde{\mathbf{p}}$ , we look at the following, for  $z \in \mathbb{R}^{C-1}$ :

$$\mathbb{P}(\sqrt{n_j}(\hat{\mathbf{p}}_j - \tilde{\mathbf{p}}) \leq z) = \int_{\mathbb{R}^{C-1}} \mathbb{P}(\sqrt{n_j}(\hat{\mathbf{p}}_j - \tilde{\mathbf{p}} - \varepsilon_j) \leq z - y | \varepsilon_j) f_{\varepsilon}(y) dy. \quad (7.6)$$

From the assumptions in part (c) of Proposition 1,  $\sqrt{n_j} \varepsilon_j$  converges in distribution to  $\mathcal{N}_{C-1}(\mathbf{0}, \Xi_{\varepsilon})$ . Then,

$$\lim_{n_j \rightarrow \infty} \mathbb{P}(\sqrt{n_j}(\hat{\mathbf{p}}_j - \tilde{\mathbf{p}} - \varepsilon_j) \leq z - y | \varepsilon_j) f_{\varepsilon}(y) = \Phi(z - y, \Xi_p) \phi(y, \Xi_{\varepsilon}). \quad (7.7)$$

A simple application of dominated convergence theorem then implies, as  $n_j \rightarrow \infty$ ,

$$\mathbb{P}(\sqrt{n_j}(\hat{\mathbf{p}}_j - \tilde{\mathbf{p}}) \leq z) \rightarrow \int_{\mathbb{R}^{C-1}} \Phi(z - y, \Xi_p) \phi(y, \Xi_{\varepsilon}) dy. \quad (7.8)$$

Note that the term on the right hand side of eq. (7.8) is the distribution function of  $\mathcal{N}_{C-1}(\mathbf{0}, \Xi_p + \Xi_\epsilon)$  at the point  $z \in \mathbb{R}^{C-1}$ . Hence we arrive at the required result, eq. (2.5).  $\square$

**Theorem 1.** Recall that the sigma field generated by the collection of the data  $\{\mathbf{L}_1, \mathbf{L}_2, \dots, \mathbf{L}_K\}$  is  $\mathcal{F}_K$ , and the posterior means of  $\boldsymbol{\mu}$  and  $\Sigma$ , conditional on  $\mathcal{F}_K$ , are  $\boldsymbol{\mu}_{PM}$  and  $\Sigma_{PM}$  respectively. Then, eq. (2.12) and eq. (2.13) can be equivalently stated as,

$$\lim_{K \rightarrow \infty} \mathbb{E} [\Pi_K (\|\boldsymbol{\mu}_{PM} - \boldsymbol{\mu}_0\| > \epsilon \mid \mathcal{F}_K)] = 0, \quad (7.9)$$

$$\lim_{K \rightarrow \infty} \mathbb{E} [\Pi_K (\|\Sigma_{PM} - \Sigma_0\| > \epsilon \mid \mathcal{F}_K)] = 0. \quad (7.10)$$

We begin with the term

$$\Pi_K (\|\boldsymbol{\mu}_{PM} - \boldsymbol{\mu}_0\| > \epsilon \mid \mathcal{F}_K). \quad (7.11)$$

We know that  $\boldsymbol{\mu}_{PM} = \mathbb{E} \left( [\Sigma_p^{-1} + N\Sigma^{-1}]^{-1} \left[ \Sigma_p^{-1} \boldsymbol{\alpha} + \Sigma^{-1} \left( \sum_{i=1}^K n_i \mathbf{L}_i \right) \right] \mid \mathcal{F}_K \right)$ , where the expectation is taken with respect to  $\Sigma$ . Since eq. (7.11) is either 1 or 0 depending on whether  $\|\boldsymbol{\mu}_{PM} - \boldsymbol{\mu}_0\| > \epsilon$  or not, on taking expectation over  $\mathcal{F}_K$ , we obtain

$$\mathbb{P} (\|\boldsymbol{\mu}_{PM} - \boldsymbol{\mu}_0\| > \epsilon \mid \mathcal{F}_K). \quad (7.12)$$

To show that eq. (7.12) goes to 0 as  $K \rightarrow \infty$ , we do the following:

$$\begin{aligned} \|\boldsymbol{\mu}_{PM} - \boldsymbol{\mu}_0\| &= \left\| \mathbb{E} \left( [\Sigma_p^{-1} + N\Sigma^{-1}]^{-1} \left[ \Sigma_p^{-1} \boldsymbol{\alpha} + \Sigma^{-1} \left( \sum_{i=1}^K n_i \mathbf{L}_i \right) \right] \mid \mathcal{F}_K \right) - \boldsymbol{\mu}_0 \right\| \\ &\leq \mathbb{E} \left( \left\| [\Sigma_p^{-1} + N\Sigma^{-1}]^{-1} \left[ \Sigma_p^{-1} (\boldsymbol{\alpha} - \boldsymbol{\mu}_0) + \Sigma^{-1} \left\{ \sum_{i=1}^K n_i (\mathbf{L}_i - \boldsymbol{\mu}_0) \right\} \right] \right\| \mid \mathcal{F}_K \right) \\ &\leq \mathbb{E} \left( \left\| [\Sigma_p^{-1} + N\Sigma^{-1}]^{-1} \Sigma_p^{-1} (\boldsymbol{\alpha} - \boldsymbol{\mu}_0) \right\| \mid \mathcal{F}_K \right) \\ &\quad + \mathbb{E} \left( \left\| [\Sigma_p^{-1} + N\Sigma^{-1}]^{-1} \Sigma^{-1} \left\{ \sum_{i=1}^K n_i (\mathbf{L}_i - \boldsymbol{\mu}_0) \right\} \right\| \mid \mathcal{F}_K \right). \end{aligned} \quad (7.13)$$

Applying the identity  $(A+B)^{-1} = A^{-1} - A^{-1}(A^{-1} + B^{-1})^{-1}A^{-1}$  on the first term of the above,

$$\left\| [\Sigma_p^{-1} + N\Sigma^{-1}]^{-1} \Sigma_p^{-1} (\boldsymbol{\alpha} - \boldsymbol{\mu}_0) \right\| = \left\| \left[ \mathbf{I} - \Sigma_p \left( \Sigma_p + \frac{\Sigma}{N} \right)^{-1} \right] (\boldsymbol{\alpha} - \boldsymbol{\mu}_0) \right\|, \quad (7.14)$$

which goes to 0 a.e., using the properties of inverse Wishart distribution and the assumptions on the prior parameters we have. Consequently,  $\mathbb{E} \left\| [\Sigma_p^{-1} + N\Sigma^{-1}]^{-1} \Sigma_p^{-1} (\boldsymbol{\alpha} - \boldsymbol{\mu}_0) \right\|$  goes to 0 and hence, we can argue that

$$\mathbb{P} \left( \mathbb{E} \left\| [\Sigma_p^{-1} + N\Sigma^{-1}]^{-1} \Sigma_p^{-1} (\boldsymbol{\alpha} - \boldsymbol{\mu}_0) \right\| > \frac{\epsilon}{2} \right) \rightarrow 0. \quad (7.15)$$

For the second term in eq. (7.13), first note that  $\sum_{i=1}^K n_i (\mathbf{L}_i - \boldsymbol{\mu}_0)$  is equivalent, in distributional sense, to  $\sqrt{N}Z$ , where  $Z \sim \mathcal{N}(\mathbf{0}, \Sigma_0)$ . On the other hand,

$$\left\| \mathbb{E} \left( [\Sigma_p^{-1} + N\Sigma^{-1}]^{-1} \Sigma^{-1} \right) \right\| \leq \left\| \Sigma_p \mathbb{E} \left[ (\Sigma + N\Sigma_p)^{-1} \right] \right\| \leq C_K \|\Sigma_p\|, \quad (7.16)$$

where  $C_K = O(1/N)$ . Using the above results, we can write

$$\mathbb{P} \left[ \left\| \mathbb{E} \left( [\Sigma_p^{-1} + N\Sigma^{-1}]^{-1} \Sigma^{-1} \sqrt{N} Z \mid \mathcal{F}_K \right) \right\| > \frac{\epsilon}{2} \right] \leq \mathbb{P} \left[ \left\| C_K \sqrt{N} Z \right\| > \frac{\epsilon}{4 \|\Sigma_p\|} \right]. \quad (7.17)$$

In light of the fact that  $\|\Sigma_p\|$  is finite and that  $C_K = O(1/N)$ , it is easy to see that the above probability goes to 0. This, along with eq. (7.15), completes the proof that eq. (7.12) goes to 0 as  $K \rightarrow \infty$ .

In order to prove eq. (7.10), we follow a similar idea as above. Note that  $\Psi$  is a constant positive definite matrix and  $\nu$  is a finite constant. Following eq. (2.17), for large  $K$ , straightforward calculations yield the following.

$$\|\Sigma_{PM} - \Sigma_0\| = \frac{1}{K} \left\| \mathbb{E} \left[ \sum_{i=1}^K n_i (\mathbf{L}_i - \boldsymbol{\mu})(\mathbf{L}_i - \boldsymbol{\mu})^T - \Sigma_0 \mid \mathcal{F}_k \right] \right\|. \quad (7.18)$$

Further, using the true value  $\boldsymbol{\mu}_0$  inside the term on the right hand side above, we get

$$\begin{aligned} \sum_{i=1}^K n_i (\mathbf{L}_i - \boldsymbol{\mu})(\mathbf{L}_i - \boldsymbol{\mu})^T - \Sigma_0 &= \sum_{i=1}^K n_i (\mathbf{L}_i - \boldsymbol{\mu}_0)(\mathbf{L}_i - \boldsymbol{\mu}_0)^T - \Sigma_0 + N(\boldsymbol{\mu} - \boldsymbol{\mu}_0)(\boldsymbol{\mu} - \boldsymbol{\mu}_0)^T \\ &\quad - 2 \sum_{i=1}^K n_i (\mathbf{L}_i - \boldsymbol{\mu}_0)(\boldsymbol{\mu} - \boldsymbol{\mu}_0)^T, \end{aligned} \quad (7.19)$$

and subsequently,

$$\begin{aligned} \mathbb{P} (\|\Sigma_{PM} - \Sigma_0\| > \epsilon \mid \mathcal{F}_k) &\leq \mathbb{P} \left( \left\| \frac{1}{K} \sum_{i=1}^K n_i (\mathbf{L}_i - \boldsymbol{\mu}_0)(\mathbf{L}_i - \boldsymbol{\mu}_0)^T - \Sigma_0 \right\| > \frac{\epsilon}{3} \right) \\ &\quad + \mathbb{P} \left( \left\| \frac{1}{K} \mathbb{E} [N(\boldsymbol{\mu} - \boldsymbol{\mu}_0)(\boldsymbol{\mu} - \boldsymbol{\mu}_0)^T \mid \mathcal{F}_k] \right\| > \frac{\epsilon}{3} \right) \\ &\quad + \mathbb{P} \left( \frac{2}{K} \left\| \mathbb{E} \left[ \sum_{i=1}^K n_i (\mathbf{L}_i - \boldsymbol{\mu}_0)(\boldsymbol{\mu} - \boldsymbol{\mu}_0)^T \mid \mathcal{F}_k \right] \right\| > \frac{\epsilon}{3} \right). \end{aligned} \quad (7.20)$$

Using similar arguments as in eq. (7.16), we can show that  $\mathbb{E}[N(\boldsymbol{\mu} - \boldsymbol{\mu}_0)(\boldsymbol{\mu} - \boldsymbol{\mu}_0)^T \mid \mathcal{F}_k]$  is bounded and therefore, the second term goes to 0. Next, taking cue from the previous part regarding the consistency of  $\boldsymbol{\mu}$ , it is also easy to see that for large  $K$ ,  $\mathbb{E}[n_i (\mathbf{L}_i - \boldsymbol{\mu}_0)(\boldsymbol{\mu} - \boldsymbol{\mu}_0)^T \mid \mathcal{F}_k]$  is small enough. Hence, the third term in eq. (7.20) also goes to 0. Finally, for the first term in that inequality, observe that  $\sqrt{n_i}(\mathbf{L}_i - \boldsymbol{\mu}_0) \sim \mathcal{N}_{C-1}(0, \Sigma_0)$  for all  $i$ . Thus,  $W = \sum_{i=1}^K n_i (\mathbf{L}_i - \boldsymbol{\mu}_0)(\mathbf{L}_i - \boldsymbol{\mu}_0)^T \sim \text{Wishart}(\Sigma_0, K)$ , and the law of large numbers implies that  $W/K - \Sigma_0 \rightarrow 0$  in probability. Consequently, we get that  $\mathbb{P} (\|\Sigma_{PM} - \Sigma_0\| > \epsilon) \rightarrow 0$  and that completes the proof.  $\square$

## Appendix

### A Additional tables

Table A1: Accuracy (in %) of predicting the category with maximum count (corresponding final margins, averaged over all repetitions, are given in parentheses) for different values of  $C$ ,  $K$ ,  $n$ , when data are generated from SEO1.

$C$	$K$	$n$	Correct (average margin)	Incorrect (average margin)	No call (average margin)
3	25	100	95.8% (1158)	0.4% (91)	3.8% (120)
3	25	1000	99.2% (11894)	0.2% (255)	0.6% (399)
3	25	5000	98.62% (54870)	0.92% (454)	0.46% (1108)
3	25	50000	100% (549287)		
3	50	100	95.75% (2384)	0.75% (104)	3.5% (125)
3	50	1000	99% (23425)		1% (346)
3	50	5000	99.5% (114905)		0.5% (1581)
3	50	50000	100% (1139338)		
5	25	100	93.75% (744)	2.25% (27)	4% (81)
5	25	1000	98.25% (7395)	0.25% (38)	1.5% (368)
5	25	5000	98.75% (38799)	0.25% (330)	1% (788)
5	25	50000	99.75% (373575)		0.25% (2292)
5	50	100	94.5% (1584)	1.5% (73)	4% (168)
5	50	1000	98.25% (14388)	0.5% (516)	1.25% (378)
5	50	5000	99.5% (71562)	0.25% (1461)	0.25% (858)
5	50	50000	100% (730428)		

Table A2: Accuracy (in %) of predicting the category with maximum count (corresponding final margins, averaged over all repetitions, are given in parentheses) for different values of  $C, K, n$ , when data are generated from SEO2.

$C$	$K$	$n$	Correct (average margin)	Incorrect (average margin)	No call (average margin)
3	25	100	72.6% (626)	8% (204)	19.4% (370)
3	25	1000	89.6% (9374)	2.8% (2174)	7.6% (2866)
3	25	5000	89.6% (54572)	3.2% (5723)	7.2% (8492)
3	25	50000	98.2% (551259)	1% (20745)	0.8% (40335)
3	50	100	77.5% (1227)	6.75% (355)	15.75% (480)
3	50	1000	88% (18912)	4% (3189)	8% (3772)
3	50	5000	93.25% (100611)	3.75% (7688)	3% (15393)
3	50	50000	96.75% (1098195)	1% (41312)	2.25% (96337)
5	25	100	62.75% (298)	13.75% (146)	23.5% (153)
5	25	1000	77.5% (5681)	8.75% (1067)	13.75% (1758)
5	25	5000	85.75% (33422)	5.5% (4861)	8.75% (7226)
5	25	50000	93% (371349)	2% (32987)	5% (50911)
5	50	100	70% (605)	8.25% (205)	21.75% (314)
5	50	1000	84.75% (10986)	4.75% (1286)	10.5% (2891)
5	50	5000	91.75% (67476)	3.25% (6624)	5% (12241)
5	50	50000	95.5% (720905)	2.25% (71467)	2.25% (77728)

Table A3: Accuracy (in %) of predicting the category with maximum count (corresponding final margins, averaged over all repetitions, are given in parentheses) for different values of  $C, K, n$ , when data are generated from SEO3.

$C$	$K$	$n$	Correct (average margin)	Incorrect (average margin)	No call (average margin)
3	25	100	43% (230)	23% (123)	34% (163)
3	25	1000	49.8% (3558)	20.2% (1948)	30% (2425)
3	25	5000	58.2% (22490)	15% (10192)	26.8% (14920)
3	25	50000	73.8% (323688)	8.8% (100474)	17.4% (187450)
3	50	100	43.2% (414)	18% (258)	38.8% (297)
3	50	1000	56.2% (6696)	13.4% (3131)	30.4% (4367)
3	50	5000	64.6% (45409)	7.2% (16465)	28.2% (27686)
3	50	50000	79.4% (700278)	5.2% (141302)	15.4% (270719)
5	25	100	40% (91)	26% (64)	34% (47)
5	25	1000	42.6% (1323)	26.2% (772)	31.2% (755)
5	25	5000	51% (8658)	24% (5588)	25% (5423)
5	25	50000	65.6% (154765)	10.6% (80310)	23.8% (81664)
5	50	100	42.4% (154)	22% (106)	35.6% (110)
5	50	1000	45% (2439)	23.2% (1423)	31.8% (1536)
5	50	5000	55.8% (18972)	14.2% (9671)	30% (10822)
5	50	50000	70% (305700)	8.8% (79016)	21.2% (184868)

Table A4: Detailed results from all constituencies in Bihar.

Constituency	True data		At the time of calling the race				Decision
	Rounds	Final margin	Round	Votes left (%)	Lead	Predicted margin	
Agiaon	28	48165	14	49.1	21498	42268	Correct
Alamnagar	37	29095	19	49.0	16498	32418	Correct
Alauli	25	2564	13	47.7	7899	14798	Correct
Alinagar	29	3370	28	0.8	3439	3450	Correct
Amarpur	31	3242	30	2.5	2341	2370	Correct
Amnour	29	3824	24	12.8	4556	5177	Correct
Amour	34	52296	17	49.3	28624	55829	Correct
Araria	35	47828	17	48.0	21851	41762	Correct
Arrah	36	3105	29	13.9	6758	7970	Correct
Arwal	30	19651	13	49.1	10167	19375	Correct
Asthawan	31	11530	16	46.9	9270	17362	Correct
Atri	34	7578	17	46.5	3218	5957	Correct
Aurai	33	48008	16	47.4	24134	45972	Correct
Aurangabad	34	2063	15	49.7	2754	5676	Correct
Babubarhi	32	12022	16	49.3	10414	20871	Correct
Bachhwara	32	737	15	49.7	5631	10832	Incorrect
Bagaha	33	30494	17	47.0	21112	39649	Correct
Bahadurganj	33	44978	17	47.4	14130	26583	Correct
Bahadurpur	34	2815	31	1.4	1965	1981	Correct
Baikunthpur	33	10805	18	43.1	5730	10274	Correct
Baisi	30	16312	15	46.6	6761	12218	Incorrect
Bajpatti	34	2325	29	11.2	4225	4778	Correct
Bakhri	28	439					No call
Bakhtiarpur	30	20694	15	47.9	7992	14667	Correct
Balrampur	35	53078	18	45.6	6706	12218	Correct
Baniapur	34	27219	17	48.7	19610	38052	Correct
Banka	27	17093	14	47.9	12133	23154	Correct
Bankipur	45	38965	22	49.3	22243	43644	Correct
Banmankhi	33	27872	16	47.7	11958	23007	Correct
Barachatti	33	6737	20	39.2	5916	9722	Correct
Barari	28	10847	14	46.7	26361	49359	Correct
Barauli	33	14493	16	47.1	4342	8183	Correct
Barbigha	25	238					No call
Barh	30	10084	21	28.5	7709	10450	Correct
Barhara	32	4849	28	12.1	5190	5932	Correct
Barharia	33	3220	16	50.0	8954	17765	Correct
Baruraj	30	43548	15	46.3	26826	49863	Correct
Bathnaha	32	47136	17	46.5	26047	48536	Correct
Begusarai	38	5392	17	50.0	4814	9165	Incorrect
Belaganj	32	23516	16	47.2	18889	35735	Correct
Beldaur	33	5289	16	48.4	5233	10518	Correct
Belhar	32	2713	17	48.6	6701	13164	Correct
Belsand	28	13685	13	49.2	6577	12898	Correct

continued ...

Table A4: Detailed results from all constituencies in Bihar.

Constituency	True data		At the time of calling the race				Decision
	Rounds	Final margin	Round	Votes left (%)	Lead	Predicted margin	
Benipatti	32	32904	16	49.6	20784	41325	Correct
Benipur	31	6793	25	18.2	6957	8499	Correct
Bettiah	30	18375	17	46.8	4802	9154	Correct
Bhabua	29	9447	15	48.8	4971	9665	Correct
Bhagalpur	36	950	20	40.6	9662	15892	Correct
Bhorey	37	1026	19	48.5	4141	7857	Correct
Bibhutipur	30	40369	14	49.4	17213	33918	Correct
Bihariganj	34	19459	18	45.0	7497	13702	Correct
Biharsharif	41	15233	19	47.5	7362	14029	Correct
Bihpur	30	6348	16	46.6	9088	16985	Correct
Bikram	35	35390	17	49.9	21569	42937	Correct
Bisfi	36	10469	19	45.6	8169	14305	Correct
Bochaha	30	11615	15	48.6	7128	13904	Correct
Bodh Gaya	34	4275	17	47.0	2704	5174	Incorrect
Brahampur	37	50537	18	47.4	23505	44079	Correct
Buxar	30	3351	28	6.8	2508	2680	Correct
Chainpur	36	23650	18	48.2	10926	21057	Correct
Chakai	32	654	17	47.5	3278	6470	Correct
Chanpatia	27	13680	14	47.2	8080	15260	Correct
Chapra	36	7222	30	14.6	6135	7234	Correct
Chenari	33	17489	16	49.6	11612	23038	Correct
Cheria Bariarpur	29	40379	15	47.4	19725	37310	Correct
Chhatapur	32	20858	16	48.6	15892	30405	Correct
Chiraiya	31	17216	15	49.4	10055	20033	Correct
Danapur	38	16005	19	47.9	18866	35408	Correct
Darauli	34	11771	16	49.8	7508	14898	Correct
Daraundha	35	11492	20	41.4	4388	7409	Correct
Darbhang Rural	30	2019	14	49.6	6509	13277	Correct
Darbhang	32	10870	16	47.4	11923	22419	Correct
Dehri	32	81	15	46.7	5497	10371	Correct
Dhaka	34	10396	17	47.5	18059	33826	Correct
Dhamdaha	35	33701	18	48.1	24815	47600	Correct
Dhauraiya	31	2687	15	49.8	9155	18294	Incorrect
Digha	51	46073	25	49.3	23335	45542	Correct
Dinara	32	7896	16	47.9	4136	8036	Incorrect
Dumraon	34	23854	17	47.4	14094	26750	Correct
Ekma	33	13683	17	48.0	5019	9433	Correct
Fatuha	29	19407	15	47.1	11926	21926	Correct
Forbesganj	36	19749	18	49.3	9320	18607	Correct
Gaighat	33	7345	27	17.4	4232	5180	Correct
Garkha	33	9746	19	39.7	6592	10794	Correct
Gaura Bauram	25	7519	13	47.7	8065	15274	Correct
Gaya Town	29	12123	15	48.7	9590	18989	Correct

continued ...

Table A4: Detailed results from all constituencies in Bihar.

Constituency	True data		At the time of calling the race				Decision
	Rounds	Final margin	Round	Votes left (%)	Lead	Predicted margin	
Ghosi	28	17804	14	49.3	17071	33251	Correct
Gobindpur	35	32776	18	48.3	21332	41204	Correct
Goh	32	35377	17	47.5	15573	29875	Correct
Gopalganj	37	36641	18	48.6	16397	31998	Correct
Gopalpur	30	24580	15	49.3	9626	19022	Correct
Goriakothi	35	12345	18	47.3	10776	20645	Correct
Govindganj	28	27924	14	46.7	10488	19611	Correct
Gurua	31	6152	16	47.9	9515	18353	Correct
Hajipur	35	3248	19	47.7	6633	12491	Correct
Harlakhi	32	17815	15	49.6	9927	19681	Correct
Harnaut	33	27050	16	48.4	11151	21626	Correct
Harsidhi	28	16071	14	49.7	11104	21881	Correct
Hasanpur	30	21039	18	39.7	8936	14705	Correct
Hathua	32	30237	16	48.1	18605	35616	Correct
Hayaghat	25	10420	23	6.4	5735	6090	Correct
Hilsa	33	13	20	33.7	4433	6678	Incorrect
Hisua	41	16775	20	48.1	15057	28784	Correct
Imamganj	32	16177	16	48.5	7363	13978	Correct
Islampur	31	3767	29	4.8	3297	3474	Correct
Jagdishpur	32	21492	16	47.8	8949	17206	Correct
Jale	33	21926	17	47.3	6953	13395	Correct
Jamalpur	33	4468	21	36.2	4626	7197	Correct
Jamui	32	41009	17	47.9	25100	48494	Correct
Jehanabad	33	33399	16	47.8	17010	32635	Correct
Jhajha	35	1779	17	47.3	8142	15372	Correct
Jhanjharpur	34	41861	17	49.7	28969	57628	Correct
Jokihat	31	7543	18	39.8	3951	6575	Correct
Kadwa	30	31919	15	47.4	13917	26993	Correct
Kahalgaon	36	42947	18	47.5	26297	49848	Correct
Kalyanpur (A)	34	10329	17	49.0	9814	18960	Correct
Kalyanpur (B)	27	852	13	48.1	4335	8519	Correct
Kanti	33	10254	25	23.2	5588	7257	Correct
Karakat	36	17819	17	48.8	5691	11004	Correct
Kargahar	35	3666	32	5.0	3995	4226	Correct
Kasba	31	17081	21	32.2	11768	17215	Correct
Katihar	29	11183	18	37.8	6557	10741	Correct
Katoria	27	6704	24	8.6	5688	6253	Correct
Keoti	31	5267	26	13.3	6814	7892	Correct
Kesaria	29	9352	14	48.8	5048	9796	Correct
Khagaria	28	2661	24	9.1	3550	3921	Correct
Khajauli	31	23037	16	48.9	18093	35361	Correct
Kishanganj	33	1221	16	48.5	7588	14366	Correct
Kochadhaman	26	36072	14	47.1	25184	46871	Correct

continued ...

Table A4: Detailed results from all constituencies in Bihar.

Constituency	True data		At the time of calling the race				Decision
	Rounds	Final margin	Round	Votes left (%)	Lead	Predicted margin	
Korha	30	29007	15	47.7	23079	43352	Correct
Kuchaikote	36	20753	17	49.4	12435	24464	Correct
Kumhrar	49	26466	26	49.4	25992	51127	Correct
Kurhani	33	480	15	50.0	6379	12687	Correct
Kurtha	26	27542	13	49.6	12732	25261	Correct
Kusheshwarasthan	27	7376	14	47.1	5444	10264	Correct
Kutumba	28	16330	15	46.7	9879	18414	Correct
Lakhisarai	45	10709	22	48.0	5317	10227	Correct
Lalганj	34	26613	17	47.9	18727	36231	Correct
Laukaha	37	9471	18	48.2	10665	20498	Correct
Lauriya	27	29172	14	47.9	14345	27468	Correct
Madhepura	35	15072	16	49.6	6466	12643	Correct
Madhuban	27	6115	18	34.1	4790	7219	Correct
Madhubani	37	6490	18	48.7	7292	13924	Correct
Maharajganj	34	1638	31	3.1	1271	1291	Correct
Mahishi	31	1972	16	48.0	9344	17525	Incorrect
Mahnar	31	7781	16	46.9	4244	7697	Correct
Mahua	30	13687	15	48.6	7468	14255	Correct
Makhdumpur	26	21694	14	46.5	13559	24842	Correct
Maner	35	32919	18	47.6	24582	46654	Correct
Manihari	31	20679	14	49.3	9831	19437	Correct
Manjhi	32	25154	16	49.1	5858	11338	Correct
Marhaura	28	10966	18	34.7	3863	5834	Correct
Masaurhi	37	32161	19	48.7	16543	32147	Correct
Matihani	38	65	19	48.3	5017	9799	Incorrect
Minapur	30	15321	15	48.8	6930	13343	Correct
Mohania	30	11100	15	46.3	5643	10447	Correct
Mohiuddinnagar	29	15195	21	23.6	6799	9229	Correct
Mokama	31	35634	16	48.1	21745	41507	Correct
Morwa	29	10550	16	40.9	5802	9608	Correct
Motihari	33	14987	18	49.8	11764	23447	Correct
Munger	37	1346	36	0.4	885	888	Correct
Muzaffarpur	34	6132	20	41.9	5124	8643	Correct
Nabinagar	30	19926	14	49.1	14704	28899	Correct
Nalanda	32	15878	15	49.6	5022	9756	Correct
Narkatia	30	27377	15	49.1	16619	32266	Correct
Narkatiaganj	28	21519	14	48.1	11391	22076	Correct
Narpatganj	34	28681	16	49.5	12358	24689	Correct
Nathnagar	36	7481	26	24.8	10089	13833	Correct
Nautan	30	26106	16	46.6	15003	28245	Correct
Nawada	37	25835	19	42.7	6251	10608	Correct
Nirmali	31	44195	16	49.7	20916	41440	Correct
Nokha	31	17212	16	48.2	9007	17107	Correct

continued ...

Table A4: Detailed results from all constituencies in Bihar.

Constituency	True data		At the time of calling the race				Decision
	Rounds	Final margin	Round	Votes left (%)	Lead	Predicted margin	
Obra	32	22233	16	49.9	8742	17256	Correct
Paliganj	30	30928	15	48.0	17531	33748	Correct
Parbatta	33	1178	32	0.2	1511	1511	Correct
Parihar	32	1729	16	48.6	3935	7704	Correct
Paroo	32	14722	16	49.5	4446	8661	Correct
Parsa	27	16947	14	46.7	8660	15974	Correct
Patepur	31	25958	16	47.7	13250	25623	Correct
Patna Sahib	39	18281	21	48.9	8388	16381	Correct
Phulparas	36	11198	18	49.3	8810	17390	Correct
Phulwari	38	13870	20	47.6	12749	24323	Correct
Pipra (A)	36	8605	18	49.6	6769	13534	Correct
Pipra (B)	30	19716	15	48.1	7511	14380	Correct
Pirpanti	35	26994	18	48.9	17639	34501	Correct
Pranpur	32	3266	16	46.7	10395	18800	Correct
Purnia	35	32288	20	48.3	24222	45472	Correct
Rafiganj	34	9219	17	48.4	10410	20108	Correct
Raghopur	38	37820	19	47.3	9179	17353	Correct
Raghunathpur	33	17578	15	48.6	13250	25591	Correct
Rajapakar	28	1497	27	0.3	1728	1744	Correct
Rajauli	37	12166	19	47.2	5413	10246	Correct
Rajgir	33	16132	17	46.5	5070	9708	Correct
Rajnagar	34	19388	17	49.0	9794	19390	Correct
Rajpur	34	20565	17	48.8	11750	23016	Correct
Ramgarh	32	305	15	47.8	4488	8404	Correct
Ramnagar	32	16087	16	47.9	11121	21162	Correct
Raniganj	36	2395	34	1.5	2550	2583	Correct
Raxaul	28	37094	14	49.4	23282	45791	Correct
Riga	33	32851	18	48.7	19030	37330	Correct
Rosera	35	35814	19	47.0	30151	56895	Correct
Runnisaidpur	30	24848	16	46.4	22558	41764	Correct
Rupauli	33	19343	16	50.0	15912	31754	Correct
Saharsa	38	20177	20	49.3	23357	46054	Correct
Sahebganj	32	15393	16	48.4	5297	10446	Correct
Sahebpur Kamal	26	13846	13	48.7	9133	17390	Correct
Sakra	28	1742	18	30.7	4361	6242	Correct
Samastipur	30	4588	26	9.1	5765	6425	Correct
Sandesh	30	50109	16	48.0	26711	51168	Correct
Sarairanjan	31	3722	15	49.6	3992	8100	Correct
Sasaram	36	25779	20	48.3	19372	37283	Correct
Shahpur	36	22384	18	49.6	5654	11182	Correct
Sheikhpura	28	6003	21	20.6	4134	5195	Correct
Sheohar	32	36461	16	47.7	16516	31584	Correct
Sherghati	30	16449	15	49.8	10533	21095	Correct

continued ...

Table A4: Detailed results from all constituencies in Bihar.

Constituency	True data		At the time of calling the race				Decision
	Rounds	Final margin	Round	Votes left (%)	Lead	Predicted margin	
Sikandra	32	5668	21	34.4	5237	7854	Correct
Sikta	30	2080	21	24.5	4218	5570	Incorrect
Sikti	29	13716	14	49.5	15465	29903	Correct
Simri Bakhtiarpur	36	1470	34	1.0	1995	2017	Correct
Singheshwar	33	4995	29	6.8	5231	5640	Correct
Sitamarhi	31	11946	27	8.7	7984	8768	Correct
Siwan	36	1561	16	48.1	5874	11223	Incorrect
Sonbarsha	32	13732	17	46.6	4935	9348	Correct
Sonepur	29	6557	15	47.3	4988	9105	Incorrect
Sugauli	30	3045	16	48.6	4810	9193	Correct
Sultanganj	35	11603	28	16.6	6231	7414	Correct
Supaul	29	28246	15	49.2	10219	20448	Correct
Surajgarha	40	9327	19	50.0	4647	9650	Incorrect
Sursand	35	9242	17	49.0	8237	15987	Correct
Taraiya	31	11542	15	49.9	9042	18196	Correct
Tarapur	34	7256	22	31.1	5871	8452	Correct
Tarari	36	10598	21	33.4	3105	4683	Correct
Teghra	31	47495	16	47.8	19584	37529	Correct
Thakurganj	31	23509	16	49.3	11008	21202	Correct
Tikari	35	2745	17	49.7	5312	10259	Incorrect
Triveniganj	29	3402	28	2.8	2499	2552	Correct
Ujiarpur	32	23010	16	48.5	16022	30998	Correct
Vaishali	34	7629	17	48.0	5278	10524	Correct
Valmikinagar	33	21825	16	47.0	15531	29394	Correct
Warisnagar	35	13913	17	47.9	6403	12099	Correct
Warsaliganj	38	9073	19	49.2	8982	17503	Correct
Wazirganj	34	22422	16	49.6	12787	25470	Correct
Ziradei	30	25156	15	47.0	13708	25846	Correct