

# The Feature-First Block Model

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**Abstract:** Labelled networks are an important class of data, naturally appearing in numerous applications in science and engineering. A typical inference goal is to determine how the vertex labels (or *features*) affect the network's structure. In this work, we introduce a new generative model, the feature-first block model (FFBM), that facilitates the use of rich queries on labelled networks. We develop a Bayesian framework and devise a two-level Markov chain Monte Carlo approach to efficiently sample from the relevant posterior distribution of the FFBM parameters. This allows us to infer if and how the observed vertex-features affect macro-structure. We apply the proposed methods to several real-world networks to extract the most important features along which the vertices are partitioned. The approach stands out from its peers in that the whole feature-space is used automatically and features can be rank-ordered implicitly by importance.

**Keywords:** Stochastic Block Model; Labelled Networks; Inference.

## 1 Introduction

Many real-world networks exhibit strong community structure, with most nodes belonging to densely connected clusters. In this work, we examine vertex-labelled networks, referring to the labels as *features*. A typical goal is to determine whether a given feature impacts graphical structure. Answering this requires a random graph model; the standard is the stochastic block model (SBM) – see Peixoto (2017).

Analysing a labelled network with a simple SBM variant, requires partitioning the graph into blocks grouped by distinct values of the feature of interest. The associated model can then be used to test for evidence of heterogeneous connectivity between the feature-grouped blocks. Nevertheless, this approach can only consider disjoint feature sets and the feature-grouped blocks are often an unnatural partition of the graph.

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## 2 The Feature-First Block Model

We would instead prefer to partition the graph into its most natural blocks and then find which of the available features – if any – best predict the resulting partition. Thus motivated, we present a novel framework for modelling labelled networks. This would not be the first extension of the SBM to labelled networks, e.g. Stanley et al (2019). However, most of the current approaches are focused on leveraging feature information to partition the graph more reliably in the presence of noise. We seek instead to develop a model well suited for inferring how vertex features impact graphical structure and to report our confidence in those conclusions.

## 2 Feature-First Block Model

We propose a novel generative model for labelled networks – which we call the feature-first block model (FFBM), illustrated in Figure 1. Let  $N$  denote the number of vertices,  $B$  the number of blocks and  $\mathcal{X}$  the set of values each feature can take. We write  $X$  for the  $N \times D$  *feature matrix* containing the feature vectors  $\{x_i\}_{i=1}^N$  as its rows. For the FFBM, we start with the feature matrix  $X$  and generate a random vector of block memberships  $b \in [B]^N$ . For each vertex  $i$ , the block membership  $b_i \in [B]$  is generated based on the feature vector  $x_i$ , independently between vertices,  $p(b|X, \theta) = \prod_{i \in [N]} \phi_{b_i}(x_i; \theta)$ .

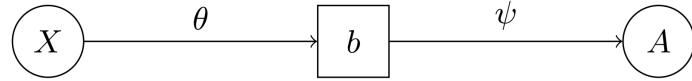


FIGURE 1. The Feature-First Block Model (FFBM)

Once the block memberships  $b$  have been generated, we then draw the adjacency matrix of the graph,  $A$ , from the microcanonical DC-SBM, Peixoto (2017), with additional parameters  $\psi$ ,  $A \sim \text{DC-SBM}_{\text{MC}}(b, \psi)$ .

## 3 Inference

We wish to leverage the FFBM for inference. Specifically, given a labelled network  $(A, X)$ , we wish to infer if and how the observed features  $X$  impact the graphical structure  $A$ . Formally, this means characterising the posterior distribution:  $p(\theta|A, X) \propto p(\theta) \cdot p(A|X, \theta)$ . Therefore, following standard Bayesian practice, instead we aim to draw samples from the posterior,  $\theta^{(t)} \sim p(\theta|A, X)$ . We propose an iterative Markov chain Monte Carlo (MCMC) approach to obtain these samples  $\{\theta^{(t)}\}$ . We first draw a sample  $b^{(t)}$  from the block membership posterior, and then use  $b^{(t)}$  to obtain a corresponding sample  $\theta^{(t)}$ :

$$b^{(t)} \stackrel{\text{distr}}{\approx} p(b|A, X) \quad \text{then} \quad \theta^{(t)} \stackrel{\text{distr}}{\approx} p(\theta|X, b^{(t)}), \quad (1)$$

where these approximations become exact as the number of MCMC iterations  $t \rightarrow \infty$ . Splitting the Markov chain into two levels side-steps the intractable summation over all latent  $b \in [B]^N$  required to directly compute the likelihood,  $p(A|X, \theta)$ . The resulting  $\theta^{(t)}$  samples are asymptotically unbiased in that the expectation of their distribution converges to the true posterior.

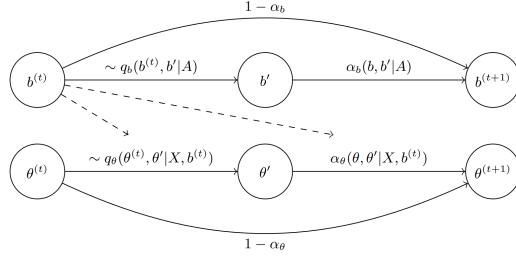


FIGURE 2.  $\theta$ -sample generation.

Figure 2 shows an overview of the proposed method, with  $q$  and  $\alpha$  denoting the Metropolis-Hastings proposal distribution and acceptance probability. Due to the formulation of the FFBM, evaluating  $p(b|X)$  does not depend on  $X$  so we do not need  $X$  to sample  $b$ . And on the other level, in order to obtain samples for  $\theta$  we use only  $b$  but not  $A$ , as  $(\theta \perp\!\!\!\perp A)|b$ .

## 4 Experimental results

We apply our proposed methods to a variety of real-world datasets. For reference, the inferred partitions  $b$  for all of these are given on Figure 3. We require metrics to assess model performance. First, the average description length per entity (nodes and edges)  $\bar{S}_e$  is used to gauge the SBM fit. Second, to assess the performance of the feature-to-block predictor, the vertex set  $[N]$  is partitioned at random into training and test sets,  $\mathcal{G}_0$  and  $\mathcal{G}_1$ . Then the average cross-entropy loss over each set is used to gauge the quality of the fit – denoted  $\bar{\mathcal{L}}_*$ .

For the higher-dimensional datasets, we also develop and apply a novel dimensionality reduction method to select only the top  $D'$  features. We then retrain the feature-block predictor using only the retained feature set, and report the cross-entropy loss over the training and test sets for the reduced classifier – denoted  $\bar{\mathcal{L}}'_*$ .

Table 1 summarises the high-level results for each experiment. We see that the dimensionality reduction procedure brings the training and test losses closer. This indicates that the retained features are indeed well correlated with the underlying graphical partition and that the approach generalises correctly.

TABLE 1. Results averaged over  $n = 10$  iterations (mean  $\pm$  std. dev.).

Dataset	$B$	$D$	$D'$	$S_e$	$\hat{L}_0$	$\hat{L}_1$	$c^*$	$\hat{L}'_0$	$\hat{L}'_1$
Polbooks	3	3	—	$2.250 \pm 0.000$	$0.563 \pm 0.042$	$0.595 \pm 0.089$	—	—	—
School	10	13	10	$1.894 \pm 0.004$	$0.787 \pm 0.127$	$0.885 \pm 0.129$	$1.198 \pm 0.249$	$0.793 \pm 0.132$	$0.853 \pm 0.132$
FB egonet	10	480	10	$1.626 \pm 0.003$	$1.326 \pm 0.043$	$1.538 \pm 0.069$	$0.94 \pm 0.019$	$1.580 \pm 0.150$	$1.605 \pm 0.106$

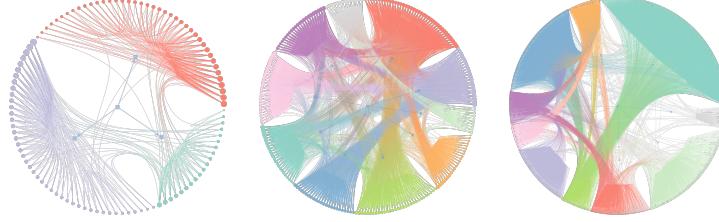


FIGURE 3. Networks laid out and coloured according to inferred block memberships. Left to right: Polbooks, Krebs (2004); Primary School, Stehle et al (2011); Facebook Egonet, Leskovec and Mcauley (2012).

## 5 Conclusion

An efficient MCMC algorithm is developed for sampling from the posterior distribution of the relevant parameters in the FFBM; the main idea is to divide up the graph into its most natural partition under the associated parameter values, and then to determine whether the vertex features can accurately explain the partition. Through several applications on empirical network data, this approach is shown to be effective at extracting and describing the most natural communities in a labelled network. Nevertheless, it can only currently explain the structure at the macroscopic scale. Future work will benefit from extending the FFBM to a further hierarchical model, so that the structure of the network can be explained at all scales of interest.

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