OSCAR.MADRID@STAT.UCLA.EDU

YI.YU.2@WARWICK.AC.UK

Change point localization in dependent dynamic nonparametric random dot product graphs

Oscar Hernan Madrid Padilla

Department of Statistics University of California Los Angeles, CA 90095-1554, USA

Yi Yu

Department of Statistics University of Warwick Coventry, CV4 7AL, UK

Carey E. Priebe

Department of Applied Mathematics and Statistics Johns Hopkins University Baltimore, MD 21218-2682, USA CEP@JHU.EDU

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Abstract

In this paper, we study the offline change point localization problem in a sequence of dependent nonparametric random dot product graphs. To be specific, assume that at every time point, a network is generated from a nonparametric random dot product graph model (see e.g. Athreya et al., 2018), where the latent positions are generated from unknown underlying distributions. The underlying distributions are piecewise constant in time and change at unknown locations, called change points. Most importantly, we allow for dependence among networks generated between two consecutive change points. This setting incorporates edge-dependence within networks and temporal dependence between networks, which is the most flexible setting in the published literature.

To accomplish the task of consistently localizing change points, we propose a novel change point detection algorithm, consisting of two steps. First, we estimate the latent positions of the random dot product model, our theoretical result being a refined version of the state-of-the-art results, allowing the dimension of the latent positions to diverge. Subsequently, we construct a nonparametric version of the CUSUM statistic (e.g. Page, 1954; Padilla et al., 2019a) that allows for temporal dependence. Consistent localization is proved theoretically and supported by extensive numerical experiments, which illustrate state-of-the-art performance. We also provide in depth discussion of possible extensions to give more understanding and insights.

Keywords: Dependent dynamic networks, Nonparametric random dot product graph models, Change point localization.

1. Introduction

Computationally-efficient and theoretically-justified change point localization methods that can handle new data types are in high demand, due to technological advances in a broad

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range of application areas including finance, biology, social sciences, to name only a few. The literature on change point detection is extensive, including the univariate mean case (e.g. Frick et al., 2014; Fryzlewicz, 2014; Wang et al., 2018b), the high-dimensional mean case (e.g. Wang and Samworth, 2016; Cho, 2016), the robust mean case (e.g. Fearnhead and Rigaill, 2018; Pein et al., 2017), the covariance case (e.g. Aue et al., 2009; Wang et al., 2017; Avanesov and Buzun, 2018), the univariate nonparametric case (e.g. Zou et al., 2014; Padilla et al., 2019a), and the multivariate nonparametric case (e.g. Arlot et al., 2012; Matteson and James, 2014; Garreau and Arlot, 2018; Padilla et al., 2019b).

In this paper we are concerned with offline change point localization in dynamic networks. Let $\{A(t)\}_{t=1}^T \subset \{0,1\}^{n \times n}$ be a sequence of adjacency matrices generated from a sequence of distributions $\{\mathcal{L}_t\}_{t=1}^T$, such that for an unknown sequence of change points $\{\eta_k\}_{k=1}^K \subset \{2,\ldots,T\}$ with $1 = \eta_0 < \eta_1 < \ldots < \eta_K \leq T < \eta_{K+1} = T + 1$, we have that

 $\mathcal{L}_{t-1} \neq \mathcal{L}_t$, if and only if $t \in \{\eta_1, \ldots, \eta_K\}$.

The goal is to estimate the change point collection $\{\eta_k\}_{k=1}^K$ accurately.

The model described can be used to study various application problems. For instance, in epidemiology, studying the dynamic networks formed by human interaction can facilitate the detection of transmissible diseases outbreaks; in finance, dynamic networks formed by within-companies transactions over time may possess abrupt changes which indicate abnormal market behaviours; in neuroscience, we provide a detailed real data example in the context of detecting changes in the neuronal activity in Section 4.2. In response to the growing demand from applications, there has been recently an increasing interest in the literature studying the model described above. Wang et al. (2018a) considered an independent sequence of inhomogeneous Bernoulli networks and presented a nearly optimal change point localization algorithm, accompanied with a phase transition phenomenon. Zhao et al. (2019) assumed an independent sequence of graphon models with independent edges and proposed consistent yet optimal localization result. Other network change point papers include Wang et al. (2014), Cribben and Yu (2017), Liu et al. (2018), Chu and Chen (2017), Mukherjee (2018), among others. We would like to mention that both Cribben and Yu (2017) and Liu et al. (2018) have exploited the eigenvectors information to conduct change point detection, but both lack theoretical results. Our paper, to the best of our knowledge, is the first to provide theoretical justifications for eigenvector-based change point detection methods. More in-depth comparisons with Wang et al. (2018a) will be conducted later in the paper.

1.1 Random dot product graph models

Different from the aforementioned papers, in order to allow for dependence among edges, we assume that at every time point, the network is generated from a random dot product graph (e.g. Young and Scheinerman, 2007; Athreya et al., 2018). We formally define the model in Definitions 1 and 2, which are both based on Athreya et al. (2018).

Definition 1 (Inner product distribution) Let F be a probability distribution whose support is given by $\mathcal{X}_F \subset \mathbb{R}^d$. We say that F is a d-dimensional inner product distribution on \mathbb{R}^d if for all $x, y \in \mathcal{X}_F$, it holds that $x^\top y \in [0, 1]$. **Definition 2 (Random dot product graph with distribution** F) Let F be an inner product distribution with $\{X_i\}_{i=1}^n \stackrel{i.i.d.}{\sim} F$. Let $X = (X_1, \ldots, X_n)^\top \in \mathbb{R}^{n \times d}$. Suppose A is a random adjacency matrix given by

$$\mathbb{P}\{A \mid X\} = \prod_{1 \le i < j \le n} (X_i^\top X_j)^{A_{ij}} (1 - X_i^\top X_j)^{1 - A_{ij}}.$$
 (1)

We write $A \sim \text{RDPG}(F, n)$.

We would like to make a few comments regarding random dot product graph models. For first time reading, one can safely skip this and jump to Section 1.2.

Equivalence of distributions

It can be seen from Definition 2 that the latent positions come into play only through their inner products, i.e. we have

$$A_{ij} \sim \operatorname{Ber}(X_i^{\top} X_j), \quad 1 \le i, j \le n.$$

This means that one can apply any orthonormal rotations to all the latent positions and retain the same distribution of A. In light of this rotational invariance, we define the equivalence of inner product distributions below, which is also from Athreya et al. (2018).

Definition 3 (Equivalence of inner product distributions) If both $F(\cdot)$ and $G(\cdot)$ are inner product distributions defined on \mathbb{R}^d , and there exists an orthogonal operator $U : \mathbb{R}^d \to \mathbb{R}^d$ such that $F = G \circ U$, then we say F and G are equivalent.

Community structures

The random dot product graph is a generalization of the stochastic block model (Holland et al., 1983), where the latent positions X are assumed to be fixed and satisfy

$$XX^{\top} = ZQZ^{\top},$$

where $Z \in \{0, 1\}^{n \times d}$ is a membership matrix, with each row consisting of one and only one entry being 1 and $Q \in [0, 1]^{d \times d}$ is a connectivity matrix encoding the edge probabilities.

One may be puzzled by the observation that under Definition 2, we have that for any $(i, j) \in \{1, ..., n\}^2, i \neq j$,

$$\mathbb{E}(A_{ij}) = \mathbb{E}(X_i^{\top} X_j) = \mathbb{E}(X_1^{\top} X_2),$$

where the second identity follows from the fact that within a network the latent positions are i.i.d., and therefore one loses the community structure and connections from the stochastic block model.

This observation is due to the randomness of the latent positions. To enforce a version of "communities" under Definition 2, one may introduce a membership vector and treat the distribution F as a mixture distribution. To be specific, we have an alternative to Definition 2 below.

Definition 4 Let τ_1, \ldots, τ_n be *i.i.d.* random variables satisfying

$$\mathbb{P}\{\tau = m\} = \pi_m, \quad \pi_m \ge 0, \, m \in \{1, \dots, M\}, \, \sum_{m=1}^M \pi_m = 1,$$

where M is a positive integer. Let $\{F_m\}_{m=1}^M$ be a sequence of d-dimensional inner product distributions. Assume that

$$X_i \mid \tau_i \stackrel{ind.}{\sim} F_{\tau_i}, \quad i = 1, \dots, n.$$

Let $X = (X_1, \ldots, X_n)^{\top} \in \mathbb{R}^{n \times d}$. Suppose A is a random adjacency matrix given by

$$\mathbb{P}\{A \mid X\} = \prod_{1 \le i < j \le n} (X_i^{\top} X_j)^{A_{ij}} (1 - X_i^{\top} X_j)^{1 - A_{ij}}$$

We write $A \sim \text{RDPG}(F, n)$, where

$$F = \sum_{m=1}^{M} \pi_m F_m.$$

We remark that Definition 4 is a special case of Definition 2. Therefore the theoretical results based on Definition 2 also hold for Definition 4. The vector τ prompts the vertex correspondence in a dynamic network. For instance, one may assume a sequence of RDPG(F, n) using Definition 4, with latent positions drawn independently and the membership vector unchanged. There are also other variants. For instance, one may also assume instead that the membership vector τ is fixed.

1.2 List of contributions

We highlight the contributions of this paper.

First of all, we propose a novel algorithm for change point localization in dependent dynamic random dot product graph models, see Algorithm 2. This proceeds by first estimating the latent positions $\{\hat{X}_i(t)\}_{i=1,t=1}^{n,T}$. However, due to the latent positions' rotationalinvariance properties discussed in Section 1.1, one pertaining challenge in the RDPG literature is to match the rotations of the latent position estimators of different networks (e.g. Athreya et al., 2018; Cape et al., 2019). We propose a novel way to get around this issue with matching. Specifically, we define $\hat{Y}_{ij}^t = (X_i(t))^\top X_j(t)$, and construct a Kolmogorov– Smirnov CUSUM statistic (Padilla et al., 2019a) based on $\{\hat{Y}_{ij}^t : (i,j) \in \{(l,n/2+l), l = 1, \ldots, n/2\}, t = 1, \ldots, T\}$. One may question the power of the Kolmogorov–Smirnov distance, but it allows for more general distributions for latent positions. Among those distributions stochastic block models are special cases. One may also question the effectiveness of using only a subset of all the possible edges, we will discuss in Section 3.3 that in terms of order, this is in fact the same as using all possible edges.

Secondly, under an appropriate signal-to-noise ratio condition, we prove that our proposed method (Algorithm 2) can estimate the number and locations of change points consistently, which will be formally stated in Section 3.2. It is worth mentioning that Theorem 9 handles the situation where there exists dependence across time and among edges. This is not shown in the existing network change point detection literature. To be more specific, the dependence among edges are imposed by assuming the latent positions are random and the edges are conditionally independent given the latent positions. Our proposed method is also robust to some model mis-specification, see the discussions following Theorem 9 for details.

Thirdly, we provide in-depth discussions on the characterization of jumps in Section 3.1. Note that the data we have are a collection of adjacency matrices. However, as stated in Definition 2, the data generating mechanism depends on latent positions' distributions Fs. A natural question is whether the changes in F will lead to the changes in the distributions of the adjacency matrices, and if so, whether we can characterize the changes. The results we developed in Section 3.1 are interesting *per se*, and can shed light on network testing problems.

Lastly, numerical experiments provide ample evidence on the strength of our proposed approach. In particular, we highlight the advantage of our method in scenarios with dependent networks.

The rest of the paper is organized as follows. Section 2 provides the formal problem setup and our proposed method in detail, including discussions on possible extensions. The characterization of the distributional changes and statistical guarantees for our approach are collected in Section 3. We conclude with numerical experiments in Section 4 and final discussions in Section 5. Technical details are deferred to the Appendix.

2. Methodology

2.1 Setup

We first formally state the full model descriptions.

Model 1 Let $\{A(1), \ldots, A(T)\} \subset \mathbb{R}^{n \times n}$ be a sequence of adjacency matrices of random dot product graphs, satisfying the following.

1. (Random dot product graphs) For any $t \in \{1, ..., T\}$, it holds that

$$\mathbb{P}\left\{A(t) \mid X(t)\right\} = \prod_{1 \le i < j \le n} (X_i(t)^\top X_j(t))^{A_{ij}(t)} (1 - X_i(t)^\top X_j(t))^{1 - A_{ij}(t)},$$

where $X(t) = (X_1(t), \dots, X_n(t))^\top \in \mathbb{R}^{n \times d}$ satisfies the following.

There exists a sequence $1 = \eta_0 < \eta_1 < \ldots < \eta_K \leq T < \eta_{K+1} = T + 1$ of time points, called change points. For $k \in \{0, \ldots, K\}$, we have that

$$X_i(\eta_k) \in \mathbb{R}^d \stackrel{\text{ind}}{\sim} F_{\eta_k}, \quad i = 1, \dots, n,$$

and for $t \in \{\eta_k + 1, \dots, \eta_{k+1} - 1\}$, we have that

$$X_{i}(t) \begin{cases} = X_{i}(t-1), & \text{with probability } \rho, \\ \underset{\sim}{\text{ind}} F_{\eta_{k}}, & \text{with probability } 1-\rho, \end{cases}$$
(2)

and with F_t 's satisfying Definition 1. Throughout, we write $P_t = X(t)X(t)^{\top}$ for the matrix of latent link probabilities at time $t \in \{1, \ldots, T\}$.

2. (Minimal spacing) The minimal spacing between two consecutive change points satisfies

$$\min_{k=1,\dots,K+1} \{\eta_k - \eta_{k-1}\} = \Delta > 0.$$

3. (Minimal jump size) For each $k \in \{0, \ldots, K\}$ and for any $X, Y \stackrel{i.i.d.}{\sim} F_{\eta_k}$, denote

$$G_{\eta_k}(z) = \mathbb{P}\left\{ X^\top Y \le z \right\}, \quad z \in [0, 1]$$

The magnitudes of the changes in the data generating distribution are such that

$$\kappa = \min_{k=1,\dots,K+1} \kappa_k = \min_{k=1,\dots,K} \sup_{z \in [0,1]} |G_{\eta_k}(z) - G_{\eta_{k-1}}(z)| > 0.$$
(3)

4. Assume that for every $k \in \{0, \ldots, K\}$ and $i \in \{1, \ldots, n\}$,

$$\mathbb{E}\left\{X_i(\eta_k)X_i(\eta_k)^{\top}\right\} = \Sigma_k \in \mathbb{R}^{d \times d},$$

where Σ_k has eigenvalues $\mu_1^k \geq \cdots \geq \mu_d^k > 0$, with $\{\mu_l^k, k = 0, \dots, K, l = 1, \dots, d\}$ all being universal constants.

In Model 1, between two consecutive change points, the latent positions are dependent with exponentially decaying correlations; and for latent positions drawn at time points separated by change points, they are independent. If $\rho = 0$ in (2), then all the latent positions are independent, which implies that the adjacency matrices are independent.

The distributional changes occurring at change points are quantified through cumulative distribution functions $\{G_{\eta_k}\}$ defined in Model 1(3). Intuitively, since the unconditional distributions of $\{A(t)\}$ are completely characterized by the joint distributions of $\{X_i(t)^{\top}X_j(t)\}$, it is natural to quantify the changes with respect $\{G_{\eta_k}\}$. (A more detailed discussion on this can be found in Section 3.1.) In particular, the changes are measured by the Kolmogorov– Smirnov distance in (3), since the Kolmogorov–Smirnov distance does not require assumptions about the moments of the distributions, or about their discrete/continuous nature. With the stochastic block model being a special case of the random dot product graph, the distributions thereof are point-mass distributions, which handicaps the adoption of other (potentially more powerful) distribution distances, including the total variation distance.

Model 1(4) is imposed to guarantee that the latent link probabilities satisfy rank(P_t) = d with high probability. Without this full-rank assumption, assuming that $r = \operatorname{rank}(\Sigma_k) < d$ implies that there exists a rank-r subspace which characterises the latent positions, and the effective dimension is r instead of d. For simplicity, we assume that Σ_k is of full rank.

2.2 Methods

To arrive at our construction, we start by defining the main statistic, and its population version. Without loss of generality, we assume that the number of nodes n is an even integer. If n is odd, then we randomly ignore a certain but fixed node and all edges connecting to it throughout the whole procedure.

Definition 5 (CUSUM statistics) Let $\mathcal{O} = \{(i, n/2 + i), i = 1, \dots, n/2\}.$

• (Sample version) With $\{A(t)\}_{t=1}^T \subset \mathbb{R}^{n \times n}$, let

$$\widehat{X}(t) = U_A(t)\Lambda_A(t)^{1/2}$$

where $U_A(t) \in \mathbb{R}^{n \times d}$ is an orthogonal matrix with columns being the leading d eigenvectors of A(t), and $\Lambda_A(t) \in \mathbb{R}^{d \times d}$ is a diagonal matrix with entries being the largest d, in absolute value, eigenvalues of A(t).

For any $t \in \{1, \ldots, T\}$ and $(i, j) \in \mathcal{O}$, let

$$\widehat{Y}_{ij}^t = \widehat{X}_i(t)^\top \widehat{X}_j(t),$$

where $\widehat{X}_i(t)^{\top}$ is the *i*th row of $\widehat{X}(t)$. For any integer triplet $(s, t, e), 0 \leq s < t < e \leq T$ and $z \in \mathbb{R}$, we define the CUSUM statistic as

$$D_{s,e}^{t}(z) = \left| \sqrt{\frac{2(e-t)}{n(e-s)(t-s)}} \sum_{k=s+1}^{t} \sum_{(i,j)\in\mathcal{O}} \mathbb{1}\{\widehat{Y}_{ij}^{k} \le z\} - \sqrt{\frac{2(t-s)}{n(e-s)(e-t)}} \sum_{k=t+1}^{e} \sum_{(i,j)\in\mathcal{O}} \mathbb{1}\{\widehat{Y}_{ij}^{k} \le z\} \right|,$$
(4)

and

$$D_{s,e}^t = \sup_{z \in [0,1]} |D_{s,e}^t(z)|.$$

• (Population version) With $\{A(t)\}_{t=1}^T \subset \mathbb{R}^{n \times n}$, recall that $P_t = X(t)X(t)^\top$ and write

$$X(t) = U_P(t)\Lambda_P(t)^{1/2}$$

where $U_P(t) \in \mathbb{R}^{n \times d}$ is an orthogonal matrix with columns being the d eigenvectors of P_t with largest absolute eigenvalues, and $\Lambda_P(t) \in \mathbb{R}^{d \times d}$ is a diagonal matrix with entries being the leading d eigenvalues of P_t .

For any $t \in \{1, \ldots, T\}$ and $(i, j) \in \mathcal{O}$, let

$$Y_{ij}^t = X_i(t)^\top X_j(t), \tag{5}$$

where $X_i(t)^{\top}$ is the *i*th row of X. For any integer triplet (s, t, e), $0 \le s < t < e \le T$ and $z \in \mathbb{R}$, we define the CUSUM statistic as

$$\begin{split} \widetilde{D}_{s,e}^t(z) &= \left| \sqrt{\frac{2(e-t)}{n(e-s)(t-s)}} \sum_{k=s+1}^t \sum_{(i,j)\in\mathcal{O}} \mathbb{E}\left(\mathbbm{1}\{Y_{ij}^k \le z\}\right) \right. \\ &- \sqrt{\frac{2(t-s)}{n(e-s)(e-t)}} \sum_{k=t+1}^e \sum_{(i,j)\in\mathcal{O}} \mathbb{E}\left(\mathbbm{1}\{Y_{ij}^k \le z\}\right) \end{split}$$

and

$$\widetilde{D}_{s,e}^t = \sup_{z \in [0,1]} |\widetilde{D}_{s,e}^t(z)|.$$

We remark that in Definition 5, if the *d*th and (d + 1)th eigenvalues share the same value, then one can randomly pick an eigenvector to construct $\hat{X}, X \in \mathbb{R}^{n \times d}$. In addition, we do not require a specific order of the eigenvectors in constructing \hat{X} and X.

Recall that the distributions of the latent positions are equivalent up to a rotation, see Definition 3. To avoid extra efforts in matching the rotations when comparing two latent position distributions, we resort to the inner products of latent positions instead of latent positions itself. We explain this via (5). For any orthogonal matrix $U \in \mathbb{R}^{d \times d}$, it holds that

$$Y_{ij}^t = (X_i(t))^{\top} X_j(t) = (UX_i(t))^{\top} UX_j(t).$$

With Definition 5, we arrive at our proposed procedure Algorithm 2 that builds on the wild binary segmentation algorithm (Fryzlewicz, 2014). The method requires first estimating the latent positions, a subroutine shown in Algorithm 1 (adjacency spectral embedding, see e.g. Sussman et al., 2012). Note that this only needs to be done once regardless of the choice of the tuning parameter τ , and is parallelizable. Since the complexity of the truncated principal component analysis is of order $O(dn^2)$, Algorithm 1 has the computational cost of order $O(Tdn^2)$. Once the latent positions are estimated, we run the remaining steps in Algorithm 2, which amounts to running Algorithm 2 in Padilla et al. (2019a). For a fixed τ which leads to \tilde{K} change points, we have the computational complexity of order $O(\tilde{K}MTn\log(n))$, which translates to $O(Tdn^2 + \tilde{K}MTn\log(n))$ for the overall cost of Algorithm 2, where M is the number of random intervals drawn in Algorithm 2.

Algorithm 1 ScaledPCA (A, d)

INPUT: Matrix $A \in \mathbb{R}^{n \times n}$, and tuning parameter $d \in \mathbb{Z}_+$. $A = (v_1, \ldots, v_n) \operatorname{diag}(\lambda_1, \ldots, \lambda_n) (v_1, \ldots, v_n)^\top$, where $|\lambda_1| \ge \ldots \ge |\lambda_n|$. $X \leftarrow (v_1, \ldots, v_d) \operatorname{diag}(|\lambda_1|^{1/2}, \ldots, |\lambda_d|^{1/2})$ **OUTPUT:** X

In every network, there are n(n-1)/2 observations, but note that in Definition 5, we in fact only use n/2 of them. This is for technical convenience, since due to the choice of \mathcal{O} , we obtain independent observations within one network. We acknowledge that there are other variants of this treatment. For instance, instead using a fixed choice of \mathcal{O} , one can do multiple random sub-samplings and combine the results; one can also gather all the observations and create a U-statistic instead. In Section 3.3, we will show that in terms of rate, using n/2 edges is as effective as using all possible edges.

3. Theory

In this section, we provide the statistical guarantees for Algorithm 2. In order to enhance the theoretical understanding, we take a step back and understand how the jump defined in (3) through the cumulative distribution functions of the inner products can be related to the jumps in terms of the distributions of the adjacency matrices. The main results which provide theoretical guarantees of our algorithm are collected in Theorem 9.

3.1 Characterizations of the changes

We summarize the notation below and consider two different sets of models.

Algorithm 2 NonPar-RDPG-CPD $((s, e), \{(\alpha_m, \beta_m)\}_{m=1}^M, \tau)$

INPUT: A sample $\{A(t)\}_{t=s+1}^e \subset \mathbb{R}^{n \times n}$, collection of intervals $\{(\alpha_m, \beta_m)\}_{m=1}^M$, tuning parameters $d \in \mathbb{Z}_+$, and $\tau > 0$. for t = s + 1, ..., e do $X(t) \leftarrow \text{ScaledPCA}(A(t), d)$ end for for $m = 1, \ldots, M$ do $(s_m, e_m) \leftarrow [s, e] \cap [\alpha_m, \beta_m]$ if $e_m - s_m > 1$ then $b_m \leftarrow \operatorname{argmax}_{s_m+1 \le t \le e_m-1} D^t_{s_m, e_m}$ $a_m \leftarrow D^{b_m}_{s_m,e_m}$ else $a_m \leftarrow -1$ end if end for $m^* \leftarrow \operatorname{argmax}_{m=1,\dots,M} a_m$ if $a_{m^*} > \tau$ then add b_{m^*} to the set of estimated change points NonPar-RDPG-CPD($(s, b_{m*}), \{(\alpha_m, \beta_m)\}_{m=1}^M, \tau$) NonPar-RDPG-CPD($(b_{m*}+1, e), \{(\alpha_m, \beta_m)\}_{m=1}^M, \tau$) end if **OUTPUT:** The set of estimated change points.

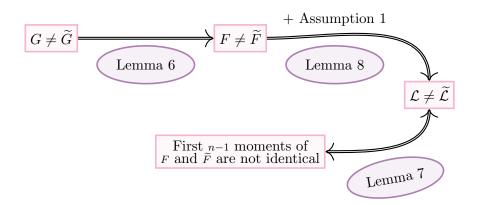


Figure 1: Flowchart of Section 3.1. The notation $A \Rightarrow B$ means A implies B.

Model 2 We assume the following two independent models:

$$\{A_{ij}, 1 \le i < j \le n\} | \{X_i\}_{i=1}^n \stackrel{\text{ind}}{\sim} \operatorname{Ber}(X_i^\top X_j), \quad X_i \stackrel{\text{ind}}{\sim} F \in \mathbb{R}^d;$$

and

$$\{\widetilde{A}_{ij}, 1 \le i < j \le n\} | \{\widetilde{X}_i\}_{i=1}^n \stackrel{\text{ind}}{\sim} \operatorname{Ber}(\widetilde{X}_i^\top \widetilde{X}_j), \quad \widetilde{X}_i \stackrel{\text{ind}}{\sim} \widetilde{F} \in \mathbb{R}^d.$$

For $i \neq j$, the cumulative distribution functions of $X_i^{\top} X_j$ and $\widetilde{X}_i^{\top} \widetilde{X}_j$ are denoted by $G(\cdot)$ and $\widetilde{G}(\cdot)$, respectively. We further write \mathcal{L} and $\widetilde{\mathcal{L}}$ for the joint unconditional distributions of $\{A_{ij}, 1 \leq i < j \leq n\}$ and $\{\widetilde{A}_{ij}, 1 \leq i < j \leq n\}$, respectively.

The rest of this subsection is summarized in Figure 1. The notation $A \Rightarrow B$ means A implies B.

Lemma 6 With the notation in Model 2, if $F = \widetilde{F}$, then $G = \widetilde{G}$.

This follows automatically from the definitions, and is equivalent to the claim that if $G \neq \tilde{G}$ then $F \neq \tilde{F}$, which implies that (3) is equivalent to

$$F_{\eta_k} \neq F_{\eta_{k-1}}, \quad k \in \{1, \dots, K\}.$$

However, $F \neq \tilde{F}$ does not imply $\mathcal{L} \neq \tilde{\mathcal{L}}$. As a simple toy example, consider F and \tilde{F} to be defined in Definition 1, with the same mean but different variances, and n = 2. Then $F \neq \tilde{F}$ but $\mathcal{L} = \tilde{\mathcal{L}}$. 7 below shows that \mathcal{L} is determined by the first n-1 moments of F.

Lemma 7 Under Model 2, we have that $\mathcal{L} = \widetilde{\mathcal{L}}$ if and only if there exists an orthogonal operator $U \in \mathbb{R}^{d \times d}$, such that if d = 1,

$$\mathbb{E}_F(X_1^k) = \mathbb{E}_{\widetilde{F}}\{(U\widetilde{X}_1)^k\}, \quad k = 1, \dots, n-1,$$

if d > 1

$$\mathbb{E}_{F}\left(\prod_{l=1}^{d} X_{1,l}^{k_{l}}\right) = \mathbb{E}_{\widetilde{F}}\left\{\prod_{l=1}^{d} (U\widetilde{X}_{1})_{l}^{k_{l}}\right\}, \quad k_{l} \in \mathbb{Z}, \quad k_{l} \ge 0, \quad \sum_{l=1}^{d} k_{l} = k, \quad k = 1, \dots, n-1,$$

where $X_{1,l}$ and $(U\widetilde{X}_1)_l$ are the *l*th coordinates of the X_1 and $U\widetilde{X}_1$.

It can be seen from Lemma 7 that the unconditional distribution of the data matrix is determined by the first n-1 moments of the underlying distribution F. Unfortunately, without additional assumptions, the first n-1 moments do not determine the distribution (e.g. Heyde, 1963)¹. This means that only assuming (3) can not guarantee that the data matrices A and \tilde{A} have different distributions.

The final claim we make in this subsection is that under some additional but weak conditions, we will be able to guarantee that $\mathcal{L} \neq \widetilde{\mathcal{L}}$.

Assumption 1 Under Model 2, let

$$\kappa_0 = \sup_{z \in [0,1]} |G(z) - G(z)|.$$

It holds that

$$\kappa_0 \sqrt{n} > 3\sqrt{\log(n)}.$$

Lemma 8 Assume that Model 2 and Assumption 1 hold. Then we have that

 $\mathcal{L} \neq \widetilde{\mathcal{L}}.$

Lemma 8 suggests that under Assumption 1, $G \neq \tilde{G}$ implies $\mathcal{L} \neq \tilde{\mathcal{L}}$. This enhances the rationale of imposing the distributional changes occurring at the change points on the differences on G, as detailed in Model 1(4). Assumption 1 is a weak assumption, which will be further elaborated in Section 3.2. The proofs of Lemmas 7 and 8 are collected in Appendix A.

3.2 Consistent estimation of change points

We first state a signal-to-noise ratio condition below.

Assumption 2 (Signal-to-noise ratio) There exists a universal constant $C_{\text{SNR}} > 0$, such that there exists a diverging sequence $a_T \to \infty$, as $T \to \infty$, satisfying

$$\kappa \sqrt{\Delta n(1-\rho)} > C_{\text{SNR}} \sqrt{T} \max\{\sqrt{d \log(n \vee T)}, \, d^{3/2}\} + a_T.$$

To better understand Assumption 2, we would like to use Assumptions 2 and 3 in Wang et al. (2018a) as benchmarks, since Wang et al. (2018a) studied a simpler problem assuming independence within and across networks, and showed a phase transition phenomenon in the minimax sense. However, we would like to emphasize that comparing Assumption 2 and Assumptions 2 and 3 in Wang et al. (2018a) is comparing apples and oranges, to some extent. Even though the jump size κ are defined differently in these two papers, both take values in (0, 1]. The parameter ρ in this paper indicates the correlation between networks, while the parameter ρ in Wang et al. (2018a) represents the entrywise sparsity. For simplicity, we let $\rho = 1$ in Wang et al. (2018a) for this discussion.

One key difference is that in Assumption 2, the required signal-to-noise ratio is inflated by $\sqrt{1-\rho}$. We might view this as the effective sample size being shrunk from Δ to $(1-\rho)\Delta$,

^{1.} We are grateful to Richard J. Samworth for this reference and constructive discussions.

due to the dependence across time. In Model 1, we do not allow $\rho = 1$, but allow $\rho \to 1$, as long as Assumption 2 holds. In the extreme case that $\rho = 1$, between two consecutive change points, there is essentially only one observation. As long as Assumption 1 holds, Lemma 8 shows that the distributions of the adjacency matrices before and after change points are different, which implies that one can identify the change points with probability 1.

Another difference is that in our paper, the signal-to-noise ratio is inflated by \sqrt{T} compared to Wang et al. (2018a). This is due to the fact that we estimate the latent positions separately for every single network, while the graphons were estimated based on a version of sample average of the adjacency matrices in Wang et al. (2018a). The reason we estimate the positions separately roots in the difficulty of deriving theoretical properties of eigenvectors of a sample average matrices. More discussions on this can be found in Section 3.3.

We allow the dimensionality d to diverge, provided that Assumption 2 holds. The dimensionality d is essentially the low rank condition imposed in Wang et al. (2018a). The upper bound on the rank r in Wang et al. (2018a) comes into play with the term \sqrt{r} , while we have $d^{3/2}$ here. The difference again is rooted in the estimation of the latent positions, although we do not claim optimality here.

The sequence a_T can diverge at any arbitrarily slow rate. We will explain the role of a_T after we state Theorem 9.

Finally, we make connections between Assumptions 1 and 2. Recall that we use Assumption 1 in Lemma 8, where only one observation is available for each distribution, i.e. $\Delta = 1$, $\rho = 0$ and T = 2. Ignoring the universal constants, the only difference left between Assumptions 1 and 2 is the term $d^{3/2}$. Of course, if d = O(1), then this is also a universal constant, and there is no difference left. The interesting thing happens when d is allowed to diverge faster than the poly-logarithm term. Assumption 1 is required to differentiate two different distributions, which roughly speaking is related to a testing task; while Assumption 2 is used below in Theorem 9 with the purpose of consistent localization, which is an estimation problem. To this end, the extra $d^{3/2}$ in Assumption 2 is a piece of evidence that estimation is a harder problem than a testing one.

Theorem 9 Let data be from Model 1 and satisfy Assumption 2. Assume the following.

• The tuning parameter τ in Algorithm 2 satisfies

$$c_{\tau,1}T^{1/2}\max\{\sqrt{d\log(n\vee T)}, d^{3/2}\} < \tau < c_{\tau,2}\kappa\sqrt{\Delta n(1-\rho)},$$
 (6)

where $c_{\tau,1}, c_{\tau,2} > 0$ are universal constants depending on all the universal constants in Model 1 and Assumption 2.

- The tuning parameter d in Algorithms 1 and 2 are the true dimension d of the latent positions.
- The intervals satisfy

$$\max_{m=1,\dots,M} (\alpha_m - \beta_m) \le C_R \Delta,\tag{7}$$

where $C_R > 3/2$ is a universal constant.

Let $\{\widehat{\eta}_k\}_{k=1}^{\widehat{K}}$ be the output of Algorithm 2. We have that

$$\mathbb{P}\left\{\widehat{K}=K, \quad |\widehat{\eta}_k - \eta_k| \le C_{\epsilon} \frac{T \max\{d \log(n \lor T), d^3\}}{\kappa_k^2 n(1-\rho)}, \forall k\right\}$$
$$\ge 1 - C(n \lor T)^{-c} - CTe^{-n} - \exp\left(\log(T/\Delta) - (4C_R)^{-1}T^{-1}M\Delta\right),$$

where C, c > 0 are universal constants depending only on the other universal constants.

The proof of Theorem 9 can be found in Appendix D, following two sets of lemmas – technical details on estimating the latent positions and on change point analysis, collected in Appendices B and C, respectively.

Suppose that $Te^{-n} \to 0$ and that M satisfies

$$\frac{T/\Delta\log\left(T/\Delta\right)}{M} \to 0.$$
(8)

Then it can be seen from Theorem 9 that with probability tending to 1, as T diverges, we have that $\hat{K} = K$ and

$$\max_{k=1,\dots,K} \frac{|\widehat{\eta}_k - \eta_k|}{\Delta} \le \max_{k=1,\dots,K} C_{\epsilon} \frac{T \max\{d \log(n \lor T), d^3\}}{\Delta \kappa_k^2 n(1-\rho)} \le C_{\epsilon} \frac{T \max\{d \log(n \lor T), d^3\}}{\Delta \kappa^2 n(1-\rho)} \to 0,$$

where the second inequality follows from the definition of κ and the convergence follows from Assumption 2, with the aid of an arbitrarily diverging sequence a_T . This implies that the change point estimators we obtain are consistent, with a vanishing localization rate. Since the quantity M only appears in the probability, we remark that the larger M is, the more likely that our estimators would perform satisfactorily, while the higher the computational cost is.

Algorithm 2 in fact can handle networks of varying size. For instance, if we do not allow for the dependence across time, then Theorem 9 holds provided that all network sizes are of the same order, which amounts to $c_1n \leq n_t \leq c_2n$, $t = 1, \ldots, T$, for universal constants $c_1, c_2 > 0$.

In view of Theorem 9, the two most important tuning parameters in Algorithm 2 are the threshold τ and the dimension d.

The threshold τ is set to satisfy (6). The upper and lower bounds in (6) are the lower bound on the signals and the upper bound on the noise, in a large probability event, respectively. If the input τ is larger than the upper bound in (6), then one may not be able to detect all change points; while if the input is smaller than the lower bound, then there is the risk of falsely detecting change points. Note that both the upper and lower bounds involve unknown constants. We will discuss the practical guidance in choosing τ in Section 4.

In Theorem 9, we assume that the input d should be the true dimension. This is a seemingly strong condition. We would like to comment on this from a few different angles.

• In the context of stochastic block models, which are simpler than the RDPG models, the parameter d is a lower bound on the number of communities. To estimate the number of communities in a stochastic block model is yet open, despite a tremendous amount of efforts (e.g. Bickel and Sarkar, 2016; Lei, 2016; Chen and Lei, 2018; Li et al., 2016; Franco Saldaña et al., 2017). We do not intend to propose a method to estimate the dimensionality here, but in practice, one could resort to the aforementioned papers.

- Without a theoretically-justified method to estimate d, we need to discuss on the potential misspecification. If one overestimates d, i.e. with an input $d_1 > d$, then our method can still consistently estimate the change points under Assumption 2, with a sufficiently large constant C_{SNR} . This is due to the fact our statistic is a function of inner products of latent position estimators. Overestimating d will only add extra noise which is in fact of the same order of the noise introduced when estimating the latent positions with true dimension d.
- Another possible misspecification is underestimating the dimension d, i.e. the input of the algorithms is $d_2 < d$. This is a more damning issue than overestimating d, however it does not necessarily lead to inconsistent change point estimators. Now we assume a toy example where the true dimension d = 3. Recall the definition on the jump size κ that

$$\kappa = \min_{k=1,...,K} \sup_{z \in [0,1]} |G_{\eta_k}(z) - G_{\eta_{k-1}}(z)|$$

=
$$\min_{k=1,...,K} \sup_{z \in [0,1]} |\mathbb{P}_{\eta_k} \left\{ X^\top Y \le z \right\} - \mathbb{P}_{\eta_{k-1}} \left\{ X^\top Y \le z \right\} |$$

=
$$\min_{k=1,...,K} \sup_{z \in [0,1]} \left| \mathbb{P}_{\eta_k} \left\{ \sum_{i=1}^3 X_i Y_i \le z \right\} - \mathbb{P}_{\eta_{k-1}} \left\{ \sum_{i=1}^3 X_i Y_i \le z \right\} \right|.$$

If we underestimate d and we miss out the third dimension, our *de facto* jump size becomes

$$\kappa_{1} = \min_{k=1,\dots,K} \sup_{z \in [0,1]} \left| \mathbb{P}_{\eta_{k}} \left\{ \sum_{i=1}^{2} X_{i} Y_{i} \le z \right\} - \mathbb{P}_{\eta_{k-1}} \left\{ \sum_{i=1}^{2} X_{i} Y_{i} \le z \right\} \right|.$$

Provided that the signal-to-noise ratio condition holds for κ_1 , i.e.

$$\kappa_1 \sqrt{\Delta n(1-\rho)} > C_{\text{SNR}} \sqrt{T} \max\{\sqrt{\log(n \vee T)}, \ d^{3/2}\} + a_T,$$

with the notation defined in Assumption 2, Theorem 9 still holds. In general, underestimating the dimension d decreases the true jump sizes at the change points. Identical arguments to those in Theorem 9 can lead to consistent detection of change points, whose decreased jump sizes satisfy the signal-to-noise ratio condition Assumption 2.

On a different note, without assuming (7), and using the trivial bound $C_R \leq T/\Delta$, it can be shown that we will achieve a larger localization error. The resulting rate inflates that of Theorem 9 by a factor of polynomials of T/Δ .

3.3 Possible extensions

There are three aspects of the methods proposed in Section 2.2 that might not seem to be satisfactory at first sight. In this subsection, we discuss possible extensions. Readers who are not familiar with the area, may safely skip this subsection during the first time reading.

From dense to sparse networks

The networks we are dealing with in this paper are dense, i.e. the average degrees are of order of the network size. In order to allow for sparse networks, one might wish to replace (1) in Definition 1 with the following

$$\mathbb{P}\left\{A \mid X\right\} = \prod_{1 \le i < j \le n} (\alpha X_i^\top X_j)^{A_{ij}} (1 - \alpha X_i^\top X_j)^{1 - A_{ij}},$$

where $\alpha = \alpha(n) \in (0, 1]$.

If α is known, then one could simply replace the definition of \widehat{X} in Definition 5 with

$$\widehat{X}(t) = \alpha^{-1/2} U_A(t) \Lambda_A(t)^{1/2}.$$

The signal-to-noise ratio and the localization errors will change correspondingly by a polynomial factor of α , following the identical derivations.

If α is unknown but satisfies $\alpha n \gtrsim \log(n)$, then one could use graphon estimation methods, e.g. the universal singular value thresholding (USVT) method (Chatterjee et al., 2015), to first produce a USVT estimator of each P(t), namely $\widehat{A}(t)$. The quantity $\widehat{X}(t)$ can be defined to be

$$\widehat{X}(t) = U_{\widehat{A}}(t)\Lambda_{\widehat{A}}(t)^{1/2}$$

and the rest of the algorithm remains the same. The localization rate would change from

$$\frac{T \max\{d \log(n \lor T), d^3\}}{\kappa^2 n(1-\rho)} \quad \text{to} \quad \frac{T \max\{\sqrt{n/\alpha}, d^3\}}{\kappa_k^2 n(1-\rho)},$$

by simply using Theorem 1 in Xu (2018) instead of Lemma 11 in controlling large probability events and following all the rest of our proofs. The term $d \log(n \vee T)$ in the upper bound is due to the fact that conditional on the latent positions, the entries in the upper triangular matrix of A(t)'s are independent. This is not true for the USVT estimator \hat{A} , and therefore the difference between the two different rates is not merely multiplying a polynomial factor of the sparsity parameter α .

From individual estimation to a bulk estimation

In Algorithm 2, we estimate every individual network separately, which results in the polynomial dependence on T in both the signal-to-noise ratio and the localization rate. One natural question would be if there is a way to conduct Algorithm 1 to a bulk of adjacency matrices at once in order to improve the statistical accuracy and computational efficiency.

There are two possible extensions. One is to use the omnibus embedding proposed in Levin et al. (2017) and the other is to conduct Algorithm 1 to a sample average of the adjacency matrices. Either way is suffered from the lack of some critical theoretical understanding. When the bulk of adjacency matrices used to construct either the omnibus matrix or the sample average matrix, are not generated from the same set of latent positions, the behaviours of the sample eigenvectors remain unknown. In change point detection, one needs to deal with intervals containing adjacency matrices coming from different latent positions. Without knowing how the eigenvectors would behave, eigenvector-based change points detection methods would not be able utilize bulk of adjacency matrices. In fact, this is also the reason that the methods proposed in Cribben and Yu (2017) and Liu et al. (2018) lack theoretical guarantees.

From using n/2 edges to all edges

In Algorithm 2, we only use n/2 out of n(n-1)/2 edges for technical convenience. In our choice, all the edges are conditionally independent given the latent positions and the concentration inequalities are easier to handle.

One extreme is to use all possible edges such that $D_{s,e}^t(z)$ defined in Definition 5 is a *U*-statistic. To be specific, (4) is replaced by

$$D_{s,e}^{t}(z) = \left| \sqrt{\frac{2(e-t)}{n(n-1)(e-s)(t-s)}} \sum_{k=s+1}^{t} \sum_{i=1}^{n-1} \sum_{j=i+1}^{n} \mathbbm{1}\{\widehat{Y}_{ij}^{k} \le z\} - \sqrt{\frac{2(t-s)}{n(n-1)(e-s)(e-t)}} \sum_{k=t+1}^{e} \sum_{i=1}^{n-1} \sum_{j=i+1}^{n} \mathbbm{1}\{\widehat{Y}_{ij}^{k} \le z\} \right|.$$

Using the Hoeffding theorem (Theorem 5.2 in Hoeffding, 1948), we can see that in our cases, the variance of using all edges and that of only using n/2 edges are of the same order. This means that using all edges will not improve the statistical accuracy (in terms of rates) but creates extra computational burden.

Generalised random dot product graph

The random dot product graph models have a generalisation, namely generalised random dot product graph models (GRDPG, Rubin-Delanchy et al., 2017). Recalling Model 1, GRDPG assumes that

$$\mathbb{P}\{A(t) \mid X(t)\} = \prod_{1 \le i < j \le n} (X_i(t)^\top I_{p,q} X_j(t))^{A_{ij}(t)} (1 - X_i(t)^\top I_{p,q} X_j(t))^{1 - A_{ij}(t)}$$

where $I_{p,q} = \text{diag}(1, \ldots, 1, -1, \ldots, -1)$ with p ones and q minus ones. We remark that the algorithms and theoretical results developed in this paper for RDPGs also hold for GRDPGs.

4. Numerical Experiments

4.1 Simulations

We now assess the performance of our proposed estimator NonPar-RDPG-CPD (Algorithm 2) in different scenarios, and compare our results with those produced by the network binary segmentation (NBS) algorithm (Wang et al., 2018a) and the modified neighbourhood smoothing (MNBS) algorithm (Zhao et al., 2019)². The measurements we adopt are the absolute error $|\hat{K} - K|$, where \hat{K} and K are the numbers of the change point estimators and the true change points, respectively, and the one-sided Hausdorff distance defined as

$$d(\widehat{\mathcal{C}}|\mathcal{C}) = \max_{\eta \in \mathcal{C}} \min_{x \in \widehat{\mathcal{C}}} |x - \eta|,$$

where \mathcal{C} is the set of true change points, and $\widehat{\mathcal{C}}$ is the set of estimated change points. We also consider the metric $d(\mathcal{C}|\widehat{\mathcal{C}})$. For Hausdorff distances, we report the medians over 100

^{2.} Code implementing our method can be found in https://github.com/hernanmp/RDPG. The algorithms are now included in the R package changepints (Xu et al., 2021).

Monte Carlo simulations, and for $|\hat{K} - K|$, we report the means over 100 Monte Carlos trials. By convention, if $\hat{\mathcal{C}} = \emptyset$, we define $d(\hat{\mathcal{C}}|\mathcal{C}) = \infty$ and $d(\mathcal{C}|\hat{\mathcal{C}}) = -\infty$.

Choice of tuning parameters. Recall that NonPar-RDPG-CPD involves three tuning parameters: (1) the threshold τ for declaring change points, (2) the number of random intervals M and (3) the dimension of the embedding d.

(1) We choose τ based on the model selection criteria from Zou et al. (2014). To be specific, we stack all the \hat{Y}_{ij}^t into one matrix $\hat{Z} \in \mathbb{R}^{T \times n/2}$. For any given τ , we let $\{\hat{\eta}_k, k = 0, \dots, \hat{K}(\tau) + 1\}$ be the set of change points estimated by NonPar-RDPG-CPD, with $\hat{\eta}_0(\tau) = 0$ and $\hat{\eta}_{\hat{K}(\tau)+1}(\tau) = T$. Define

$$BIC_{\tau} = \sum_{l=1}^{n/2} \left[\sum_{k=0}^{\hat{K}(\tau)} \sum_{t=2}^{T-1} \{ \hat{\eta}_{k+1}(\tau) - \hat{\eta}_{k}(\tau) \} \frac{\widehat{H}_{ktl} \log(\widehat{H}_{ktl}) + (1 - \widehat{H}_{ktl}) \log(1 - \hat{H}_{ktl})}{t(T - t)} + \xi \widehat{K}(\tau) \right],$$

where $\widehat{H}_{ktl} = \widehat{H}_{\widehat{\eta}_k(\tau):\widehat{\eta}_{k+1}(\tau)}^l(\widehat{Z}_{t,l})$ with $\widehat{H}_{\widehat{\eta}_k(\tau):\widehat{\eta}_{k+1}(\tau)}^l$ being the empirical cumulative distribution function of the observations $\{\widehat{Z}_{t,l}\}_{\widehat{\eta}_k(\tau)\leq t\leq \widehat{\eta}_{k+1}(\tau)-1}$ and $\xi = \log^{2.1}(n)/5$. The metric $\operatorname{BIC}_{\tau}$ is constructed by calculating along each column of \widehat{Z} the BIC-type scores defined in Equation (2.4) in Zou et al. (2014), and then aggregating the scores to produce $\operatorname{BIC}_{\tau}$. We select the model with τ that minimizes $\operatorname{BIC}_{\tau}$.

- (2) As for input representing the dimension of the latent positions d, we set d = 10 throughout this section, with varying d scenarios discussed in Appendix E. In general, we find the procedure very robust to the choice of d provided it is no smaller than the true dimension of the latent positions. This supports our discussions on misspecification after Theorem 9.
- (3) We also set M = 120, which is large enough for the various settings considered to perform well.

As for the competitor NBS, we follow the proposal by the authors in Wang et al. (2018a) setting τ to be of order $n \log^2(T)$. For the competitor MNBS, we use the default choice of its tuning parameters with code generously provided by the authors of Zhao et al. (2019). To be specific, the scaling parameter for the threshold is set as $D_0 = 0.25$, the constant B_0 for the neighborhood size is chosen as $B_0 = 3$, the threshold size is set as $\delta_0 = 0.1$ and the local window size is set as $h = \sqrt{T}$.

Disclaimer: We would like to emphasize that the comparisons to the competitors might not be fair, due to the fact that the tuning parameter choosing schemes in Zhao et al. (2019) and Wang et al. (2018a) are not meant for dependent networks.

We construct four different models, in each of which, T = 150 and K = 2. The locations of the change points are evenly spaced, giving rise to three disjoint intervals $\mathcal{A}_1 = [1, 50], \mathcal{A}_2 = [51, 100]$ and $\mathcal{A}_3 = [101, 150]$. As for the sizes of networks, we consider $n \in \{100, 200, 300\}$.

Scenario 1. Stochastic block models. We construct two matrices of probabilities, $P, Q \in \mathbb{R}^{n \times n}$. The matrix P satisfies

$$P_{i,j} = \begin{cases} 0.5, & i, j \in \mathcal{B}_l, \ l \in \{1, \dots, 4\}, \\ 0.3, & \text{otherwise}, \end{cases}$$

where $\mathcal{B}_1, \ldots, \mathcal{B}_4$ are evenly sized communities of nodes that form a partition of $\{1, \ldots, n\}$. The matrix Q satisfies

$$Q_{i,j} = \begin{cases} 0.45, & i, j \in \mathcal{B}_l, \ l \in \{1, \dots, 4\}, \\ 0.2, & \text{otherwise.} \end{cases}$$

We then construct a sequence of matrices $\{E(t)\}_{t=1}^T \subset \mathbb{R}^{n \times n}$ such that

$$E_{i,j}(t) = \begin{cases} P_{i,j}, & t \in \mathcal{A}_1 \cup \mathcal{A}_3, \\ Q_{i,j}, & \text{otherwise,} \end{cases}$$

for every $i, j \in \{1, ..., n\}$.

The data are then generated with a correlation parameter $\rho \in \{0, 0.5, 0.9\}$. Specifically, for any ρ , we have $A_{i,j}(1) \sim \text{Ber}(P_{i,j}(1))$, and between two consecutive change points,

$$A_{i,j}(t+1) \sim \begin{cases} \operatorname{Ber}((1 - E_{i,j}(t+1))\rho + E_{i,j}(t+1)), & A_{i,j}(t) = 1, \\ \operatorname{Ber}((E_{i,j}(t+1))(1-\rho)), & A_{i,j}(t) = 0, \end{cases}$$

for $1 \leq i < j \leq n$.

Scenario 2. We first generate

$$X_i(t) \stackrel{\text{ind}}{\sim} \text{Uniform}[0.2, 0.8], \quad i = 1, \dots, n, \ t \in \mathcal{A}_1 \cup \mathcal{A}_3.$$

Then for any $\varepsilon \in \{0.05, 0.15, 0.3\}$, we generate

$$X_i(t) = \begin{cases} Z_i(t) + 0.2, & i \in \{1, \dots, \lfloor n\varepsilon \rfloor\}, \\ Z_i(t), & \text{otherwise,} \end{cases}$$

where $Z_i(t) \stackrel{\text{ind}}{\sim} \text{Uniform}[0.2, 0.8]$ for $i \in \{1, \ldots, n\}$ and $t \in \mathcal{A}_2$. Then we generate $A_{i,j}(t) \sim \text{Ber}(X_i(t)X_j(t))$.

Scenario 3. For $t \in \{1, 101\}$, we generate $Z_i(t) \stackrel{\text{ind}}{\sim} \mathcal{N}(0, I_3)$, and for $t \in \mathcal{A}_1 \cup \mathcal{A}_3 \setminus \{1, 101\}$, we generate

$$Z_i(t) \begin{cases} \stackrel{\text{ind}}{\sim} \mathcal{N}(0, I_3), & \text{with probability } 0.9, \\ = Z_i(t-1), & \text{with probability } 0.1. \end{cases}$$

We then set

$$P_{i,j}(t) = \frac{\exp\{Z_i(t)^{\top} Z_j(t)\}}{1 + \exp\{Z_i(t)^{\top} Z_j(t)\}}$$

Furthermore, we generate $P_{i,j}(51) \sim \text{Beta}(100, 100)$, and for $t \in \{52, \ldots, 100\}$ we generate

$$P(t) \begin{cases} = P(t-1), & \text{with probability 0.9,} \\ \sim \text{Beta}(100, 100), & \text{with probability 0.1.} \end{cases}$$

Once the mean matrices $\{P(t)\}_{t=1}^T \mathbb{R}^{n \times n}$ have been constructed, we independently draw $A_{i,j}(t) \sim \text{Ber}(P_{i,j}(t))$, for all $i, j \in \{1, \ldots, n\}$ and $t \in \{1, \ldots, T\}$.

Scenario 4. For $t \in \{1, 101\}$ we generate $X_t \in \mathbb{R}^5$ as

$$X_i(t) \sim \text{Dirichlet}(1, 1, 1, 1, 1),$$

for all $i \in \{1, \ldots, n\}$. Then for $t \in \mathcal{A}_1 \cup \mathcal{A}_3 \setminus \{1, 101\}$,

 $X_i(t) \begin{cases} = X_i(t-1), & \text{with probability 0.9,} \\ \sim \text{Dirichlet}(1, 1, 1, 1, 1) & \text{otherwise,} \end{cases}$

for all $i \in \{1, \ldots, n\}$. We also have

$$X_i(51) \sim \begin{cases} \text{Dirichlet}(500, 500, 500, 500, 500), & i \in \{1, \dots, \lfloor n\varepsilon \rfloor\},\\ \text{Dirichlet}(1, 1, 1, 1, 1), & i \in \{\lfloor n\varepsilon \rfloor + 1, \dots, n\}, \end{cases}$$

and for $t \in \mathcal{A}_2 \setminus \{51\}$,

$$X_{i}(t) \begin{cases} = X_{i}(t-1), & \text{with probability 0.9,} \\ \sim \text{Dirichlet}(500, 500, 500, 500, 500), & \text{with probability 0.1 if } i \in \{1, \dots, \lfloor n\varepsilon \rfloor\}, \\ \sim \text{Dirichlet}(1, 1, 1, 1, 1), & \text{with probability 0.1, if } i \in \{\lfloor n\varepsilon \rfloor + 1, \dots, n\}, \end{cases}$$

for all $i \in \{1, \ldots, n\}$, where $\varepsilon \in \{0.05, 0.15, 0.3\}$.

Examples of matrices A(t) generated in each scenario are depicted in Figures 2-3. We can see qualitative differences among **Scenarios 1-4**. In particular, **Scenario 1** produces adjacency matrices with block structure. Interpretation is less clear for the other models, but we see that **Scenario 3** seems to generate more dense graphs than **Scenarios 2** and **4**.

Results comparing NonPar-RDPG-CPD with NBS are provided in Tables 1-4. We observe that, overall, NonPar-RDPG-CPD provides generally reliable estimation of the number of change points and their locations.

In **Scenario 1** with $\rho = 0$, a model where the marginal distributions of A(t) only change in mean, we see from Table 1 that NBS outperforms our proposed approach. This does not come as a surprise since NBS is designed to detect change points in mean. However, as ρ increases and the number of samples decreases, the most robust method seems to be NonPar-RDPG-CPD.

Scenario 2 poses an interesting example where the behaviour of only a fraction of nodes in the network changes at the change points. Furthermore, the data are generated under an RDPG model. As shown in Table 2, NonPar-RDPG-CPD seems to be the best method for estimating the number of change points. A possible explanation is that the underlying

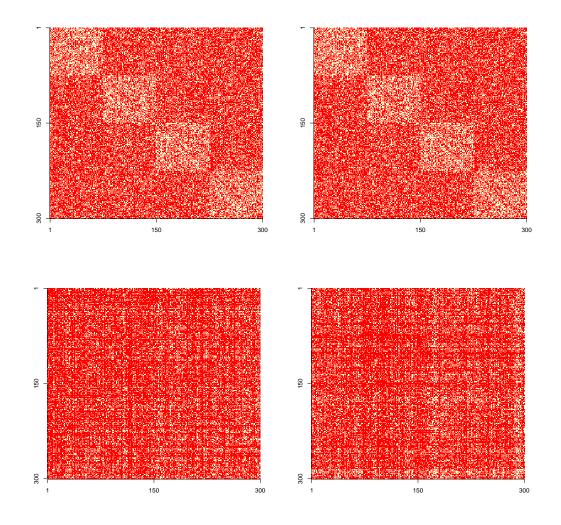


Figure 2: The top row shows two instances of data generated in **Scenario 1**. The left panel corresponds to A(t) for t before the first change point, and the right panel to A(t) between the first and second change points. The bottom row shows the corresponding plots for **Scenario 2** with $\varepsilon = 0.05$.

changes in the distributions of A(t) not only occur at the level of the means, and hence the NBS might not be the ideal for this scenario even though it outperforms MNBS in this framework. Our method was constructed under the assumption of the RDPG model.

To assess the robustness of our method to misspecification, we can look at the performance of our method in the context of **Scenario 3** which is not an RDPG. Interestingly, Table 3 shows that NonPar-RDPG-CPD is the best in this model with MNBS coming in second. In contrast, NBS suffers greatly, overestimating the number of change points. This makes sense since between change points, the latent positions X(t) remain constant with

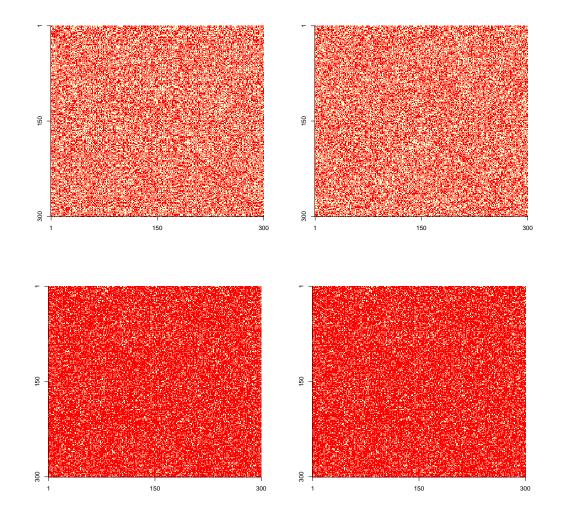


Figure 3: The top row shows two instances of data generated in **Scenario 3**. The left panel corresponds to A(t) for t before the first change point, and the right panel to A(t) between the first and second change points. The bottom row shows the corresponding plots for **Scenario 4** with $\varepsilon = 0.05$.

probability 0.9 and change with probability 0.1. Hence, some of these changes in X(t) could be confused as change points by NBS.

Finally, Scenario 4 consists of an example of Model 1. However, similarly as Scenario 2, the change points correspond to shifts in the behaviour of only some of the nodes in the network. In particular, Table 4 suggests that our method performs reasonably well, improving its performance when the signal-to-noise ratio increases. This is different from the NBS which once again tends to overestimate the number of change points. As for the MNBS, we see that this method is unable to detect the change points in this example.

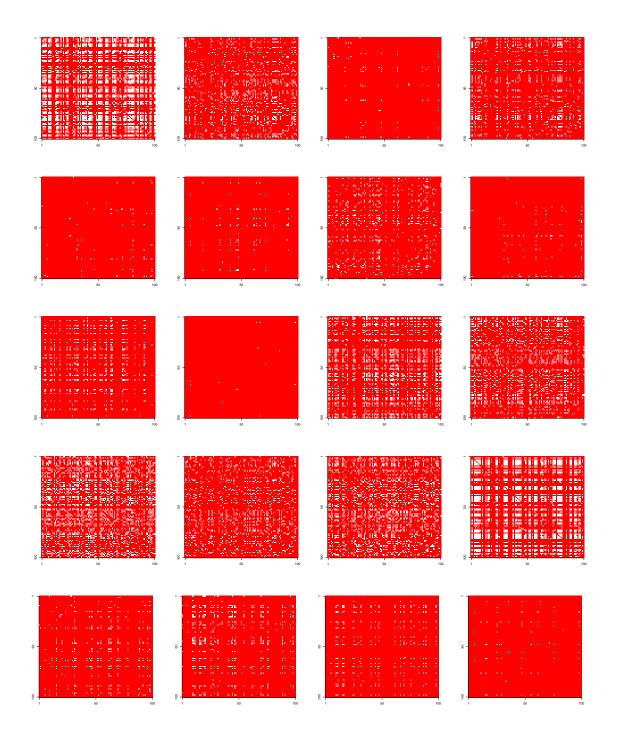


Figure 4: Examples of adjacency matrices, down-sampled to a 100×100 , between the change points estimated by NonPar-RDPG-CPD in the zebrafish example. From left to right and from top to bottom, the first two rows of panels correspond to t = 3, 7, 15, 32, 40, 45, 52, 60, 65, 75, 80 and 87. From left to right and from top to bottom, the last two rows correspond to t = 6, 8, 9, 10, 11, 12 and 13.

				$1(\widehat{a} a)$	$1(a \hat{a})$
Method	<u>n</u>	ρ	K-K	$\frac{d(\widehat{\mathcal{C}} \mathcal{C})}{1}$	$\frac{d(\mathcal{C} \widehat{\mathcal{C}})}{1-2}$
NonPar-RDPG-CPD	300	0	0.1	1.0	1.0
NBS	300	0	0.0	1.0	1.0
MNBS	300	0	1.16	50.0	0.0
NonPar-RDPG-CPD	200	0	0.0	1.0	1.0
NBS	200	0	0.0	1.0	1.0
MNBS	200	0	1.92	\inf	$-\inf$
NonPar-RDPG-CPD	100	0	0.2	1.0	1.0
NBS	100	0	0.0	1.0	1.0
MNBS	100	0	0.84	50.0	0.0
NonPar-RDPG-CPD	300	0.5	0.0	0.0	0.0
NBS	300	0.5	21.2	1.0	43.0
MNBS	300	0.5	0.0	0.0	0.0
NonPar-RDPG-CPD	200	0.5	0.04	0.0	0.0
NBS	200	0.5	21.3	1.0	4.30
MNBS	200	0.5	0.0	0.0	0.0
NonPar-RDPG-CPD	100	0.5	0.16	0.0	0.0
NBS	100	0.5	21.3	1.0	42.0
MNBS	100	0.5	0.12	0.0	0.0
NonPar-RDPG-CPD	300	0.9	0.0	0.0	0.0
NBS	300	0.9	21.0	1.0	43.0
MNBS	300	0.9	3.12	0.0	36.0
NonPar-RDPG-CPD	200	0.9	0.0	0.0	0.0
NBS	200	0.9	21.0	1.0	43.0
MNBS	200	0.9	2.88	0.0	35.0
NonPar-RDPG-CPD	100	0.9	0.0	1.0	1.0
NBS	100	0.9	21.04	1.0	43.0
MNBS	100	0.9	3.28	0.0	35.0

Table 1: Scenario 1

4.2 Real data

Our goal is to estimate change points in the context of the neuronal activity in larval zebrafish. The data consist of simultaneous whole-brain neuronal activity data at near single cell resolution (Prevedel et al., 2014). The original data format is a matrix of size 5379×5000 . This corresponds to the neural activity of 5379 neurons over 5000 frames, where one second in time corresponds to 20 frames.

To construct the final sequence of networks, we proceed as in Park et al.. Specifically, we first remove artificial neurons leaving us with a 5105×5000 matrix. Then we bin the data into 100 non-overlapping periods. Each period corresponds to 2.5 seconds of the original data. The resulting time series is then $Z(t) \in \mathbb{R}^{5105 \times 50}$ for $t \in \{1, \ldots, 100\}$. Following Lyzinski et al. (2017), we finally construct the adjacency matrices $A(t) \in \mathbb{R}^{5105 \times 5105}$ as

$$A_{i,j}(t) = \mathbb{1}\{\operatorname{corr}(Z_i(t), Z_j(t)) > 0.7\}, \quad t = 1, \dots, T,$$

where T = 100.

With the time series $\{A(t)\}_{t=1}^{T}$ in hand, we proceed to run change point detection with Algorithm 2. The implementation details are the same as those in Section 4.1. However,

Method	n	ε	$ K - \widehat{K} $	$d(\widehat{\mathcal{C}} \mathcal{C})$	$d(\mathcal{C} \widehat{\mathcal{C}})$
NonPar-RDPG-CPD	300	0.3	0.04	0.0	0.0
NBS	300	0.3	0.28	1.0	1.0
MNBS	300	0.3	0.76	0.0	21.0
NonPar-RDPG-CPD	200	0.3	0.0	0.0	0.0
NBS	200	0.3	0.32	1.0	1.0
MNBS	200	0.3	0.48	0.0	1.0
NonPar-RDPG-CPD	100	0.3	0.08	3.0	3.0
NBS	100	0.3	0.08	1.0	1.0
MNBS	100	0.3	0.64	0.0	18.0
NonPar-RDPG-CPD	300	0.15	0.0	2.0	2.0
NBS	300	0.15	0.4	1.0	1.0
MNBS	300	0.15	0.76	0.0	21.0
NonPar-RDPG-CPD	200	0.15	0.04	3.0	3.0
NBS	200	0.15	0.28	1.0	1.0
MNBS	200	0.15	0.76	0.0	20.0
NonPar-RDPG-CPD	100	0.15	0.28	4.0	10.0
NBS	100	0.15	0.32	1.0	1.0
MNBS	100	0.15	0.48	1.0	5.0
NonPar-RDPG-CPD	300	0.05	0.72	36.0	5.0
NBS	300	0.05	0.84	1.0	9.0
MNBS	300	0.05	1.24	1.0	21.0
NonPar-RDPG-CPD	200	0.05	0.64	37.0	6.0
NBS	200	0.05	0.76	3.0	11.0
MNBS	200	0.05	0.6	4.0	8.0
NonPar-RDPG-CPD	100	0.05	0.72	19.0	15.0
NBS	100	0.05	1.4	\inf	$-\inf$
MNBS	100	0.05	1.88	\inf	$-\inf$

Table 2:Scenario 2

Table 3: Scenario 3

Method	n	$ K - \widehat{K} $	$d(\widehat{\mathcal{C}} \mathcal{C})$	$d(\mathcal{C} \widehat{\mathcal{C}})$
NonPar-RDPG-CPD	300	0.24	0.0	0.0
NBS	300	15.04	1.0	43.0
MNBS	300	0.84	25	36
NonPar-RDPG-CPD	200	0.08	0.0	0.0
NBS	200	14.4	43.0	1.0
MNBS	200	0.84	23	36
NonPar-RDPG-CPD	100	0.52	3.0	5.0
NBS	100	13.96	1.0	43.0
MNBS	100	1.16	23	35

to facilitate computations at every instance of time we randomly sample 800 nodes in the network and work with a down-sampled version of A(t). After running our method, we estimate change points at t = 5, 10, 29, 36, 42, 50, 57, 62, 71, 79, 85, and 89. In the original 250 seconds time stamp, the changes correspond to 12.5 25.0, 72.5, 90.0, 105.0, 125.0, 142.5, 155.0, 177.5, 197.5, 212.5, and 222.5 seconds. Simple inspection suggests that our estimated

Method	n	ε	$ K - \widehat{K} $	$d(\widehat{\mathcal{C}} \mathcal{C})$	$d(\mathcal{C} \widehat{\mathcal{C}})$
NonPar-RDPG-CPD	300	0.3	0.72	35.0	12.0
NBS	300	0.3	19.4	1.0	43.0
MNBS	300	0.3	2.0	\inf	$-\inf$
NonPar-RDPG-CPD	200	0.3	0.84	40.0	10.0
NBS	200	0.3	19.4	1.0	43.0
MNBS	200	0.3	2.0	\inf	$-\inf$
NonPar-RDPG-CPD	100	0.3	1.0	30.0	20.0
NBS	100	0.3	9.44	3.0	41.0
MNBS	100	0.3	2.0	\inf	$-\inf$
NonPar-RDPG-CPD	300	0.15	0.8	34.0	17.0
NBS	300	0.15	20.24	1.0	43.0
MNBS	300	0.15	2.0	\inf	$-\inf$
NonPar-RDPG-CPD	200	0.15	0.96	40.0	11.0
NBS	200	0.15	17.0	1.0	43.0
MNBS	200	0.15	2.0	\inf	$-\inf$
NonPar-RDPG-CPD	100	0.15	0.84	34	18.0
NBS	100	0.15	10.64	1.0	41.0
MNBS	100	0.15	2.0	\inf	$-\inf$
NonPar-RDPG-CPD	300	0.05	0.80	33.0	17.0
NBS	300	0.05	20.48	1.0	43.0
MNBS	300	0.05	2.0	\inf	$-\inf$
NonPar-RDPG-CPD	200	0.05	0.88	38.0	19.0
NBS	200	0.05	17.56	1.0	43.0
MNBS	200	0.05	2.0	\inf	$-\inf$
NonPar-RDPG-CPD	100	0.05	1.04	32.0	16.0
NBS	100	0.05	11.48	3.0	41.0
MNBS	100	0.05	2.0	\inf	$-\inf$

Table 4: Scenario 4

change points are in agreement with the extracted intensity signal of Ca2+ fluorescence using spatial filters in Figure 3 (c) in Prevedel et al. (2014). As remarked in Park et al., a lab scientist induced a change-point at the 16th second, by giving an olfactory stimulus to the zebrafish. In the scale of our time series $\{A(t)\}_{t=1}^{T}$, this change corresponds to t = 6which seems to be captured by our algorithm that detected a change point at t = 5.

We also considered change point detection with the algorithm NBS (Wang et al., 2018a). The set of estimated change points is roughly the same to that estimated by NonPar-RDPG-CPD: 10, 14, 22, 26, 32, 36, 42, 50, 58, 62, 66, 72, 80, and 90. One important difference, however, is that NBS did not detect a change point near t = 6, the change point created by the lab scientist. We also tried the MNBS method (Zhao et al., 2019), but this only detected changes at 14, 45, 66, 80. Tuning parameters of both NBS and MNBS are chosen as described in Section 4.1.

Finally, we have included Figure 4 which shows down-sampled versions of A(t) for values of t between estimated change points. This reinforces our intuition that the structural breaks estimated with NonPar-RDPG-CPD are meaningful.

5. Discussions

In this paper, we have studied the offline change point localization problem in a sequence of dependent nonparametric random dot product graphs. We allow for a weakly dependent process along the time and introduce the dependence within networks via latent positions. In fact, conditional on the latent positions, the edges within a network are independent and one may wish to further allow for dependence among edges conditional on the latent positions. We remark that this is technically feasible - one can incorporate a weak dependence version of Bernstein's inequality (Lemma 14) in the estimation of the latent positions (Lemma 11). Such deployment requires a data generating mechanism characterizing the dependence among edges, but without a natural distance among edges, we refrain our pursuit on this direction.

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Appendix A. Technical details of Section 3.1

Proof [Proof of Lemma 7] For any $i, j \in \{1, ..., n\}, i \neq j$, it holds that

$$\mathbb{P}\{A_{ij}|X_i, X_j\} = X_i^\top X_j = X_i^\top U^\top U X_j,$$

for any orthogonal operator $U \in \mathbb{R}^{d \times d}$. In this proof, by the equivalence in terms of the distributions F and \widetilde{F} , we mean the equivalence up to a rotation, which is detailed in Definition 3. Without loss of generality, if a rotation is needed, we omit it in the notation.

We divide this proof into two cases: (a) d = 1 and (b) d > 1.

(a) p = 1.

Since the entries of A and \widetilde{A} are Bernoulli random variables, they only take values in $\{0,1\}^{n\times n}$. For any symmetric matrix $v \in \{0,1\}^{n\times n}$, we have

$$\mathbb{P}\{A = v\} = \mathbb{E}\left\{\mathbb{E}\left(\prod_{i=1}^{n-1} \prod_{j=i+1}^{n} \mathbb{1}\{A_{ij} = v_{ij}\} \middle| \{X_l\}_{l=1}^n\right)\right\}$$
$$= \mathbb{E}\left[\prod_{i=1}^{n-1} \prod_{j=i+1}^{n} \{(X_i X_j) v_{ij} + (1 - X_i X_j)(1 - v_{ij})\}\right].$$
(9)

If $\mathcal{L} = \widetilde{\mathcal{L}}$, then we have the following.

• If $v_{ij} \equiv 1$, then

(9) =
$$\mathbb{E}\left[\prod_{i=1}^{n-1}\prod_{j=i+1}^{n} (X_i X_j)\right] = \left\{\mathbb{E}(X_1^{n-1})\right\}^n$$
,

which implies that $\mathbb{E}_F(X_1^{n-1}) = \mathbb{E}_{\widetilde{F}}(\widetilde{X}_1^{n-1})$. Note that in order to have an edge, $n \ge 2$, which implies that $n-1 \ge 1$.

• If there is one and only one pair (i, j), i < j, such that $v_{ij} = v_{ji} = 0$, and $v_{kl} = 1$, $(k, l) \notin \{(i, j), (j, i)\}$, then without loss of generality, we let (i, j) = (1, 2). If n = 2, then

$$(9) = 1 - \{\mathbb{E}(X_1)\}^2,\$$

which implies that $\mathbb{E}_F(X_1^{n-1}) = \mathbb{E}_{\widetilde{F}}(\widetilde{X}_1^{n-1}).$ If $n \geq 3$, then

$$(9) = \mathbb{E}\left[\prod_{i=3}^{n-1} \prod_{j=i+1}^{n} (X_i X_j) \cdot \prod_{r=1}^{2} \prod_{l=3}^{n} (X_r X_l) \cdot (1 - X_1 X_2)\right]$$
$$= \left\{\mathbb{E}(X_1^{n-2})\right\}^2 \left\{\mathbb{E}(X_1^{n-1})\right\}^{n-2} - \left\{\mathbb{E}(X_1^{n-1})\right\}^n,$$

which implies $\mathbb{E}_F(X_1^{n-2}) = \mathbb{E}_{\widetilde{F}}(\widetilde{X}_1^{n-2}).$

• If $n \ge 3$, then for $k \in \{2, \ldots, n-1\}$, without loss of generality, let $v_{1j} = v_{j1} = 0$, $j \in \{2, \ldots, k+1\}$, and $v_{rs} = v_{sr} = 1$ otherwise. We have that

$$(9) = \mathbb{E}\left[\prod_{i=k+2}^{n-1}\prod_{j=i+1}^{n}(X_{i}X_{j})\cdot\prod_{l=1}^{k+1}\prod_{i=k+2}^{n}(X_{l}X_{i})\cdot\prod_{r=2}^{k+1}(1-X_{1}X_{r})\right]$$
$$= \left\{\mathbb{E}(X_{1}^{n-1})\right\}^{n-k-1}\mathbb{E}\left[\prod_{l=1}^{k+1}X_{l}^{n-k-1}\cdot\prod_{r=2}^{k+1}(1-X_{1}X_{r})\right]$$
$$= \left\{\mathbb{E}(X_{1}^{n-1})\right\}^{n-k-1}\sum_{r=0}^{k}\binom{k}{r}(-1)^{r}\mathbb{E}(X_{1}^{n-k-1+r})\left[\mathbb{E}(X_{1}^{n-k})\right]^{r}.$$
(10)

Note that, if k = 2, then the summands in (10) include moments n-1, n-2 and n-3. We have already shown that $\mathbb{E}_F(X_1^{n-1}) = \mathbb{E}_{\widetilde{F}}(\widetilde{X}_1^{n-1})$ and $\mathbb{E}_F(X_1^{n-2}) = \mathbb{E}_{\widetilde{F}}(\widetilde{X}_1^{n-2})$, therefore (10) implies that $\mathbb{E}_F(X_1^{n-3}) = \mathbb{E}_{\widetilde{F}}(\widetilde{X}_1^{n-3})$.

• By induction, for $n > k_0$ and $k_0 \ge 3$, if it holds that $\mathbb{E}_F(X_1^{n-s}) = \mathbb{E}_{\widetilde{F}}(\widetilde{X}_1^{n-s})$, $s = 1, \ldots, k_0$, then we have $\mathbb{E}_F(X_1^{n-k_0-1}) = \mathbb{E}_{\widetilde{F}}(\widetilde{X}_1^{n-k_0-1})$, due to the fact that the summands in (10) include moment $n-s, s = 1, \ldots, k_0 + 1$.

We conclude that if $\mathcal{L} = \widetilde{\mathcal{L}}$, then $\mathbb{E}_F(X_1^k) = \mathbb{E}_{\widetilde{F}}(\widetilde{X}_1^k), k = 1, \dots, n-1$.

If $\mathbb{E}_F(X_1^k) = \mathbb{E}_{\widetilde{F}}(\widetilde{X}_1^k)$, $k = 1, \ldots, n-1$, then it follows from that for any v,

$$(9) = \sum_{l=0}^{\sum_{i$$

which is a function solely of $\mathbb{E}_F(X_1^k)$, k = 1, ..., n - 1. We, therefore, have that $\mathcal{L} = \widetilde{\mathcal{L}}$. (b) d > 1.

Since the entries of A and \tilde{A} are Bernoulli random variables, they only take values in $\{0,1\}^{n\times n}$. For any symmetric matrix $v \in \{0,1\}^{n\times n}$, we have

$$\mathbb{P}\{A=v\} = \mathbb{E}\left\{\mathbb{E}\left(\prod_{i=1}^{n-1}\prod_{j=i+1}^{n}\mathbb{1}\{A_{ij}=v_{ij}\}\Big|\{X_l\}_{l=1}^n\right)\right\}$$
$$= \mathbb{E}\left[\prod_{i=1}^{n-1}\prod_{j=i+1}^{n}\left\{\left(\sum_{k=1}^d X_{i,k}X_{j,k}\right)v_{ij} + \left(1 - \sum_{k=1}^d X_{i,k}X_{j,k}\right)(1 - v_{ij})\right\}\right].$$
(11)

If $\mathcal{L} = \widetilde{\mathcal{L}}$, then we have the following.

• If $v_{ij} \equiv 1$, then

$$(11) = \mathbb{E}\left[\prod_{i=1}^{n-1} \prod_{j=i+1}^{n} \left(\sum_{k=1}^{d} X_{i,k} X_{j,k}\right)\right] \\ = \mathbb{E}\left[\prod_{j=2}^{n} \left(\sum_{k=1}^{p} X_{1,k} X_{j,k}\right) \cdot \prod_{i=2}^{n-1} \prod_{j=i+1}^{n} \left(\sum_{k=1}^{d} X_{i,k} X_{j,k}\right)\right] \\ = \mathbb{E}\left\{\left[\sum_{k_{2},\dots,k_{n}=1}^{d} \left(\prod_{l=2}^{n} X_{1,k_{l}}\right) \cdot \left(\prod_{l=2}^{n} X_{l,k_{l}}\right)\right] \cdot \left[\prod_{i=2}^{n-1} \prod_{j=i+1}^{n} \left(\sum_{k=1}^{d} X_{i,k} X_{j,k}\right)\right]\right\} \\ = \sum_{k_{2},\dots,k_{n}=1}^{d} \mathbb{E}\left(\prod_{l=2}^{n} X_{1,k_{l}}\right) \mathbb{E}\left\{\left(\prod_{l=2}^{n} X_{l,k_{l}}\right) \cdot \left[\prod_{i=2}^{n-1} \prod_{j=i+1}^{n} \left(\sum_{k=1}^{d} X_{i,k} X_{j,k}\right)\right]\right\}, (12)$$

where the third identity follows from the independence assumption. Note that for any $(k_2, \ldots, k_n) \in \{1, \ldots, p\}^{\otimes (n-1)}$, the term

$$\mathbb{E}\left\{\left(\prod_{l=2}^{n} X_{l,k_l}\right) \cdot \left[\prod_{i=2}^{n-1} \prod_{j=i+1}^{n} \left(\sum_{k=1}^{d} X_{i,k} X_{j,k}\right)\right]\right\}$$

in (12) does not involve X_1 , and the term

$$\mathbb{E}\left(\prod_{l=2}^{n} X_{1,k_l}\right)$$

includes all possible terms of the form

$$\mathbb{E}\left(\prod_{l=1}^{d} X_{1,l}^{k_l}\right), \quad \sum_{l=1}^{d} k_l = n-1, \quad k_l \ge 0, \ l \in \{1, \dots, d\}.$$
(13)

Due to the exchangeablility, we conclude that (11) is solely a function of polynomials of (13).

If n = 2, then due to Definition 1, we have that $\mathcal{L} = \widetilde{\mathcal{L}}$ implies that

$$\mathbb{E}\left(\prod_{l=1}^{d} X_{1,l}^{k_l}\right) = \mathbb{E}\left(\prod_{l=1}^{d} \widetilde{X}_{1,l}^{k_l}\right), \quad \sum_{l=1}^{d} k_l = n-1, \quad k_l \ge 0, \ l \in \{1, \dots, d\}.$$

• If $n \ge 3$, then we prove by induction. Assume that

$$\mathbb{E}\left(\prod_{l=1}^{p} X_{1,l}^{k_l}\right) = \mathbb{E}\left(\prod_{l=1}^{p} \widetilde{X}_{1,l}^{k_l}\right), \quad \sum_{l=1}^{p} k_l = n - k, \dots, n - 1, \quad k_l \ge 0, \ l \in \{1, \dots, p\},$$

where $n-1 \ge n-k \ge 2$. We now proceed to prove that

$$\mathbb{E}\left(\prod_{l=1}^{p} X_{1,l}^{k_l}\right) = \mathbb{E}\left(\prod_{l=1}^{p} \widetilde{X}_{1,l}^{k_l}\right), \quad \sum_{l=1}^{p} k_l = n - k - 1, \dots, n - 1, \quad k_l \ge 0, \ l \in \{1, \dots, d\}.$$
(14)

To show this, we assume that $v_{1j} = v_{j1} = 0, j \in \{2, \ldots, k+1\}$, and $v_{rs} = 1$ otherwise. We have that

$$(11) = \mathbb{E}\left[\prod_{j=2}^{k+1} \left(1 - \sum_{s=1}^{s} X_{1,s} X_{j,s}\right) \cdot \prod_{l=k+2}^{n} \left(\sum_{s=1}^{d} X_{1,s} X_{l,s}\right) \cdot \prod_{i=2}^{n-1} \prod_{r=i+1}^{n} \left(\sum_{s=1}^{d} X_{i,s} X_{r,s}\right)\right)\right]$$
$$= (-1)^{k} \mathbb{E}\left\{\prod_{l=k+2}^{n} \left(\sum_{s=1}^{d} X_{1,s} X_{l,s}\right) \cdot \prod_{i=2}^{n-1} \prod_{r=i+1}^{n} \left(\sum_{s=1}^{d} X_{i,s} X_{r,s}\right)\right)\right\} + f(X)$$
$$= (-1)^{k} \sum_{s_{k+2},\dots,s_{n}=1}^{d} \mathbb{E}\left(\prod_{l=k+2}^{n} X_{1,s_{l}}\right) \mathbb{E}\left\{\prod_{l=k+2}^{n} \left(\sum_{s=1}^{d} X_{l,s}\right) \times \prod_{i=2}^{n-1} \prod_{r=i+1}^{n} \left(\sum_{s=1}^{d} X_{i,s} X_{r,s}\right)\right\} + f(X),$$

where f(X) is solely a function of

$$\mathbb{E}\left(\prod_{l=1}^{d} X_{1,l}^{k_l}\right), \quad \sum_{l=1}^{d} k_l = n - k, \dots, n - 1, \quad k_l \ge 0, \ l \in \{1, \dots, d\}.$$

Note that

$$\sum_{s_{k+2},\dots,s_n=1}^d \mathbb{E}\left(\prod_{l=k+2}^n X_{1,s_l}\right)$$

is a function of

$$\mathbb{E}\left(\prod_{l=1}^{d} X_{1,l}^{k_l}\right), \quad \sum_{l=1}^{d} k_l = n - k - 1, \quad k_l \ge 0, \, l \in \{1, \dots, d\}.$$

Therefore we have shown (14).

To this end, we have that $\mathcal{L} = \widetilde{\mathcal{L}}$ implies that

$$\mathbb{E}\left(\prod_{l=1}^{d} X_{1,l}^{k_l}\right) = \mathbb{E}\left(\prod_{l=1}^{d} \widetilde{X}_{1,l}^{k_l}\right), \quad \sum_{l=1}^{d} k_l = 1, \dots, n-1, \quad k_l \ge 0, \ l \in \{1, \dots, d\}.$$
(15)

To show that (15) implies that $\mathcal{L} = \widetilde{\mathcal{L}}$, we notice that for any v,

$$(11) = \sum_{l=0}^{\sum_{i$$

which is solely a function of

$$\mathbb{E}\left(\prod_{l=1}^{d} X_{1,l}^{k_l}\right), \quad \sum_{l=1}^{d} k_l = 1, \dots, n-1, \quad k_l \ge 0, \ l \in \{1, \dots, d\}.$$

The final claim holds.

Proof [Proof of Lemma 8] For simplicity, we assume n is an even number. Let $\mathcal{O} = \{(i, n/2 + i), i = 1, \dots, n/2\}$. Let

$$z_* \in \underset{z \in [0,1]}{\operatorname{argsup}} |G(z) - \widetilde{G}(z)|.$$

Note that

$$\begin{aligned} \left| \sqrt{\frac{2}{n}} \sum_{(i,j)\in\mathcal{O}} \left(\mathbbm{I}\{Y_{ij} \le z_*\} - \mathbbm{I}\{\tilde{Y}_{ij} \le z_*\} \right) \right| \\ &= \left| \sqrt{\frac{2}{n}} \sum_{(i,j)\in\mathcal{O}} \left\{ (\mathbbm{I}\{Y_{ij} \le z_*\} - \mathbbm{E}[\mathbbm{I}\{Y_{ij} \le z_*\}]) - \left(\mathbbm{I}\{\tilde{Y}_{ij} \le z_*\} - \mathbbm{E}[\mathbbm{I}\{\tilde{Y}_{ij} \le z_*\}] \right) \right\} \\ &+ \sqrt{\frac{n}{2}} \left\{ \mathbbm{E}[\mathbbm{I}\{Y_{ij} \le z_*\}] - \mathbbm{E}[\mathbbm{I}\{\tilde{Y}_{ij} \le z_*\}] \right\} \\ &\geq \sqrt{\frac{n}{2}} \left| \mathbbm{E}[\mathbbm{I}\{Y_{ij} \le z_*\}] - \mathbbm{E}[\mathbbm{I}\{\tilde{Y}_{ij} \le z_*\}] \right| - \left| \sqrt{\frac{2}{n}} \sum_{(i,j)\in\mathcal{O}} (\mathbbm{I}\{Y_{ij} \le z_*\}] \right) \right| \\ &- \left| \sqrt{\frac{2}{n}} \sum_{(i,j)\in\mathcal{O}} (\mathbbm{I}\{\tilde{Y}_{ij} \le z_*\} - \mathbbm{E}[\mathbbm{I}\{\tilde{Y}_{ij} \le z_*\}]) \right| \\ &- \left| \sqrt{\frac{2}{n}} \sum_{(i,j)\in\mathcal{O}} (\mathbbm{I}\{\tilde{Y}_{ij} \le z_*\} - \mathbbm{E}[\mathbbm{I}\{Y_{ij} \le z_*\}]) \right| \\ &- \left| \sqrt{\frac{2}{n}} \sum_{(i,j)\in\mathcal{O}} (\mathbbm{I}\{\tilde{Y}_{ij} \le z_*\} - \mathbbm{E}[\mathbbm{I}\{\tilde{Y}_{ij} \le z_*\}]) \right| . \end{aligned}$$
(16)

Next, it follows from Hoeffding's inequality that

$$\mathbb{P}\left\{ \left. \max\left\{ \left| \sqrt{\frac{2}{n}} \sum_{(i,j)\in\mathcal{O}} \left(\mathbbm{1}\left\{Y_{ij} \le z_*\right\} - \mathbb{E}\left[\mathbbm{1}\left\{Y_{ij} \le z_*\right\}\right] \right) \right|, \\ \left| \sqrt{\frac{2}{n}} \sum_{(i,j)\in\mathcal{O}} \left(\mathbbm{1}\left\{\widetilde{Y}_{ij} \le z_*\right\} - \mathbb{E}\left[\mathbbm{1}\left\{\widetilde{Y}_{ij} \le z_*\right\}\right] \right) \right| \right\} > \sqrt{\log(n)} \right\} \le 2n^{-4}.$$
(17)

Combining (16) and (17), we have that with probability at least $1 - 2n^{-4}$,

$$\left| \sqrt{\frac{2}{n}} \sum_{(i,j)\in\mathcal{O}} \left(\mathbb{1}\{Y_{ij} \le z_*\} - \mathbb{1}\{\widetilde{Y}_{ij} \le z_*\} \right) \right| \ge \kappa_0 \sqrt{n/2} - 2\sqrt{\log(n)}.$$
(18)

We then prove by contradiction. If $\mathcal{L} = \widetilde{\mathcal{L}}$, then it follows from Hoeffding's inequality that

$$\mathbb{P}\left\{ \left| \sqrt{\frac{2}{n}} \sum_{(i,j)\in\mathcal{O}} \left(\mathbb{1}\left\{ Y_{ij} \le z_* \right\} - \mathbb{1}\left\{ \widetilde{Y}_{ij} \le z_* \right\} \right) \right| \le \sqrt{\log(n)} \right\} \ge 1 - 2n^{-4}.$$
(19)

Due to Assumption 1, (18) and (19) contradict with each other, which implies that $\mathcal{L} \neq \widetilde{\mathcal{L}}$.

Appendix B. Large probability events

Define

$$\Delta_{s,e}^t(z) = \sum_{k=s+1}^e w_k \sum_{(i,j)\in\mathcal{O}} \left(\mathbb{1}\{\widehat{Y}_{ij}^k \le z\} - \mathbb{E}\left\{ \mathbb{1}\{Y_{ij}^k \le z\}\right\} \right),$$

where

$$w_{k} = \begin{cases} \sqrt{\frac{2}{n}} \sqrt{\frac{e-t}{(e-s)(t-s)}}, & k = s+1, \dots, t, \\ -\sqrt{\frac{2}{n}} \sqrt{\frac{t-s}{(e-s)(e-t)}}, & k = t+1, \dots, e. \end{cases}$$

In this section, we are to show the following two events hold with probability tending to 1, as $(n \lor T) \to \infty$,

$$\mathcal{B}_{1} = \left\{ \max_{0 \le s < t < e \le T} \Delta_{s,e}^{t} \le C_{9} \sqrt{\frac{T}{1-\rho}} \max\{d \log(n \lor T), d^{3/2} \sqrt{\log(n \lor T)}\} \right\}$$

and

$$\mathcal{B}_{2} = \left\{ \max_{0 \le s < t < e \le T} \sup_{z \in [0,1]} \left| \sqrt{\frac{2}{n(e-s)}} \sum_{k=s+1}^{e} \sum_{(i,j) \in \mathcal{O}} \left(\mathbb{1}\{\widehat{Y}_{ij}^{k} \le z\} - \mathbb{E}\left\{ \mathbb{1}\{Y_{ij}^{k} \le z\} \right\} \right) \right. \\ \left. \le C_{9} \sqrt{\frac{T}{1-\rho}} \max\{d \log(n \lor T), \ d^{3/2} \sqrt{\log(n \lor T)}\} \right\}.$$

This is formally stated in Lemma 16. To reach there, we denote

$$\mathcal{E}_{1} = \left\{ \max_{t=1,\dots,T} \| U_{P_{t}}^{\top}(A(t) - P_{t})U_{P_{t}} \|_{\mathrm{F}} \le C_{1}\sqrt{\log(n \vee T)} \right\},\$$

$$\mathcal{E}_{2} = \left\{ \max_{t=1,\dots,T} \| (A(t) - P_{t})U_{P_{t}} \|_{2 \to \infty} \le C_{2}\sqrt{d\log(n \vee T)} \right\},\$$

$$\mathcal{E}_3 = \left\{ \max_{t=1,\dots,T} \|A(t) - P_t\|_{\text{op}} \le C_3 \sqrt{n} \right\}$$

and

$$\mathcal{E}_4 = \left\{ 2^{-1} n \min_{k=1,\dots,K} \mu_d^k \le \min_{t=1,\dots,T} \lambda_d(P_t) \le \max_{t=1,\dots,T} \lambda_1(P_t) \le (3/2) n \max_{k=1,\dots,K} \mu_1^k \right\},$$

where $C_1 > 4\sqrt{6}$, $C_2 > 4\sqrt{6}$, $C_3 > 0$ are universal constants. Throughout, $\|\cdot\|_{2\to\infty}$ denotes the two-to-infinity norm. To be specific, for any matrix $M \in \mathbb{R}^{m_1 \times m_2}$,

$$||M||_{2\to\infty} = \max_{x\in\mathbb{R}^{m_2}:\,||x||_2=1} ||Ax||_{\infty},$$

where $||Ax||_{\infty}$ denotes the largest absolute value of the entries in Ax.

Lemma 10 Under Model 1, for any $t \in \{1, ..., T\}$, it holds that

$$\mathbb{P}\{\lambda_{d+1}(P_t)=0\}=1$$

Proof For any $t \in \{1, \ldots, T\}$, we have that

$$P_t = X(t)(X(t))^{\top}.$$

For any realisation of $X(t) \in \mathbb{R}^{n \times d}$, $\lambda_{d+1}(P_t) = 0$. Thus the final claim holds.

Lemma 11 Under Model 1, we have that

$$\max\left\{\mathbb{P}\left\{\mathcal{E}_{1} \mid \{X(t)\}_{t=1}^{T}\right\}, \mathbb{P}\left\{\mathcal{E}_{1}\right\}\right\} \ge 1 - (n \lor T)^{-c_{1}},$$
(20)

$$\max\left\{\mathbb{P}\left\{\mathcal{E}_{2} \mid \{X(t)\}_{t=1}^{T}\right\}, \mathbb{P}\left\{\mathcal{E}_{2}\right\}\right\} \ge 1 - (n \lor T)^{-c_{2}}$$
(21)

and

$$\max\left\{\mathbb{P}\left\{\mathcal{E}_{3} \mid \{X(t)\}_{t=1}^{T}\right\}, \mathbb{P}\left\{\mathcal{E}_{3}\right\}\right\} \ge 1 - 4Te^{-n},$$
(22)

where $c_1, c_2 > 0$ are universal constants depending on C_1 and C_2 , respectively.

Proof We start with $\mathbb{P}\left\{\mathcal{E}_1 \mid \{X(t)\}_{t=1}^T\right\}$. For any $(i,j) \in \{1,\ldots,d\}^{\otimes 2}$ and any $t \in \{1,\ldots,T\}$, it satisfies that

$$[U_{P_t}^{\top}(A(t)-P_t)U_{P_t}]_{ij} = 2\sum_{k=1}^{n-1}\sum_{l=k+1}^{n}(U_{P_t})_{li}(A(t)-P_t)_{lk}(U_{P_t})_{kj} + \sum_{k=1}^{n}(U_{P_t})_{ki}(A(t)-P_t)_{kk}(U_{P_t})_{kj}.$$
(23)

For any $\varepsilon > 0$, there exists an absolute constant c > 0 such that

$$\mathbb{P}\left\{\left|2\sum_{k=1}^{n-1}\sum_{l=k+1}^{n} (U_{P_t})_{li} (A(t) - P_t)_{lk} (U_{P_t})_{kj}\right| > \varepsilon \left|\{X(t)\}_{t=1}^T\right\}\right\}$$

$$\leq 2 \exp\left\{-\frac{c\varepsilon^{2}}{\sum_{k=1}^{n-1}\sum_{l=k+1}^{n}(U_{P_{l}})_{li}^{2}(U_{P_{l}})_{kj}^{2}}\right\}$$
$$\leq 2 \exp\left\{-\frac{c\varepsilon^{2}}{\sqrt{\sum_{k=1}^{n}\sum_{l=1}^{n}(U_{P_{l}})_{li}^{2}(U_{P_{l}})_{kj}^{2}}}\right\} = 2 \exp\{-c\varepsilon^{2}\},$$
(24)

where the first inequality follows from Theorem 2.6.3 in Vershynin (2018), and the identity follows from the definitions of U_P . Moreover,

$$\left|\sum_{k=1}^{n} (U_{P_t})_{ki} (A(t) - P_t)_{kk} (U_{P_t})_{kj}\right| \le \sum_{k=1}^{n} |(U_P)_{ki} (U_P)_{kj}| \le \sqrt{\sum_{k=1}^{n} (U_P)_{ki}^2 \sum_{k=1}^{n} (U_P)_{kj}^2} = 1.$$
(25)

Combining (23), (24) and (25), and taking ε to be $(C_1/2)\sqrt{\log(n \vee T)}$, we have that

$$\mathbb{P}\{\mathcal{E}_{1}^{c} \mid \{X(t)\}_{t=1}^{T}\} \le 2Td^{2} \exp\left\{-\frac{C_{1}^{2}}{32}\log(n \lor T)\right\} \le (n \lor T)^{-c_{1}},$$

where $c_1 > 0$ depends on C_1 .

In addition, it holds that

$$\mathbb{P}\{\mathcal{E}_1\} = \mathbb{E}\left\{\mathbb{P}\{\mathcal{E}_1 \mid \{X(t)\}_{t=1}^T\}\right\} \ge 1 - (n \lor T)^{-c_1},$$

therefore, (20) follows.

We then show that (21) holds. For $i \in \{1, ..., n\}$ and $j \in \{1, ..., d\}$, we have that

$$[\{A(t) - P_t\}U_{P_t}]_{ij} = \sum_{l \in \{1, \dots, n\} \setminus \{i\}} \{(A(t) - P_t\}_{il}(U_{P_t})_{lj} + \{A(t) - P_t\}_{ii}(U_{P_t})_{ij}.$$

Since

$$|\{A(t) - P_t\}_{ii}(U_{P_t})_{ij}| \le 1$$

and by Hoeffding's inequality that there exists a universal constant c > 0, for any $\varepsilon > 0$,

$$\mathbb{P}\left\{\left|\sum_{l\in\{1,\dots,n\}\setminus\{i\}} \{A(t) - P_t\}_{il}(U_{P_t})_{lj}\right| > \varepsilon \left|\{X(t)\}_{t=1}^T\right\}\right\} \\
\leq 2\exp\left\{-\frac{c\varepsilon^2}{\sum_{l\in\{1,\dots,n\}\setminus\{i\}}(U_{P_t})_{lj}^2}\right\} \leq 2\exp\{-c\varepsilon^2\},.$$
(26)

we have that

$$\mathbb{P}\left\{\max_{t=1,\dots,T} \|\{A(t) - P_t\}U_t\|_{2\to\infty}^2 > (\varepsilon + \sqrt{d})^2\right\}$$
$$= \mathbb{P}\left\{\max_{t=1,\dots,T}\max_{i=1,\dots,n}\sum_{j=1}^d \left[\sum_{l=1}^n \{A(t) - P_t\}_{il}(U_{P_t})_{lj}\right]^2 > (\varepsilon + \sqrt{d})^2\right\}$$

$$\begin{split} &\leq nTd \max_{t=1,\dots,T} \max_{i=1,\dots,n} \max_{j=1,\dots,d} \mathbb{P}\left\{ \left| \sum_{l=1}^{n} \{A(t) - P_t\}_{il} (U_{P_t})_{lj} \right|^2 > \frac{(\varepsilon + \sqrt{d})^2}{d} \right\} \\ &\leq nTd \max_{t=1,\dots,T} \max_{i=1,\dots,n} \max_{j=1,\dots,d} \mathbb{P}\left\{ \left| \sum_{l\in\{1,\dots,n\}\setminus\{i\}} \{A(t) - P_t\}_{il} (U_{P_t})_{lj} \right| > \frac{\varepsilon}{\sqrt{d}} \right\} \\ &\leq 2nTd \exp\left\{ -\frac{c\varepsilon^2}{d} \right\}, \end{split}$$

and (21) follows by taking $\varepsilon = C_2/c\sqrt{d\log(n \vee T)}$.

Lastly, it follows from Eq.(4.18) in Vershynin (2018) that there exists a universal constant $C_3 > 0$, such that

$$\mathbb{P}\{\|A(t) - P_t\|_{\text{op}} > C\sqrt{n} \mid \{X(t)\}_{t=1}^T\} \le 4e^{-n},$$

which leads to (22).

Lemma 12 Under Model 1, it holds that

$$\mathbb{P}\left\{2^{-1}n\min_{k=1,\dots,K}\mu_d^k \le \min_{t=1,\dots,T}\lambda_d(P_t) \le \max_{t=1,\dots,T}\lambda_1(P_t) \le (3/2)n\max_{k=1,\dots,K}\mu_1^k\right\} > 1 - (n \lor T)^{-c_5},$$

Proof We first fix $t \in \{1, ..., T\}$ and for simplicity drop the dependence on t notationally. For $i \in \{1, ..., n\}$, let $Y_i = X_i \Sigma^{-1/2}$ and $Y = (Y_1, ..., Y_n)^{\top} = X \Sigma^{-1/2}$, satisfying $\mathbb{E}\{n^{-1}Y^{\top}Y\} = I_d$.

It follows from Lemma 4.1.5 in Vershynin (2018) that for any $\varepsilon > 0$, if

$$||n^{-1}Y^{\top}Y - I||_{\text{op}} \le \max\{\varepsilon, \varepsilon^2\},$$
(27)

then the eigenvalues of $n^{-1}Y^{\top}Y$ satisfy

$$(1 - \max\{\varepsilon, \varepsilon^2\})^2 \le \lambda_{\min}(n^{-1}Y^{\top}Y) \le \lambda_{\max}(n^{-1}Y^{\top}Y) \le (1 + \max\{\varepsilon, \varepsilon^2\})^2,$$

which implies that

$$n(1 - \max\{\varepsilon, \varepsilon^2\})^2 \le \lambda_{\min}(\Sigma^{-1/2}X^\top X \Sigma^{-1/2})$$
$$\le \lambda_{\max}(\Sigma^{-1/2}X^\top X \Sigma^{-1/2}) \le n(1 + \max\{\varepsilon, \varepsilon^2\})^2$$

Denote $S = \Sigma^{-1/2} X^{\top} X \Sigma^{-1/2}$. We then have

$$\lambda_1(P) = \lambda_{\max}(X^\top X) = \lambda_{\max}(\Sigma^{1/2} S \Sigma^{1/2}) \le n(1 + \max\{\varepsilon, \varepsilon^2\})^2 \max_{k=1,\dots,K} \mu_1^k$$

and

$$\lambda_d(P) = \lambda_{\min}(X^{\top}X) = \lambda_{\min}(\Sigma^{1/2}S\Sigma^{1/2}) = \max_{\dim(E)=d} \min_{v \in \mathcal{S}_E} \langle \Sigma^{1/2}S\Sigma^{1/2}v, v \rangle$$
$$= \max_{\dim(E)=d} \min_{v \in \mathcal{S}_E} \langle S\Sigma^{1/2}v, \Sigma^{1/2}v, \Sigma^{1/2}v \rangle = \max_{\dim(E)=d} \min_{v \in \mathcal{S}_E} \|\Sigma^{1/2}v\|^2 \left\langle S\frac{\Sigma^{1/2}v}{\|\Sigma^{1/2}v\|}, \frac{\Sigma^{1/2}v}{\|\Sigma^{1/2}v\|} \right\rangle$$

$$\geq \max_{\dim(E)=d} \min_{v \in \mathcal{S}_E} \left\langle S \frac{\Sigma^{1/2} v}{\|\Sigma^{1/2} v\|}, \frac{\Sigma^{1/2} v}{\|\Sigma^{1/2} v\|} \right\rangle \min_{k=1,\dots,K} \mu_d^k \geq \max_{\dim(E)=d} \min_{v \in \mathcal{S}_E} \langle Sv, v \rangle \min_{k=1,\dots,K} \mu_d^k \\ \geq n(1 - \max\{\varepsilon, \varepsilon^2\})^2 \min_{k=1,\dots,K} \mu_d^k.$$

Now it suffices to investigate (27). Since

$$\|n^{-1}Y^{\top}Y - I\|_{\text{op}} = \sup_{v \in \mathcal{S}^{d-1}} \left| \frac{1}{n} \sum_{i=1}^{n} \left\{ (Y_i^{\top}v)^2 - 1 \right\} \right|,$$

taking \mathcal{N} to be a 1/4-net on \mathcal{S}^{d-1} , it holds that

$$\begin{split} \mathbb{P}\left\{ \|n^{-1}Y^{\top}Y - I\|_{\mathrm{op}} > C\sqrt{\frac{\log(n \vee T)}{n}} \right\} \\ & \leq 9^{d} \max_{v \in \mathcal{N}} \mathbb{P}\left\{ \left| \frac{1}{n} \sum_{i=1}^{n} \left\{ (Y_{i}^{\top}v)^{2} - 1 \right\} \right| > C\sqrt{\frac{\log(n \vee T)}{n}} \right\} \\ & \leq 2 \times 9^{d} \exp\{-c \log(n \vee T)\}, \end{split}$$

where C, c > 0 are universal constants.

Thus we have that

$$\mathbb{P}\left\{2^{-1}n\min_{k=1,\dots,K}\mu_d^k \le \min_{t=1,\dots,T}\lambda_d(P_t) \le \max_{t=1,\dots,T}\lambda_1(P_t) \le (3/2)n\max_{k=1,\dots,K}\mu_1^k\right\} > 1 - (n \lor T)^{-c_4},$$

where $c_4 > 0$ is a universal constant.

Lemma 13 is adapted from Theorem 8 in Athreya et al. (2018).

Lemma 13 It holds that

$$\mathbb{P}\left\{\max_{t=1,\dots,T}\min_{W\in\mathbb{O}_{d}}\|\widehat{X}(t)-X(t)W\|_{2\to\infty} > C_{W}\frac{\sqrt{d\log(n\vee T)}\vee d^{3/2}}{n^{1/2}}\right\}$$

$$\leq 1-(n\vee T)^{-c_{1}}-(n\vee T)^{-c_{2}}-4Te^{-n}-(n\vee T)^{-c_{4}}.$$

Proof [Proof of Lemma 13] We first work on a fixed $t \in \{1, ..., T\}$, and then use union bounds arguments to reach the final conclusion. For simplicity, we drop the dependence on t for now. Recall that

$$\widehat{X} = U_A S_A^{1/2}$$
 and $X = U_P S_P^{1/2}$.

Define $W^* = W_1 W_2^{\top}$, where W_1 and W_2 are the left and right singular vectors of $U_P^{\top} U_A$, that $U_P^{\top} U_A = W_1 \Lambda_1 W_2^{\top}$. Since $W^* \in \mathbb{O}_d$, we have that

$$\min_{W \in \mathbb{O}_d} \|\widehat{X} - XW\|_{2 \to \infty} \le \|\widehat{X} - XW^*\|_{2 \to \infty}.$$

In the rest of this proof, denote by $\lambda_1, \ldots, \lambda_n$ as the eigenvalues of P, with $|\lambda_1| \ge \cdots |\lambda_n|$; denote by $\hat{\lambda}_1, \ldots, \hat{\lambda}_n$ the eigenvalues of A, with $|\hat{\lambda}_1| \ge \cdots \ge |\hat{\lambda}_n|$.

Step 1. We first provide a deterministic upper bound for $||W^*S_A^{1/2} - S_P^{1/2}W^*||_F$. We have,

$$\begin{split} W^*S_A &= (W^* - U_P^{\top}U_A)S_A + U_P^{\top}U_AS_A = (W^* - U_P^{\top}U_A)S_A + U_P^{\top}AU_A \\ &= (W^* - U_P^{\top}U_A)S_A + U_P^{\top}(A - P)U_A + U_P^{\top}PU_A \\ &= (W^* - U_P^{\top}U_A)S_A + U_P^{\top}(A - P)(U_A - U_PU_P^{\top}U_A) + U_P^{\top}(A - P)U_P + S_PU_P^{\top}U_A \\ &= (W^* - U_P^{\top}U_A)S_A + U_P^{\top}(A - P)(U_A - U_PU_P^{\top}U_A) + U_P^{\top}(A - P)U_P \\ &+ S_P(U_P^{\top}U_A - W^*) + S_PW^*, \end{split}$$

where the second and the fourth inequalities are due to

$$AU_A = U_A S_A U_A^{\top} U_A = U_A S_A$$
 and $U_P^{\top} P = U_P^{\top} U_P S_P U_P^{\top} = S_P U_P^{\top}$,

respectively. Therefore,

$$\|W^*S_A - S_PW^*\|_{\mathbf{F}} \leq \|W^* - U_P^{\top}U_A\|_{\mathbf{F}}(\|S_A\|_{\mathrm{op}} + \|S_P\|_{\mathrm{op}}) + \|U_P^{\top}(A - P)(U_A - U_PU_P^{\top}U_A)\|_{\mathbf{F}} + \|U_P^{\top}(A - P)U_P\|_{\mathbf{F}} \leq \|I_n - \Lambda_1\|_{\mathbf{F}}\|W_1\|_{\mathrm{op}}\|W_2\|_{\mathrm{op}}(\|S_A\|_{\mathrm{op}} + \|S_P\|_{\mathrm{op}}) + \|A - P\|_{\mathrm{op}}\|U_A - U_PU_P^{\top}U_A\|_{\mathbf{F}} + \|U_P^{\top}(A - P)U_P\|_{\mathbf{F}} \leq \|I_n - \Lambda_1\|_{\mathbf{F}}(2\lambda_1 + \|A - P\|_{\mathrm{op}}) + \|A - P\|_{\mathrm{op}}\|U_A - U_PU_P^{\top}U_A\|_{\mathbf{F}} + \|U_P^{\top}(A - P)U_P\|_{\mathbf{F}} = (I) + (II) + (III),$$
(28)

where λ_1 is the largest singular value of P and the last inequality is due to Weyl's inequality.

In addition, let $\{\theta_1, \ldots, \theta_d\}$ be the principal angles between the column spaces spanned by U_A and U_P . We thus have

$$\|I_n - \Lambda_1\|_{\mathrm{F}} = \sqrt{\sum_{i=1}^d (1 - \cos\theta_i)^2} \le \sqrt{d}(1 - \cos^2\theta_1) = \sqrt{d}\sin^2\theta_1 = \sqrt{d}\min_{W \in \mathbb{O}_d} \|U_A - U_P W\|_{\mathrm{op}}^2$$
$$\le \sqrt{d}\min_{W \in \mathbb{O}_d} \|U_A - U_P W\|_{\mathrm{F}}^2 \le \frac{4d^{3/2} \|A - P\|_{\mathrm{op}}^2}{\lambda_d^2}, \tag{29}$$

where the first and second inequalities are due to $\cos \theta_i$, $\sin \theta_i \in [0, 1]$, and the last inequality is due to Theorem 2 in Yu et al. (2014) and the fact that $\lambda_{d+1} = 0$.

As for term (II), there exists $W \in \mathbb{O}_d$ such that

$$\|U_A - U_P U_P^\top U_A\|_{\mathrm{F}} = \sqrt{\operatorname{tr}(U_A U_A^\top - U_A U_A^\top U_P U_P^\top)} = \sqrt{d - \operatorname{tr}(U_A^\top U_P W W^\top U_P^\top U_A)}$$
$$= \sqrt{\sum_{i=1}^d (1 - \cos^2 \theta_i)} = \sqrt{\sum_{i=1}^d \sin^2 \theta_i} \le \frac{2\sqrt{d} \|A - P\|_{\mathrm{op}}}{\lambda_d}.$$
(30)

Term (III) is dealt in Lemma 11. As for $||W^*S_A^{1/2} - S_P^{1/2}W^*||_F$, we note that the *ij*-th entry of $W^*S_A^{1/2} - S_P^{1/2}W^*$ satisfies that

$$|W_{ij}^*(\hat{\lambda}_j^{1/2} - \lambda_i^{1/2})| = \left|\frac{W_{ij}^*(\hat{\lambda}_j - \lambda_i)}{\hat{\lambda}_j^{1/2} + \lambda_i^{1/2}}\right| = \left|\frac{(W^*S_A - S_PW^*)_{ij}}{\hat{\lambda}_j^{1/2} + \lambda_i^{1/2}}\right| \le \frac{|(W^*S_A - S_PW^*)_{ij}|}{\lambda_d^{1/2}},$$

which means

$$\|W^* S_A^{1/2} - S_P^{1/2} W^*\|_{\mathcal{F}} \le \frac{\|W^* S_A - S_P W^*\|_{\mathcal{F}}}{\lambda_d^{1/2}} \le \frac{8d^{3/2} \|A - P\|_{\mathrm{op}}^2 \lambda_1}{\lambda_d^{5/2}} + \frac{4d^{3/2} \|A - P\|_{\mathrm{op}}^3}{\lambda_d^{5/2}} + \frac{2d^{1/2} \|A - P\|_{\mathrm{op}}^2}{\lambda_d^{3/2}} + \frac{\|U_P^{\top} (A - P)U_P\|_{\mathcal{F}}}{\lambda_d^{1/2}}.$$
 (31)

Step 2. We then provide an upper bound for $\min_{W \in \mathbb{O}_d} \|\widehat{X} - XW\|_{2 \to \infty}$. Since

$$\min_{W \in \mathbb{O}_d} \|\widehat{X} - XW\|_{2 \to \infty} \le \|\widehat{X} - XW^*\|_{2 \to \infty}$$

in the rest of this step, we work on $\|\widehat{X} - XW^*\|_{2\to\infty}$. We have that

$$\begin{split} \|\widehat{X} - XW^*\|_{2\to\infty} &= \|U_A S_A^{1/2} - U_P S_P^{1/2} W^*\|_{2\to\infty} \\ &= \|U_A S_A^{1/2} - U_P W^* S_A^{1/2} + U_P (W^* S_A^{1/2} - S_P^{1/2} W^*)\|_{2\to\infty} \\ &\leq \|(U_A - U_P U_P^\top U_A) S_A^{1/2}\|_{2\to\infty} + \|U_P (U_P^\top U_A - W^*) S_A^{1/2}\|_{2\to\infty} \\ &\quad + \|U_P (W^* S_A^{1/2} - S_P^{1/2} W^*)\|_{2\to\infty} \\ &= (I) + (II) + (III). \end{split}$$
(32)

As for term (I), it holds that

$$(U_A - U_P U_P^{\top} U_A) S_A^{1/2} = (A - P) U_A S_A^{-1/2} - U_P U_P^{\top} (A - P) U_A S_A^{-1/2}$$

= $(A - P) U_P W^* S_A^{-1/2} - U_P U_P^{\top} (A - P) U_P W^* S_A^{-1/2}$
+ $(I - U_P U_P^{\top}) (A - P) (U_A - U_P W^*) S_A^{-1/2}$,

which satisfies

$$\|(A-P)U_PW^*S_A^{-1/2}\|_{2\to\infty} \le \|(A-P)U_P\|_{2\to\infty}(\widehat{\lambda}_d)^{-1/2},\\ \|U_PU_P^\top(A-P)U_pW^*S_A^{-1/2}\|_{2\to\infty} \le \|U_P^\top(A-P)U_P\|_{\mathbf{F}}(\widehat{\lambda}_d)^{-1/2}$$

and

$$\|(I - U_P U_P^{\top})(A - P)(U_A - U_P W^*) S_A^{-1/2} \|_{2 \to \infty}$$

$$\leq \|A - P\|_{\rm op} \|U_A - U_P W^*\|_{\rm F} (\widehat{\lambda}_d)^{-1/2} \leq \frac{4d^{3/2} \|A - P\|_{\rm op}^3 (\widehat{\lambda}_d)^{-1/2}}{\lambda_d^2},$$

which is due to (30). Therefore we have that

$$\|(I)\|_{2\to\infty} \le \|(A-P)U_P\|_{2\to\infty}(\widehat{\lambda}_d)^{-1/2} + \|U_P^{\top}(A-P)U_P\|_{\mathbf{F}}(\widehat{\lambda}_d)^{-1/2} + \frac{4d^{3/2}\|A-P\|_{\mathrm{op}}^3(\widehat{\lambda}_d)^{-1/2}}{\lambda_d^2}.$$
(33)

As for term (II), it holds that

$$\|U_P(U_P^{\top}U_A - W^*)S_A^{1/2}\|_{2 \to \infty} \le \|I - \Lambda_1\|_{\mathrm{F}}(\lambda_1 + \|A - P\|_{\mathrm{op}})^{1/2} \le \frac{4d^{3/2}\|A - P\|_{\mathrm{op}}^2}{\lambda_d^2}\sqrt{\frac{3\lambda_1}{2}}.$$
 (34)

As for term (III), it holds that

$$\|U_P(W^*S_A^{1/2} - S_P^{1/2}W^*)\|_{2\to\infty} \le \|(W^*S_A^{1/2} - S_P^{1/2}W^*)\|_{\mathrm{F}}.$$
(35)

Combining (31), (32), (33), (34) and (35), we have that

$$\begin{split} \min_{W \in \mathbb{O}_d} \| \hat{X} - XW \|_{2 \to \infty} &\leq \| (A - P)U_P \|_{2 \to \infty} (\hat{\lambda}_d)^{-1/2} + \| U_P^\top (A - P)U_P \|_F (\hat{\lambda}_d)^{-1/2} \\ &+ \frac{4d^{3/2} \|A - P\|_{\mathrm{op}}^3 (\hat{\lambda}_d)^{-1/2}}{\lambda_d^2} + \frac{4d^{3/2} \|A - P\|_{\mathrm{op}}^3}{\lambda_d^2} \sqrt{\frac{3\lambda_1}{2}} \\ &+ \frac{8d^{3/2} \|A - P\|_{\mathrm{op}}^2 \lambda_1}{\lambda_d^{5/2}} + \frac{4d^{3/2} \|A - P\|_{\mathrm{op}}^3}{\lambda_d^{5/2}} + \frac{2d^{1/2} \|A - P\|_{\mathrm{op}}^2}{\lambda_d^{3/2}} + \frac{\| U_P^\top (A - P)U_P \|_F}{\lambda_d^{1/2}} \\ &\leq \frac{\| (A - P)U_P \|_{2 \to \infty}}{\sqrt{\lambda_d} - \| A - P \|_{\mathrm{op}}} + \frac{\| U_P^\top (A - P)U_P \|_F}{\sqrt{\lambda_d} - \| A - P \|_{\mathrm{op}}} \\ &+ \frac{4d^{3/2} \|A - P\|_{\mathrm{op}}^3}{\lambda_d^2 \sqrt{\lambda_d} - \| A - P \|_{\mathrm{op}}^2} + \frac{4d^{3/2} \|A - P \|_{\mathrm{op}}^2}{\lambda_d^2} \sqrt{\frac{3\lambda_1}{2}} \\ &+ \frac{8d^{3/2} \|A - P \|_{\mathrm{op}}^2 \lambda_1}{\lambda_d^{5/2}} + \frac{4d^{3/2} \|A - P \|_{\mathrm{op}}^3}{\lambda_d^{5/2}} + \frac{2d^{1/2} \|A - P \|_{\mathrm{op}}^2}{\lambda_d^{3/2}} + \frac{\| U_P^\top (A - P)U_P \|_F}{\lambda_d^{1/2}} \end{split}$$

where the second inequality follows from that $\widehat{\lambda}_d \geq \lambda_d - \|A - P\|_{\text{op}}$. It holds on the event $\mathcal{E}_1 \cap \mathcal{E}_2 \cap \mathcal{E}_3 \cap \mathcal{E}_4$, that

$$\mathbb{P}\left\{\max_{t=1,\dots,T}\min_{W\in\mathbb{O}_d}\|\widehat{X}(t)-X(t)W\|_{2\to\infty} > C_W \frac{\sqrt{d\log(n\vee T)}\vee d^{3/2}}{n^{1/2}}\right\}$$

$$\leq 1-(n\vee T)^{-c_1}-(n\vee T)^{-c_2}-(n\vee T)^{-c_4}-4Te^{-n},$$

where $C_W > 0$ is a universal constant depending only on $C_1, C_2, C_3, \max_{k=1,\dots,K} \mu_1^k$ and $\min_{k=1,\dots,K} \mu_d^k$.

We first state a weakly dependent version of Bernstein inequality. This is in fact Theorem 4 in Delyon (2009). The notation in Lemma 14 only applies within Lemma 14. **Lemma 14** Let $\{X_1, \ldots, X_T\}$ be centred random variables. Define

$$g = \sum_{t=2}^{T} \sum_{s=1}^{t-1} \|X_s\|_{\infty} \|\mathbb{E}(X_t \mid \mathcal{F}_s)\|_{\infty}, \quad v = \sum_{t=1}^{T} \|\mathbb{E}(X_t^2 \mid \mathcal{F}_{t-1})\|_{\infty}$$

and

$$m = \max_{t=1,\dots,T} \|X_t\|_{\infty},$$

where $F_s = \sigma\{X_1, \ldots, X_s\}$, $s \ge 1$, is the natural σ -field generated by $\{X_i\}_{i=1}^s$. For any $\varepsilon > 0$, it holds that

$$\mathbb{P}\left\{\left|\sum_{t=1}^{T} X_t\right| > \epsilon\right\} \le 2\exp\left(-\frac{\varepsilon^2}{2(v+2g) + 2\varepsilon m/3}\right).$$

Lemma 15 Under Model 1, it holds that for any $z \in \mathbb{R}$,

$$\mathbb{P}\left\{\max_{0 \le s < t < e \le T} \left|\Delta_{s,e}^{t}(z)\right| \ge C_8 \sqrt{T} \max\{\sqrt{d \log(n \lor T)}, \ d^{3/2}\}\right\} \le 4(n \lor T)^{-c} + 4Te^{-n},$$

where $c = \min\{c_1, c_2, c_4, c_5\} - 1 > 0$ is a universal constant.

In addition,

$$\mathbb{P}\left\{\max_{0\leq s$$

where $c = \min\{c_1, c_2, c_4, c_5\} - 1 > 0$ is a universal constant.

Proof For any $(i, j) \in \mathcal{O}$ and $t \in \{1, \ldots, T\}$, it holds that

$$\begin{split} \left| \widehat{Y}_{ij}^{t} - Y_{ij}^{t} \right| &= \left| (\widehat{X}_{i}(t))^{\top} \widehat{X}_{j}(t) - (X_{i}(t))^{\top} X_{j}(t) \right| = \left| (\widehat{X}_{i}(t))^{\top} \widehat{X}_{j}(t) - (W_{t} X_{i}(t))^{\top} W_{t} X_{j}(t) \right| \\ &\leq \left| (\widehat{X}_{i}(t) - W_{t} X_{i}(t))^{\top} W_{t} X_{j}(t) \right| + \left| (\widehat{X}_{i}(t) - W_{t} X_{i}(t))^{\top} (W_{t} X_{j}(t) - \widehat{X}_{j}(t)) \right| \\ &+ \left| (\widehat{X}_{j}(t) - W_{t} \widehat{X}_{j}(t))^{\top} W_{t} X_{i}(t) \right| \\ &\leq 2 \max_{t=1,\dots,T} \min_{W \in \mathbb{O}_{d}} \left\| \widehat{X}(t) - X(t) W^{\top} \right\|_{2 \to \infty} \max_{\substack{t=1,\dots,T \\ i=1,\dots,n}} \left\| X_{i}(t) \right\| \\ &+ \left(\max_{t=1,\dots,T} \min_{W \in \mathbb{O}_{d}} \left\| \widehat{X}(t) - X(t) W^{\top} \right\|_{2 \to \infty} \right)^{2} \\ &\leq 2 \max_{t=1,\dots,T} \min_{W \in \mathbb{O}_{d}} \left\| \widehat{X}(t) - X(t) W^{\top} \right\|_{2 \to \infty} + \left(\max_{t=1,\dots,T} \min_{W \in \mathbb{O}_{d}} \left\| \widehat{X}(t) - X(t) W^{\top} \right\|_{2 \to \infty} \right)^{2}, \end{split}$$

where $W_t \in \mathbb{O}_d$ satisfies

$$\|\widehat{X}(t) - X(t)W_t^{\top}\|_{2 \to \infty} = \min_{W \in \mathbb{O}_d} \|\widehat{X}(t) - X(t)W^{\top}\|_{2 \to \infty}.$$

We fix the chosen pairs $\mathcal{O} \subset \{1, \ldots, n\}^{\otimes 2}$ with $|\mathcal{O}| = n/2$, which is assumed to be an integer. As for the sequence $\{w_k\}$, it holds that

$$\sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} w_k^2 = 1.$$
(37)

We have for any $z \in \mathbb{R}$, it holds that

$$\begin{split} \left| \Delta_{s,e}^{t}(z) \right| &\leq \left| \sum_{k=s+1}^{e} w_{k} \sum_{(i,j) \in \mathcal{O}} \left(\mathbb{1}\{\widehat{Y}_{ij}^{k} \leq z\} - \mathbb{1}\{Y_{ij}^{k} \leq z\} \right) \right| \\ &+ \left| \sum_{k=s+1}^{e} w_{k} \sum_{(i,j) \in \mathcal{O}} \left(\mathbb{1}\{Y_{ij}^{k} \leq z\} - \mathbb{E}\left\{ \mathbb{1}\{Y_{ij}^{k} \leq z\} \right\} \right) \right| = (I) + (II). \end{split}$$

Term (II). As for (II), notice that

$$\mathbb{E}\left(\mathbb{1}\{Y_{ij}^k \le z\} - \mathbb{E}\left\{\mathbb{1}\{Y_{ij}^k \le z\}\right\}\right) = 0.$$

In order to apply Lemma 14, we let

$$V_i(k) = w_k \mathbb{1}\{Y_{ij}^k \le z\} - w_k \mathbb{E}\left\{\mathbb{1}\{Y_{ij}^k \le z\}\right\},\$$

with i = 1, ..., n/2, k = 1, ..., T. We order $\{V_i(k)\}$ as

$$V_1(1), \dots, V_1(T), V_2(1), \dots, V_2(T), \dots, V_{n/2}(1), \dots, V_{n/2}(T).$$
 (38)

Denote $\mathcal{F}_{i,t}$ as the natural σ -field generated by $V_i(t)$ and all the random variables before it in the order of (38), and denote $\mathcal{F}_{i,t,-}$ as the natural σ -field generated by all the random variables before $Y_i(t)$ in the order of (38) excluding $Y_i(t)$. If (i,t) = (1,1), then $\mathcal{F}_{i,t,-}$ is the σ -field generated by constants.

In addition, for the notation in Lemma 14, we have that

$$\begin{aligned} v &= \sum_{i=1}^{n/2} \sum_{t=s+1}^{e} \left\| \mathbb{E}(V_{i}(t)^{2} \mid \mathcal{F}_{i,t,-}) \right\|_{\infty} \\ &= \sum_{i=1}^{n/2} \sum_{k: \ \eta_{k} \in (s,e)} \left[(w_{\eta_{k}+1})^{2} \mathbb{E} \left\{ \mathbb{1}\{Y_{ij}^{\eta_{k}+1} \leq z\} \right\} (1 - \mathbb{E} \left\{ \mathbb{1}\{Y_{ij}^{\eta_{k}+1} \leq z\} \right\}) \right] \\ &+ \sum_{i=1}^{n/2} \sum_{\substack{t \in (s,e] \\ t \notin \{\eta_{k}+1\}}} (1 - \rho)(w_{t})^{2} \mathbb{E} \left\{ \mathbb{1}\{Y_{ij}^{t} \leq z\} \right\} (1 - \mathbb{E} \left\{ \mathbb{1}\{Y_{ij}^{t} \leq z\} \right\}) \\ &+ \sum_{i=1}^{n/2} \sum_{\substack{t \in (s,e] \\ t \notin \{\eta_{k}+1\}}} \rho(w_{t})^{2} \| (\mathbb{1}\{Y_{ij}^{t-1} \leq z\} - \mathbb{E} \left\{ \mathbb{1}\{Y_{ij}^{t-1} \leq z\} \right\})^{2} \|_{\infty} \end{aligned}$$

$$\leq 1 + \rho, \tag{39}$$

where the last inequality is due to (37),

$$m \le \max_{t=1,\dots,T} |w_t|,\tag{40}$$

and

$$g = (n/2) \sum_{k:\eta_k \in (s,e)} \left(\sum_{t=\eta_k+2}^{\min\{\eta_{k+1},e\}} \sum_{u=\eta_k+1}^t + \sum_{t=s+1}^{\eta_{k_0+1}} \sum_{u=s+2}^{t-1} \right) |w_t w_u| \rho^{t-u}.$$
(41)

Combining (39), (40), (41) and Lemma 14, we have for any $\varepsilon > 0$, it holds that

$$\mathbb{P}\left((II) \ge \varepsilon\right) \le 2\exp\left\{-C\varepsilon^2/((1-\rho)^{-1}+\varepsilon)\right\}.$$

We thus denote

$$\mathcal{E}_5 = \left\{ \max_{1 < s < t < e \le T} \left| \sum_{k=s+1}^e w_k \sum_{(i,j) \in \mathcal{O}} \left(\mathbbm{1}\{Y_{ij}^k \le z\} - \mathbbm{1}\{\mathbbm{1}\{Y_{ij}^k \le z\}\} \right) \right| \ge C_5 \sqrt{\frac{\log(n \lor T)}{1-\rho}} \right\},$$

where $C_5 > 0$ is a universal constant, and therefore it holds that

$$\mathbb{P}\{\mathcal{E}_5\} \le (n \lor T)^{-c_5},$$

where $c_5 > 0$ is a universal constant.

Term (I). As for (I), we have that

$$\mathbb{E}\left\{ \left| \mathbb{1}\{\widehat{Y}_{ij}^k \leq z\} - \mathbb{1}\{Y_{ij}^k \leq z\} \right| \right\}$$

$$\leq \max\left\{ \mathbb{P}\left\{ \left(\widehat{Y}_{ij}^k \leq z\right) \cap \left(Y_{ij}^k > z\right) \right\}, \mathbb{P}\left\{ \left(\widehat{Y}_{ij}^k > z\right) \cap \left(Y_{ij}^k \leq z\right) \right\} \right\} = \max\{(I.1), (I.2)\}.$$

Let

Let

$$\mathcal{E}_6 = \left\{ \max_{t=1,\dots,T} \min_{W \in \mathbb{O}_d} \|\widehat{X}_t - X_t W\| \le C_W \frac{\sqrt{d \log(n \vee T)} \vee d^{3/2}}{n^{1/2}} \right\}.$$

On the event \mathcal{E}_6 , it holds that

$$\max_{\substack{t=1,\dots,T\\(i,j)\in\mathcal{O}}} \left| \widehat{Y}_{ij}^t - Y_{ij}^t \right| \le 3C_W \frac{\sqrt{d\log(n \vee T)} \vee d^{3/2}}{n^{1/2}} = \delta$$

and

$$\mathbb{P}\left\{\max_{\substack{t=1,...,T\\(i,j)\in\mathcal{O}}} \left|\widehat{Y}_{ij}^{t} - Y_{ij}^{t}\right| \le \delta\right\}$$

$$\ge 1 - (n \lor T)^{-c_{1}} - (n \lor T)^{-c_{2}} - (n \lor T)^{-c_{4}} - 4Te^{-n} = 1 - p_{\delta}.$$

Therefore,

$$(I.1) = \mathbb{P}\left\{\left(\widehat{Y}_{ij}^k \le z\right) \cap \left(Y_{ij}^k > z\right) | Y_{ij}^k > z + \delta\right\} \mathbb{P}\left\{Y_{ij}^k > z + \delta\right\} \\ + \mathbb{P}\left\{\left(\widehat{Y}_{ij}^k \le z\right) \cap \left(Y_{ij}^k > z\right) | Y_{ij}^k < z + \delta\right\} \mathbb{P}\left\{Y_{ij}^k < z + \delta\right\} \\ \le p_{\delta}(1 - F_k(z + \delta)) + F_k(z + \delta) - F_k(z) \le p_{\delta} + \delta C_F$$

and

$$(I.2) = \mathbb{P}\left\{\left(\widehat{Y}_{ij}^k > z\right) \cap \left(Y_{ij}^k \le z\right) \middle| Y_{ij}^k \le z - \delta\right\} \mathbb{P}\left\{Y_{ij}^k \le z - \delta\right\} \\ + \mathbb{P}\left\{\left(\widehat{Y}_{ij}^k > z\right) \cap \left(Y_{ij}^k \le z\right) \middle| Y_{ij}^k > z - \delta\right\} \mathbb{P}\left\{Y_{ij}^k > z - \delta\right\} \\ \le p_{\delta}F_k(z - \delta) + F_k(z) - F_k(z - \delta) \le p_{\delta} + \delta C_F.$$

Then we have,

$$\mathbb{E}\left|\sum_{k=s+1}^{e} w_k \sum_{(i,j)\in\mathcal{O}} \left(\mathbbm{1}\{\widehat{Y}_{ij}^k \le z\} - \mathbbm{1}\{Y_{ij}^k \le z\}\right)\right| \le 2\sqrt{\frac{n}{2}}\sqrt{\frac{(e-t)(t-s)}{e-s}}(p_\delta + \delta C_F) \le 2\sqrt{\frac{n}{2}}\min\{\sqrt{e-t}, \sqrt{t-s}\}(p_\delta + \delta C_F).$$

Therefore, following from similar arguments as those used in bounding (II), we have that for any $\varepsilon > 0$, it holds that

$$\mathbb{P}\left\{\left|\sum_{k=s+1}^{e} w_k \sum_{(i,j)\in\mathcal{O}} \left(\mathbbm{1}\{\widehat{Y}_{ij}^k \le z\} - \mathbbm{1}\{Y_{ij}^k \le z\}\right) - \mathbbm{E}\left\{\sum_{k=s+1}^{e} w_k \sum_{(i,j)\in\mathcal{O}} \left(\mathbbm{1}\{\widehat{Y}_{ij}^k \le z\} - \mathbbm{1}\{Y_{ij}^k \le z\}\right)\right\}\right| > \varepsilon\right\}$$
$$\leq 2\exp\left\{-C\varepsilon^2/((1-\rho)^{-1}+\varepsilon)\right\},$$

which implies that

$$\mathbb{P}\left\{\left|\sum_{k=s+1}^{e} w_k \sum_{(i,j)\in\mathcal{O}} \left(\mathbbm{1}\{\widehat{Y}_{ij}^k \le z\} - \mathbbm{1}\{Y_{ij}^k \le z\}\right)\right| \\> \mathbb{E}\left|\sum_{k=s+1}^{e} w_k \sum_{(i,j)\in\mathcal{O}} \left(\mathbbm{1}\{\widehat{Y}_{ij}^k \le z\} - \mathbbm{1}\{Y_{ij}^k \le z\}\right)\right| + \varepsilon/2\right\}$$
$$\leq \mathbb{P}\left\{\left|\sum_{k=s+1}^{e} w_k \sum_{(i,j)\in\mathcal{O}} \left(\mathbbm{1}\{\widehat{Y}_{ij}^k \le z\} - \mathbbm{1}\{Y_{ij}^k \le z\}\right)\right| \\> 2\sqrt{\frac{n}{2}}\min\{\sqrt{e-t}, \sqrt{t-s}\}(p_{\delta} + \delta C_F) + \varepsilon/2\right\}$$

$$\leq 2\exp\left\{-C\varepsilon^2/((1-\rho)^{-1}+\varepsilon)\right\}+p_{\delta}.$$

Lastly, we have that

$$\mathbb{P}\left\{\max_{0\leq s< t< e\leq T} \left|\Delta_{s,e}^{t}(z)\right| \geq C_{8}\sqrt{\frac{T}{1-\rho}}\max\{\sqrt{d\log(n\vee T)}, d^{3/2}\}\right\}$$
$$\leq \mathbb{P}\left\{\left|\Delta_{s,e}^{t}(z)\right| > C_{5}\sqrt{\frac{\log(n\vee T)}{1-\rho}} + \sqrt{2n}\min\{\sqrt{e-t}, \sqrt{t-s}\}(p_{\delta}+\delta C_{F})\right\}$$
$$\leq 4(n\vee T)^{-c} + 4Te^{-n},$$

where $c = \min\{c_1, c_2, c_4, c_5\} - 1 > 0$ is a universal constant.

The result (36) follows from the identical arguments.

Lemma 16 Let

$$\Delta_{s,e}^t = \sup_{z \in \mathbb{R}} |\Delta_{s,e}^t(z)|$$

It holds that

$$\mathbb{P}\left\{\max_{0 \le s < t < e \le T} \Delta_{s,e}^t > C_9 T^{1/2} (1-\rho)^{-1/2} \max\{\sqrt{d\log(n \lor T)}, d^{3/2}\}\right\} \le 11 (n \lor T)^{-c} + 8Te^{-n}.$$

In addition,

$$\mathbb{P}\left\{ \max_{0 \le s < t < e \le T} \sup_{z \in \mathbb{R}} \left| \sqrt{\frac{2}{n(e-s)}} \sum_{k=s+1}^{e} \sum_{(i,j) \in \mathcal{O}} \left(\mathbb{1}\{\widehat{Y}_{ij}^k \le z\} - \mathbb{E}\left\{ \mathbb{1}\{Y_{ij}^k \le z\} \right\} \right) \right| \\
\le C_9 T^{1/2} (1-\rho)^{-1/2} \max\{\sqrt{d \log(n \lor T)}, d^{3/2}\} \right\} \le 11 (n \lor T)^{-c} + 8Te^{-n}.$$
(42)

 $\mathbf{Proof} \ \ \mathrm{Let}$

$$\delta = 3C_W \frac{\sqrt{d \log(n \vee T)} \vee d^{3/2}}{n^{1/2}}.$$
(43)

Let $z_m = m\delta$, $m = 1, \ldots, \lfloor 1/\delta \rfloor$. Let $I_m = [z_m - \delta, z_m + \delta]$, for $m = 1, \ldots, \lfloor 1/\delta \rfloor - 1$, and $I_{\lfloor 1/\delta \rfloor} = [z_{\lfloor 1/\delta \rfloor - 1}, 1]$. Let $M = \lfloor 1/\delta \rfloor$. Then

$$\sup_{z\in\mathbb{R}} |\Delta_{s,e}^t(z)| \le \max_{j=1,\dots,M} \left\{ |\Delta_{s,e}^t(z_j)| + \sup_{z\in I_j} |\Delta_{s,e}^t(z_j) - \Delta_{s,e}^t(z)| \right\}.$$
(44)

It follows from Lemma 15 that

$$\mathbb{P}\left\{\max_{j=1,\dots,M} |\Delta_{s,e}^t(z_j)| \ge C_8 \sqrt{T} (1-\rho)^{-1/2} \max\{\sqrt{d\log(n \vee T)}, \, d^{3/2}\}\right\} \le 4(n \vee T)^{-c} + 4Te^{-n}.$$
(45)

For every $z \in \mathbb{R}$, on the event

$$\left\{ \max_{\substack{k=1,\dots,T\\(i,j)\in\mathcal{O}}} \left| \widehat{Y}_{ij}^k - Y_{ij}^k \right| \le \delta \right\},\,$$

it holds that

$$\left| \mathbb{1}\{\widehat{Y}_{ij}^k \le z\} - \mathbb{1}\{Y_{ij}^k \le z\} \right| \le \mathbb{1}\{Y_{ij}^k \in [z - \delta, z + \delta]\}.$$

For any $z \in \mathbb{R}$, there exist z_m and z_{m+1} , $m \in \{1, \ldots, M = 1\}$, such that

$$[z-\delta, z+\delta] \subset [z_m-\delta, z_m+\delta] \cup [z_{m+1}-\delta, z_{m+1}+\delta].$$

Let

$$B_m = \sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} \mathbb{1}\{Y_{i,j}^k \in I_m\}, \ m = 1, \dots, M.$$

Therefore

$$\begin{split} |\Delta_{s,e}^{t}(z_{m}) - \Delta_{s,e}^{t}(z)| \\ &\leq \left|\sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} w_{k} \{\mathbbm{1}\{\widehat{Y}_{i,j}^{k} \leq z_{m}\} - \mathbbm{1}\{Y_{i,j}^{k} \leq z_{m}\}\}\right| \\ &+ \left|\sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} w_{k} \{\mathbbm{1}\{\widehat{Y}_{i,j}^{k} \leq z\} - \mathbbm{1}\{Y_{i,j}^{k} \leq z\}\}\right| \\ &+ \left|\sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} w_{k} \{\mathbbm{1}\{Y_{i,j}^{k} \leq z_{m}\} - \mathbbm{1}\{Y_{i,j}^{k} \leq z\}\}\right| \\ &+ \left|\sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} w_{k} \{\mathbbm{1}\{Y_{i,j}^{k} \leq z_{m}\} - \mathbbm{1}\{Y_{i,j}^{k} \leq z\}\}\right| \\ &+ \left|\sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} w_{k} \{\mathbbm{1}\{Y_{i,j}^{k} \leq z_{m}\} - \mathbbm{1}\{Y_{i,j}^{k} \leq z\}\}\right| \\ &+ \left|\sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} \|\mathbbm{1}\{\widehat{Y}_{i,j}^{k} \leq z\} - \mathbbm{1}\{Y_{i,j}^{k} \leq z\}\right| \left| + \left|\sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} \mathbbm{1}\{Y_{i,j}^{k} \leq I_{m}\}\right|\right) \\ &+ \left|\sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} w_{k} \{G_{k}(z_{m}) - G_{k}(z)\}\right| \\ &\leq \left(\sqrt{\frac{2(e-t)}{n(e-s)(t-s)}} \lor \sqrt{\frac{2(t-s)}{n(e-s)(e-t)}}\right) \left(\left|\sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} \mathbbm{1}\{Y_{i,j}^{k} \leq I_{m}\}\right|\right) \\ &+ \left|\sum_{m=1,\dots,M-1}^{e} \left|\sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} \mathbbm{1}\{Y_{i,j}^{k} \in [z_{m} - \delta, z_{m+1} + \delta]\}\right| + \left|\sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} \mathbbm{1}\{Y_{i,j}^{k} \in I_{m}\}\right|\right) \\ &+ \left(\sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} |w_{k}|\right) \max_{k=s+1,\dots,e} |G_{k}(z) - G_{k}(z_{m})| \end{aligned}$$

$$\leq 4\left(\sqrt{\frac{2(e-t)}{n(e-s)(t-s)}} \vee \sqrt{\frac{2(t-s)}{n(e-s)(e-t)}}\right) \max_{m=1,\dots,M} B_m + \sqrt{\frac{2n(e-t)(t-s)}{e-s}} \delta C_G.$$
(46)

Since

$$\max_{m=1,...,M} B_{m} \leq \left| \sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} (\mathbb{1}\{Y_{i,j}^{k}\in I_{m}\} - \mathbb{P}\{\mathbb{1}\{Y_{i,j}^{k}\in I_{m}\}\}) \right| + \left| \sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} \mathbb{P}\{\mathbb{1}\{Y_{i,j}^{k}\in I_{m}\}\} \right| \\
\leq \left| \sum_{k=s+1}^{e} \sum_{(i,j)\in\mathcal{O}} (\mathbb{1}\{Y_{i,j}^{k}\in I_{m}\} - \mathbb{P}\{\mathbb{1}\{Y_{i,j}^{k}\in I_{m}\}\}) \right| + (e-s)n\delta C_{G},$$

and

$$\mathbb{P}\left\{\max_{m=1,\dots,M}\left|\sum_{k=s+1}^{e}\sum_{(i,j)\in\mathcal{O}}(\mathbb{1}\{Y_{i,j}^{k}\in I_{m}\}-\mathbb{P}\{\mathbb{1}\{Y_{i,j}^{k}\in I_{m}\}\})\right|\leq C_{9}\sqrt{\frac{n(e-s)\log(n\vee T)}{1-\rho}}\right\}$$

$$\geq 1-(n\vee T)^{-c_{9}},$$

where $C_9, c_9 > 0$ are universal constants, we have that

$$\mathbb{P}\left\{\left(\sqrt{\frac{2(e-t)}{n(e-s)(t-s)}} \lor \sqrt{\frac{2(t-s)}{n(e-s)(e-t)}}\right) \max_{m=1,\dots,M} B_m \\ \ge C_{10}T^{1/2}(1-\rho)^{-1/2}(\sqrt{d\log(n\vee T)}\lor d^{3/2})\right\} \le (n\lor T)^{-c_{10}}, \qquad (47)$$

where $C_{10}, c_{10} > 0$ are universal constants.

Combining (43), (44), (45), (46) and (47), the proof is complete.

Appendix C. Change point analysis lemmas

Lemma 17 Under Model 1, for any pair $(s, e) \subset (0, T)$ satisfying

$$\eta_{k-1} \le s \le \eta_k \le \ldots \le \eta_{k+q} \le e \le \eta_{k+q+1}, \quad q \ge 0,$$

let

$$b_1 \in \underset{b=s+1,\ldots,e-1}{\operatorname{arg\,max}} \widetilde{D}^b_{s,e}.$$

Then $b_1 \in \{\eta_1, \ldots, \eta_K\}$. Let $z \in \operatorname{argmax}_{x \in \mathbb{R}} |\widetilde{D}_{s,e}^b(x)|$. If $\widetilde{D}_{s,e}^t(z) > 0$ for some $t \in (s,e)$, then $\widetilde{D}_{s,e}^t(z)$ is either monotonic or decreases and then increases within each of the interval (s,η_k) , (η_k, η_{k+1}) , ..., $(\eta_{k+q}, e).$

This is identical to Lemma 7 in Padilla et al. (2019a) and we omit the proof here.

Lemma 18 Under Model 1, let $0 \le s < \eta_k < e \le T$ be any interval satisfying

$$\min\{\eta_k - s, e - \eta_k\} \ge c_1 \Delta_2$$

with $c_1 > 0$. Then we have that

$$\max_{t=s+1,\ldots,e-1} \widetilde{D}_{s,e}^t \ge \frac{2^{-3/2}c_1\kappa\Delta\sqrt{n}}{\sqrt{e-s}}.$$

Proof Recall that

$$G_{\eta_k}(z) = \mathbb{P}\left\{ (X_1(\eta_k))^\top X_2(\eta_k) \le z \right\}.$$

Let

$$z_0 \in \operatorname*{argmax}_{z \in [0,1]} |G_{\eta_k}(z) - G_{\eta_{k+1}}(z)|.$$

Without loss of generality, assume that $F_{\eta_k}(z_0) > F_{\eta_{k+1}}(z_0)$. For s < t < e, note that

$$\begin{split} \widetilde{D}_{s,e}^t(z_0) &= \left| \sqrt{\frac{n(e-t)}{2(e-s)(t-s)}} \sum_{k=s+1}^t G_k(z_0) - \sqrt{\frac{n(t-s)}{2(e-s)(e-t)}} \sum_{k=t+1}^e G_k(z_0) \right| \\ &= \left| \sqrt{\frac{n(e-s)}{2(t-s)(e-t)}} \sum_{k=s+1}^t \widetilde{G}_k(z_0) \right|, \end{split}$$

where $\widetilde{G}_k(z_0) = G_k(z_0) - (e-s)^{-1} \sum_{k=s+1}^e G_k(z_0)$. Under Model 1, it holds that $\widetilde{G}_{\eta_k}(z_0) > \kappa/2$. Therefore

$$\sum_{k=s+1}^{\eta_k} \tilde{G}_k(z_0) \ge (c_1/2)\kappa\Delta, \text{ and } \sqrt{\frac{n(e-s)}{2(t-s)(e-t)}} \ge \sqrt{\frac{n}{2(e-s)}}.$$

Then

$$\max_{t=s+1,\dots,e-1} \widetilde{D}_{s,e}^t \ge \frac{2^{-3/2}c_1\kappa\Delta\sqrt{n}}{\sqrt{e-s}}$$

Lemma 19 Under Model 1, if η_k is the only change point in (s, e), then

$$\widetilde{D}_{s,e}^{\eta_k} \le \kappa_k \sqrt{n/2} \min\{\sqrt{\eta_k - s}, \sqrt{e - \eta_k}\};$$
(48)

if $(s,e) \subset (0,T)$ contain two and only two change points η_k and η_{k+1} , then we have

$$\max_{t=s+1,\dots,e-1} \widetilde{D}_{s,e}^{\eta_k} \le \sqrt{n/2}\sqrt{e-\eta_{k+1}}\kappa_{k+1} + \sqrt{n/2}\sqrt{\eta_k-s}\kappa_k;$$
(49)

if $(s, e) \subset (0, T)$ contains two or more change points, including η_k and η_{k+1} , which satisfy that $\eta_k - s \leq c_1 \Delta$, for $c_1 > 0$, then

$$\widetilde{D}_{s,e}^{\eta_k} \le \sqrt{c_1} \widetilde{D}_{s,e}^{\eta_{k+1}} + \sqrt{2(\eta_k - s)n} \kappa_k.$$
(50)

Proof As for (48), it is due to that

$$\widetilde{D}_{s,e}^{\eta_k} = \sqrt{\frac{n(\eta_k - s)(e - \eta_k)}{2(e - s)}} \sup_{z \in \mathbb{R}} |G_{\eta_k}(z) - G_{\eta_k + 1}(z)| \le \kappa_k \sqrt{n/2} \min\{\sqrt{\eta_k - s}, \sqrt{e - \eta_k}\}.$$

Eq. (49) follows similarly.

As for (50), we consider the distribution sequence $\{H_t\}_{t=s+1}^e$ be such that

$$H_t = \begin{cases} G_{\eta_k+1}, & t = s + 1, \dots, \eta_k, \\ G_t, & t = \eta_k + 1, \dots, e. \end{cases}$$

For any s < t < e, define

$$\mathcal{H}_{s,e}^{t} = \sup_{z \in \mathbb{R}} \left| \mathcal{H}_{s,e}^{t}(z) \right|,$$

where

$$\mathcal{H}_{s,e}^{t}(z) = \sqrt{\frac{n(t-s)(e-t)}{2(e-s)}} \left\{ \frac{1}{t-s} \sum_{l=s+1}^{t} H_{l}(z) - \frac{1}{e-t} \sum_{l=t+1}^{e} H_{l}(z) \right\}.$$

For any $t \geq \eta_k$ and $z \in \mathbb{R}$, it holds that

$$\left| \widetilde{D}_{s,e}^{t}(z) - \mathcal{H}_{s,e}^{t}(z) \right| = \sqrt{\frac{2(e-t)}{n(e-s)(t-s)}} \frac{n(\eta_{k}-s)}{2} \left| G_{\eta_{k+1}}(z) - G_{\eta_{k}}(z) \right| \le \sqrt{\frac{n(\eta_{k}-s)}{2}} \kappa_{k}.$$

Thus we have

$$\begin{split} \widetilde{D}_{s,e}^{\eta_{k}} &= \sup_{z \in \mathbb{R}} \left| \widetilde{D}_{s,e}^{\eta_{k}}(z) - \mathcal{H}_{s,e}^{\eta_{k}}(z) + \mathcal{H}_{s,e}^{\eta_{k}}(z) \right| \leq \sup_{z \in \mathbb{R}} \left| \widetilde{D}_{s,e}^{\eta_{k}}(z) - \mathcal{H}_{s,e}^{\eta_{k}}(z) \right| + \mathcal{H}_{s,e}^{\eta_{k}} \\ &\leq \mathcal{H}_{s,e}^{\eta_{k}} + \sqrt{\frac{n(\eta_{k} - s)}{2}} \kappa_{k} \leq \sqrt{\frac{(\eta_{k} - s)(e - \eta_{k+1})}{(\eta_{k+1} - s)(e - \eta_{k+1})}} \mathcal{H}_{s,e}^{\eta_{k+1}} + \sqrt{\frac{n(\eta_{k} - s)}{2}} \kappa_{k} \\ &\leq \sqrt{c_{1}} \widetilde{D}_{s,e}^{\eta_{k+1}} + \sqrt{2n(\eta_{k} - s)} \kappa_{k}. \end{split}$$

Lemma 20 For any $z_0 \in \mathbb{R}$ and $(s, e) \subset (0, T)$ satisfying the following: there exits a true change point $\eta_k \in (s, e)$ such that

$$\min\{\eta_k - s, \, e - \eta_k\} \ge c_1 \Delta,\tag{51}$$

$$\widetilde{D}_{s,e}^{\eta_k}(z_0) \ge (c_1/2)\sqrt{n/2} \frac{\kappa \Delta}{\sqrt{e-s}},\tag{52}$$

where $c_1 > 0$ is a sufficiently small constant, and that

$$\max_{t=s+1,\dots,e} |\widetilde{D}_{s,e}^t(z_0)| - \widetilde{D}_{s,e}^{\eta_k}(z_0) \le 2^{-3/2} c_1^3 (e-s)^{-7/2} \Delta^4 \kappa \sqrt{n},$$
(53)

for all $d \in (s, e)$ satisfying

$$|d - \eta_k| \le c_1 \Delta/32,\tag{54}$$

it holds that

$$\widetilde{D}_{s,e}^{\eta_k}(z_0) - \widetilde{D}_{s,e}^d(z_0) > c |d - \eta_k| \Delta \widetilde{D}_{s,e}^{\eta_k}(z_0) (e - s)^{-2},$$

where c > 0 is a sufficiently small constant.

Proof The proof is identical to the proof of Lemma 11 in Padilla et al. (2019a) after letting $n_{\min} = n_{\max} = n/2$.

Lemma 21 Under Model 1, consider any generic $(s, e) \subset (0, T)$, satisfying

$$\min_{l=1,\dots,K} \min\{\eta_l - s, e - \eta_l\} \ge \Delta/16, \quad \eta_k \in (s, e).$$

and

 $e - s \le C_R \Delta.$

Let

$$\kappa_{s,e}^{\max} = \max_{\substack{l=1,\dots,K\\\eta_l \in (s,e)}} \kappa_l,$$

and $b \in \operatorname{argmax}_{s < t < e} D_{s,e}^t$. For some $c_1 > 0$ and $\gamma > 0$, suppose that

$$D_{s,e}^b \ge c_1 \kappa_{s,e}^{\max} \sqrt{\Delta n},\tag{55}$$

$$\max_{t=s+1,\ldots,e-1} \sup_{z\in\mathbb{R}} \left|\Delta_{s,e}^t(z)\right| \le \gamma,\tag{56}$$

and

$$\max_{0 \le s < e \le T} \sup_{z \in \mathbb{R}} \left| \sqrt{\frac{2}{n(e-s)}} \sum_{t=s+1}^{e} \sum_{i,j \in \mathcal{O}} \left(\mathbb{1}\{\widehat{Y}^t i, j \le z\} - G_t(z) \right) \right| \le \gamma.$$
(57)

If there exits a sufficiently small $0 < c_2 < c_1/2$ such that

$$\gamma \le c_2 \kappa_{s,e}^{\max} \sqrt{\Delta n},\tag{58}$$

then there exists a change point $\eta_k \in (s, e)$ such that

$$\min\{e - \eta_k, \eta_k - s\} \ge \Delta/4 \quad and \quad |\eta_k - b| \le C_\epsilon \frac{\gamma^2}{\kappa_k^2 n},$$

where $C_{\epsilon} > 0$ is a sufficiently large constant.

Proof

Without loss of generality, assume that $\widetilde{D}_{s,e}^b > 0$ and that $\widetilde{D}_{s,e}^t$ is locally decreasing at b. Observe that there has to be a change point $\eta_k \in (s, b)$, or otherwise $\widetilde{D}_{s,e}^b > 0$ implies that $\widetilde{D}_{s,e}^t$ is decreasing, as a consequence of Lemma 17. Thus, if $s \leq \eta_k \leq b \leq e$, then

$$\widetilde{D}_{s,e}^{\eta_k} \ge \widetilde{D}_{s,e}^b \ge D_{s,e}^b - \gamma \ge (c_1 - c_2) \kappa_{s,e}^{\max} \sqrt{\Delta n/2} \ge 2^{-3/2} c_1 \kappa_{s,e}^{\max} \sqrt{\Delta n},$$
(59)

where the second inequality follows from (56), and the third inequality follows from (55) and (58). Observe that $e - s \leq C_R \Delta$ and that (s, e) contains at least one change point.

Step 1. In this step, we are to show that

$$\min\{\eta_k - s, \, e - \eta_k\} \ge \min\{1, c_1^2\}\Delta/16.$$
(60)

Suppose that η_k is the only change point in (s, e). Then (60) must hold or otherwise it follows from (48) in Lemma 19, we have

$$D_{s,e}^{\eta_k} \le \kappa_k \sqrt{\Delta n} \frac{c_1}{4},$$

which contradicts (59).

Suppose (s, e) contains at least two change points. Then $\eta_k - s < \min\{1, c_1^2\}\Delta/16$ implies that η_k is the most left change point in (s, e). Therefore it follows from (50) that

$$\widetilde{D}_{s,e}^{\eta_k} \leq \frac{c_1}{4} \widetilde{D}_{s,e}^{\eta_{k+1}} + \sqrt{2n(\eta_k - s)} \kappa_k \leq \frac{c_1}{4} \max_{t=s+1,\dots,e} \widetilde{D}_{s,e}^t + \frac{c_1 \kappa_k \sqrt{n\Delta}}{4\sqrt{2}}$$

$$\leq \frac{c_1}{4} \max_{t=s+1,\dots,e} D_{s,e}^t + \frac{c_1}{4} \gamma + \frac{c_1 \kappa_k \sqrt{n\Delta}}{4\sqrt{2}}$$

$$< \max_{t=s+1,\dots,e} D_{s,e}^t - \gamma,$$
(61)

where the last inequality follows from that

$$\max_{t=s+1,\dots,e} D_{s,e}^t = D_{s,e}^b \ge 2^{-3/2} c_1 \kappa_{s,e}^{\max} \sqrt{\Delta n},$$

as implied by (59). Therefore, (61) contradicts

$$\widetilde{D}_{s,e}^{\eta_k} \ge \widetilde{D}_{s,e}^b - \gamma,$$

which is also implied by (59).

Step 2. It follows from Lemma 20 that

$$\widetilde{D}_{s,e}^{\eta_k} - \widetilde{D}_{s,e}^{\eta_k + c_1 \Delta/32} \ge c \frac{c_1 \Delta}{32} \Delta \widetilde{D}_{s,e}^{\eta_k} (e-s)^2 \ge \frac{cc_1}{32C_R^2} (c_1 \kappa \sqrt{\Delta n} - 2\gamma) \ge 2\gamma.$$
(62)

We claim that $b \in (\eta_k, \eta_k + c_1 \Delta/32)$. By contradiction, suppose that $b \ge \eta_k + c_1 \Delta/32$. Then

$$\widetilde{D}_{s,e}^{b} \le \widetilde{D}_{s,e}^{\eta_{k}+c_{1}\Delta/32} < \widetilde{D}_{s,e}^{\eta_{k}} - 2\gamma \le \max_{t=s+1,\dots,e} \widetilde{D}_{s,e}^{t} - 2\gamma \le \max_{t=s+1,\dots,e} D_{s,e}^{t} - \gamma = D_{s,e}^{b} - \gamma, \quad (63)$$

where the first inequality follows from Lemma 17, the second follows from (62), and the fourth follows from (56). Note that (63) shows that

$$\widetilde{D}_{s,e}^b < D_{s,e}^b - \gamma,$$

which is a contradiction with (59) showing that

$$\widetilde{D}_{s,e}^b \ge \widetilde{D}_{s,e}^b - \gamma.$$

Therefore we have $b \in (\eta_k, \eta_k + c_1 \Delta/32)$.

Step 3. This follows from the identical arguments as those in **Step 3** in the proof of Lemma 15 in Padilla et al. (2019a) by letting $n_{\min} = n_{\max} = n/2$ and translating notation appropriately. We have that

$$|b - \eta_k| \le C_\epsilon \frac{\gamma^2}{n\kappa_k^2}$$

where $C_{\epsilon} > 0$ is a universal constant.

Appendix D. Proof of Theorem 9

Proof [Proof of Theorem 9] Since ϵ is the upper bound of the localisation error, by induction, it suffices to consider any interval $(s, e) \subset (1, T)$ that satisfies

$$\eta_{k-1} \le s \le \eta_k \le \ldots \le \eta_{k+q} \le e \le \eta_{k+q+1}, \quad q \ge -1,$$

and

$$\max\{\min\{\eta_k - s, s - \eta_{k-1}\}, \min\{\eta_{k+q+1} - e, e - \eta_{k+q}\}\} \le \epsilon,$$

where q = -1 indicates that there is no change point contained in (s, e).

By Assumption 2, it holds that

$$\epsilon = C_{\epsilon} \frac{T \max\{d \log(n \lor T), d^3\}}{\kappa^2 n(1-\rho)} < \Delta/4.$$

It has to be the case that for any change point $\eta_k \in (0, T)$, either $|\eta_k - s| \leq \epsilon$ or $|\eta_k - s| \geq \Delta - \epsilon \geq 3\Delta/4$. This means that $\min\{|\eta_k - s|, |\eta_k - e|\} \leq \epsilon$ indicates that η_k is a detected change point in the previous induction step, even if $\eta_k \in (s, e)$. We refer to $\eta_k \in (s, e)$ an undetected change point if $\min\{|\eta_k - s|, |\eta_k - e|\} \geq 3\Delta/4$.

In order to complete the induction step, it suffices to show that we (i) will not detect any new change point in (s, e) if all the change points in that interval have been previous detected, and (ii) will find a point $b \in (s, e)$ such that $|\eta_k - b| \leq \epsilon$ if there exists at least one undetected change point in (s, e).

Recall the definitions $Y_{ij}^k = (X_i(k))^\top X_j(k)$ and $\hat{Y}_{ij}^k = (\hat{X}_i(k))^\top \hat{X}_j(k)$. For j = 1, 2, define the events

$$\mathcal{B}_{j}(\gamma) = \left\{ \max_{1 \le s < b < e \le T} \sup_{z \in [0,1]} \left| \sum_{k=s+1}^{e} w_{k}^{(j)} \sum_{(i,j) \in \mathcal{O}} \left(\mathbb{1}\{\widehat{Y}_{ij}^{k} \le z\} - \mathbb{E}\left\{ \mathbb{1}\{Y_{ij}^{k} \le z\} \right\} \right) \right| \le \gamma \right\},$$

where

and

$$w_k^{(1)} = \begin{cases} \sqrt{\frac{2}{n}} \sqrt{\frac{(e-b)}{(b-s)(e-s)}}, & k = s+1, \dots, b, \\ -\sqrt{\frac{2}{n}} \sqrt{\frac{(b-s)}{(e-b)(e-s)}}, & k = b+1, \dots, e, \end{cases}, \quad w_k^{(2)} = \sqrt{\frac{2}{n}} \frac{1}{\sqrt{e-s}},$$

$$\gamma = C_{\gamma} T^{1/2} \max\{\sqrt{d \log(n \vee T)}, d^{3/2}\},\$$

with a sufficiently large constant $C_{\gamma} > 0$.

Define

$$\mathcal{S} = \bigcap_{k=1}^{K} \left\{ \alpha_s \in [\eta_k - 3\Delta/4, \eta_k - \Delta/2], \, \beta_s \in [\eta_k + \Delta/2, \eta_k + 3\Delta/4], \text{ for some } s = 1, \dots, S \right\}.$$

It follows from Lemma 16 that for j = 1, 2, it holds that

$$\mathbb{P}\{\mathcal{B}_i\} \ge 1 - 11(n \lor T)^{-c} - 8Te^{-n}$$

The event S is studied in Lemma 13 in Wang et al. (2018b). The rest of the proof assumes the the event $\mathcal{B}_1(\gamma) \cap \mathcal{B}_2(\gamma) \cap S$.

Step 1. In this step, we will show that we will consistently detect or reject the existence of undetected change points within (s, e). Let a_m , b_m and m^* be defined as in Algorithm 2. Suppose there exists a change point $\eta_k \in (s, e)$ such that $\min\{\eta_k - s, e - \eta_k\} \ge 3\Delta/4$. In the event S, there exists an interval (α_m, β_m) selected such that $\alpha_m \in [\eta_k - 3\Delta/4, \eta_k - \Delta/2]$ and $\beta_m \in [\eta_k + \Delta/2, \eta_k + 3\Delta/4]$.

Following Algorithm 2, $(s_m, e_m) = (\alpha_m, \beta_m) \cap (s, e)$. We have that $\min\{\eta_k - s_m, e_m - \eta_k\} \ge (1/4)\Delta$ and (s_m, e_m) contains at most one true change point.

It follows from Lemma 18, with c_1 there chosen to be 1/4, that

$$\max_{s_m < t < e_m} \widetilde{D}_{s_m, e_m}^t \ge \frac{2^{-7/2} \kappa \Delta \sqrt{n}}{\sqrt{e-s}},$$

Therefore

$$a_m = \max_{s_m < t < e_m} D_{s_m, e_m}^t \ge \max_{s_m < t < e_m} \widetilde{D}_{s_m, e_m}^t - \gamma \ge 2^{-7/2} C_R^{-1/2} \kappa \sqrt{\Delta n} - \gamma.$$

Thus for any undetected change point $\eta_k \in (s, e)$, it holds that

$$a_{m^*} = \sup_{1 \le m \le S} a_m \ge 2^{-7/2} C_R^{-1/2} \kappa \sqrt{\Delta n} - \gamma \ge c_{\tau,2} \kappa \sqrt{\Delta n}, \tag{64}$$

where the last inequality is from the choice of γ and $c_{\tau,2} > 0$ is achievable with a sufficiently large C_{SNR} in Assumption 2. This means we accept the existence of undetected change points.

Suppose that there are no undetected change points within (s, e), then for any (s_m, e_m) , one of the following situations must hold.

(a) There is no change point within (s_m, e_m) ;

- (b) there exists only one change point $\eta_k \in (s_m, e_m)$ and $\min\{\eta_k s_m, e_m \eta_k\} \leq \epsilon_k$; or
- (c) there exist two change points $\eta_k, \eta_{k+1} \in (s_m, e_m)$ and $\eta_k s_m \leq \epsilon_k, e_m \eta_{k+1} \leq \epsilon_{k+1}$.

Observe that if (a) holds, then we have

$$\max_{s_m < t < e_m} D^t_{s_m, e_m} \le \max_{s_m < t < e_m} \widetilde{D}^t_{s_m, e_m} + \gamma = \gamma < \tau,$$

so no change points are detected.

Cases (b) and (c) are similar, and case (b) is simpler than (c), so we will only focus on case (c). It follows from Lemma 19 that

$$\max_{s_m < t < e_m} \widetilde{D}_{s_m, e_m}^t \leq \sqrt{n/2} \sqrt{e_m - \eta_{k+1}} \kappa_{k+1} + \sqrt{n/2} \sqrt{\eta_k - s_m} \kappa_k$$
$$\leq \sqrt{2C_\epsilon} T^{1/2} \max\{\sqrt{d\log(n \vee T)}, d^{3/2}\},$$

therefore

$$\max_{s_m < t < e_m} D^t_{s_m, e_m} \le \max_{s_m < t < e_m} \widetilde{D}^t_{s_m, e_m} + \gamma \le 2\gamma < \tau.$$

Under (6), we will always correctly reject the existence of undetected change points.

Step 2. Assume that there exists a change point $\eta_k \in (s, e)$ such that $\min\{\eta_k - s, \eta_k - e\} \ge 3\Delta/4$. Let s_m , e_m and m^* be defined as in Algorithm 2. To complete the proof it suffices to show that, there exists a change point $\eta_k \in (s_{m*}, e_{m*})$ such that $\min\{\eta_k - s_{m*}, \eta_k - e_{m*}\} \ge \Delta/4$ and $|b_{m*} - \eta_k| \le \epsilon$.

To this end, we are to ensure that the assumptions of Lemma 20 are verified. Note that (55) follows from (64), (56) and (57) follow from the definitions of events $\mathcal{B}_1(\gamma)$ and $\mathcal{B}_2(\gamma)$, and (58) follows from Assumption 2.

Thus, all the conditions in Lemma 20 are met. Therefore, we conclude that there exists a change point η_k , satisfying

$$\min\{e_{m^*} - \eta_k, \eta_k - s_{m^*}\} > \Delta/4 \tag{65}$$

and

$$|b_{m*} - \eta_k| \le C_\epsilon \frac{\gamma^2}{n\kappa_k^2} \le \epsilon,$$

where the last inequality holds from the choice of γ and Assumption 2.

The proof is completed by noticing that (65) and $(s_{m^*}, e_{m^*}) \subset (s, e)$ imply that

$$\min\{e - \eta_k, \eta_k - s\} > \Delta/4 > \epsilon.$$

As discussed in the argument before **Step 1**, this implies that η_k must be an undetected change point.

d	n	$ K - \hat{K} $	$d(\widehat{\mathcal{C}} \mathcal{C})$	$d(\mathcal{C} \widehat{\mathcal{C}})$
7	300	0.3	0.0	0.0
9	300	0.2	0.0	0.0
11	300	0.2	0.0	0.0
13	300	0.2	0.0	0.0
15	300	0.3	0.0	0.0
17	300	0.2	0.0	0.0
7	200	0.2	0.0	0.0
9	200	0.2	0.0	0.0
11	200	0.1	0.0	0.0
13	200	0.1	1.0	0.0
15	200	0.1	1.0	0.0
17	200	0.1	1.0	1.0
7	100	0.3	2.0	2.0
9	100	0.2	1.0	3.0
11	100	0.5	5.0	5.0
13	100	0.5	3.0	5.0
15	100	0.5	6.0	7.0
17	100	0.6	10.0	10.0

Table 5: Performance of NonPar-RDPG-CPD with data generated under Scenario 3 for varying values of d.

Appendix E. Sensitivity analysis of the input d

We proceed with the same setting as in Section 4.1, focusing on Scenario 3. The only difference with Section 4.1 is that now we explore the sensitivity of NonPar-RDPG-CPD to the choice d, by considering the performance of our algorithm for $d \in \{7, 9, 11, 13, 15, 17\}$. The results in Table 5 show that, overall, NonPar-RDPG-CPD is not sensitive to d, when it is not smaller than the true dimension of the latent positions.

Appendix F. Additional experiment on community structure changes only

In this section we consider an additional scenario to the ones described in Section 4. Keeping everything as in Section 4, with $\rho = 0.5$, we modify Scenario 1 by setting the matrix Q as

$$Q_{i,j} = \begin{cases} 0.5, & i, j \in \mathcal{C}_l, \ l \in \{1, 2, 3\}, \\ 0.3, & \text{otherwise.} \end{cases}$$

where $C_1 = B_1$, $C_2 = B_2$ and $C_3 = B_3 \cup B_4$, with B_1, \ldots, B_4 as in Scenario 1. We refer to the resulting model as Scenario 5, which consists of an example where the change happens in the community structure.

The results in Table 6 show that in the setting of Scenario 5 our proposed approach still outperforms the competing methods.

Method	n	$ K - \widehat{K} $	$d(\widehat{\mathcal{C}} \mathcal{C})$	$d(\mathcal{C} \widehat{\mathcal{C}})$
NonPar-RDPG-CPD	300	0.0	1.0	1.0
NBS	300	21.6	1.0	43
MNBS	300	0.9	0.0	20.0
NonPar-RDPG-CPD	200	0.1	3.0	3.0
NBS	200	21.2	1.0	43
MNBS	200	1.1	4.0	19.0
NonPar-RDPG-CPD	100	0.6	15.0	15.0
NBS	100	22.0	2.0	44.0
MNBS	100	1.1	5.0	20.0

Table 6: Scenario 5

Appendix G. Additional simulation results on varying κ

We now consider the setting of Scenario 3 and allow for an extra parameter σ^2 . Specifically, the data are now generated as follows. For $t \in \{1, 101\}$, we generate $Z_i(t) \stackrel{\text{ind}}{\sim} \mathcal{N}(0, \sigma^2 I_3)$, and for $t \in \mathcal{A}_1 \cup \mathcal{A}_3 \setminus \{1, 101\}$, we generate

$$Z_i(t) \begin{cases} \stackrel{\text{ind}}{\sim} \mathcal{N}(0, \sigma^2 I_3), & \text{with probability } 0.9, \\ = Z_i(t-1), & \text{with probability } 0.1. \end{cases}$$

We then set

$$P_{i,j}(t) = \frac{\exp\{Z_i(t)^{\top} Z_j(t)\}}{1 + \exp\{Z_i(t)^{\top} Z_j(t)\}}$$

Furthermore, we generate $P_{i,j}(51) \sim \text{Beta}(100, 100)$, and for $t \in \{52, \ldots, 100\}$ we generate

$$P(t) \begin{cases} = P(t-1), & \text{with probability 0.9,} \\ \sim \text{Beta}(100, 100), & \text{with probability 0.1.} \end{cases}$$

Once the mean matrices $\{P(t)\}_{t=1}^T \mathbb{R}^{n \times n}$ have been constructed, we independently draw $A_{i,j}(t) \sim \text{Ber}(P_{i,j}(t))$, for all $i, j \in \{1, \ldots, n\}$ and $t \in \{1, \ldots, T\}$. We consider experiments with $\sigma^2 \in \{1.5, 2, 2.5\}$. This additional parameter is meant to capture different levels of jump sizes κ .

For the model above and with the same setting from Section 4.1, the results in Tables 7–9 show that our method once again outperforms the competing approaches.

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Table 7: $\sigma^2 = 1.5$.

Method	n	$ K - \widehat{K} $	$d(\widehat{\mathcal{C}} \mathcal{C})$	$d(\mathcal{C} \widehat{\mathcal{C}})$
NonPar-RDPG-CPD	300	0.2	0.0	0.0
NBS	300	15.1	1.0	43.0
MNBS	300	1.1	28.0	36.0
NonPar-RDPG-CPD	200	0.52	0.0	0.0
NBS	200	14.0	1.0	44.0
MNBS	200	1.0	25.0	35.0
NonPar-RDPG-CPD	100	0.32	1.0	1.0
NBS	100	14.0	1.0	45.0
MNBS	100	1.0	25.0	34.0

Table 8: $\sigma^2 = 2.0$.

Method	n	$ K - \widehat{K} $	$d(\widehat{\mathcal{C}} \mathcal{C})$	$d(\mathcal{C} \widehat{\mathcal{C}})$
NonPar-RDPG-CPD	300	0.2	0.0	0.0
NBS	300	14.9	1.0	43.0
MNBS	300	1.2	21.0	37.0
NonPar-RDPG-CPD	200	0.3	0.0	0.0
NBS	200	14.4	1.0	43.0
MNBS	200	0.9	26.0	35.0
NonPar-RDPG-CPD	100	0.3	1.0	2.0
NBS	100	13.8	1.0	45.0
MNBS	100	0.9	22.0	35.0

Table 9: $\sigma^2 = 2.5$.

Method	n	$ K - \widehat{K} $	$d(\widehat{\mathcal{C}} \mathcal{C})$	$d(\mathcal{C} \widehat{\mathcal{C}})$
NonPar-RDPG-CPD	300	0.2	0.0	0.0
NBS	300	15.4	2.0	43.0
MNBS	300	1.2	21.0	35.0
NonPar-RDPG-CPD	200	0.4	0.0	0.0
NBS	200	14.7	1.0	43.0
MNBS	200	0.9	21.0	35.0
NonPar-RDPG-CPD	100	0.3	1.0	1.0
NBS	100	13.8	1.0	43.0
MNBS	100	0.9	27.0	36.0

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