IMPERIAL

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8th October, 2024

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Joint work with:

- Dynamic multiplex graphs
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 - Axel Gandy (Chair in Statistics, Imperial College London)
- Panels of multivariate time series / vector autoregressive processes
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 - Mihai Cucuringu (Associate Professor of Statistics, University of Oxford)
 - Alessandra Luati (Chair in Statistics, Imperial College London & University of Bologna)

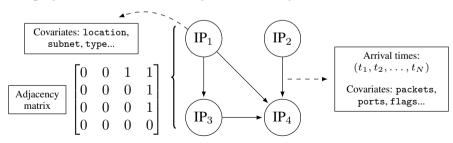
Acknowledgments:

 UK Research and Innovation / EPSRC ECR International Collaboration Grant 2024-2026 on "Spectral embedding methods and subsequent inference tasks on dynamic multiplex graphs" (EP/Y002113/1) in collaboration with Professor Carey Priebe (Johns Hopkins University)

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UNDIRECTED GRAPHS

- Undirected graph $\mathbb{G} = (V, E)$ where:
 - V is the **node set**, with cardinality n = |V|,
 - $E \subseteq V \times V$ is the **edge set**, containing dyads $(i, j) \in V \times V$.
- An edge is drawn if a node $i \in V$ connects to $j \in V$, written $(i, j) \in E$.
- From \mathbb{G} , an adjacency matrix $\mathbf{A} \in \{0,1\}^{n \times n}$, can be obtained via $A_{i,j} = \mathbb{1}_E\{(i,j)\}$.
- Real-world graphs tend to be more complex. For example:



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LATENT POSITION MODELS FOR GRAPHS

• Latent position models (Hoff, Raftery, and Handcock, 2002) for adjacency matrices:

$$oldsymbol{x}_i \overset{iid}{\sim} \mathcal{F} \quad o \quad \mathbb{P}(A_{i,j} = 1 \mid oldsymbol{x}_i, oldsymbol{x}_j) = \kappa(oldsymbol{x}_i, oldsymbol{x}_j) \quad o$$

- LPMs illustrate a powerful idea for network modelling: expressing edge-specific quantities through unobserved node features $x_i \in \mathcal{X} \subseteq \mathbb{R}^d$ sampled from a distribution \mathcal{F} .
- Node features are "linked" to link probabilities via a **kernel function** $\kappa: \mathcal{X} \times \mathcal{X} \to [0,1]$.
- $\bullet \ \, \textbf{Inner product kernels} \rightarrow \textbf{random dot product graph} \ (\textbf{RDPG}, \textbf{Athreya et al.}, 2018) : \\$

$$A_{i,j} \mid \boldsymbol{x}_1, \dots, \boldsymbol{x}_n \sim \text{Bernoulli}(\boldsymbol{x}_i^\intercal \boldsymbol{x}_j).$$

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RANDOM DOT PRODUCT GRAPHS

- RDPGs (and their generalisation, GRDPG, see Rubin-Delanchy et al., 2022) include:
 - Stochastic blockmodels (Holland, Laskey, and Leinhardt, 1983): $x_i = \mu_{z_i}$ for a community $z_i \in \{1,\ldots,K\}$, giving a between-community constant connection probability $B_{k\ell} = \mu_k^\mathsf{T} \mu_\ell$;
 - Degree-corrected stochastic blockmodels (Karrer and Newman, 2011): $x_i = \rho_i \mu_{z_i}$ for community $z_i \in \{1, \dots, K\}$ and degree-correction parameter $\rho_i \in (0, 1)$.
- The latent positions can be **estimated via the spectral decomposition** of **A**.

Definition (ASE - Adjacency spectral embedding)

For an integer $d \in \{1,\dots,n\}$ and a binary *symmetric* adjacency matrix $\mathbf{A} \in \{0,1\}^{n \times n}$, the d-dimensional adjacency spectral embedding (ASE) $\hat{\mathbf{X}} = [\hat{x}_1,\dots,\hat{x}_n]^{\mathsf{T}}$ of \mathbf{A} is

$$\hat{\mathbf{X}} = \mathbf{\Gamma} |\mathbf{\Lambda}|^{1/2} \in \mathbb{R}^{n \times d},$$

where Λ is a $d \times d$ diagonal matrix containing the d largest eigenvalues in magnitude, and Γ is a $n \times d$ matrix containing corresponding orthonormal eigenvectors.

MULTIPLEX GRAPHS

Introduction

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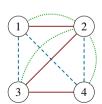
 In many real-world applications, edges can have different types. For example, links in cyber-security applications occur on different ports. In transportation networks, there are different means of transport between two locations. Edge types are usually called layers.

Vector autoregressive processes

• Undirected multiplex graph $\mathbb{G} = (V, \{E_1, \dots, E_K\})$ where:

Layer 1

- V is the **shared node set** across layers, with cardinality n = |V|,
- $E_k \subset V \times V$ is the **edge set** for the k-th layer, containing dyads $(i, j) \in V \times V$.
- Denote the adjacency matrix for the k-th layer as A_k [It is **not assumed** that $E_k \cap E_\ell = \emptyset$].



$$\mathbf{A}_2 = \left[egin{array}{cccc} 0 & 0 & 1 & 1 \\ 0 & 0 & 0 & 1 \\ 1 & 0 & 0 & 0 \\ 1 & 1 & 0 & 0 \end{array}
ight.$$

$$\mathbf{A}_1 = \left[\begin{array}{cccc} 0 & 1 & 0 & 0 \\ 1 & 0 & 1 & 0 \\ 0 & 1 & 0 & 1 \\ 0 & 0 & 1 & 0 \end{array} \right] \quad \mathbf{A}_2 = \left[\begin{array}{cccc} 0 & 0 & 1 & 1 \\ 0 & 0 & 0 & 1 \\ 1 & 0 & 0 & 0 \\ 1 & 1 & 0 & 0 \end{array} \right] \quad \mathbf{A}_3 = \left[\begin{array}{cccc} 0 & 1 & 0 & 0 \\ 1 & 0 & 1 & 1 \\ 0 & 1 & 0 & 0 \\ 0 & 1 & 0 & 0 \end{array} \right]$$

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SPECTRAL EMBEDDING OF MULTIPLEX GRAPHS

- Multiplex graphs can mainly be spectrally embedded via two methods: OMNI and UASE.
- Omnibus embedding (OMNI, Levin et al., 2017) Take ASE of the following block-matrix:

$$\tilde{\mathbf{A}} = \begin{bmatrix} \mathbf{A}_1 & (\mathbf{A}_1 + \mathbf{A}_2)/2 & \cdots & (\mathbf{A}_1 + \mathbf{A}_K)/2 \\ (\mathbf{A}_2 + \mathbf{A}_1)/2 & \mathbf{A}_2 & \cdots & (\mathbf{A}_2 + \mathbf{A}_K)/2 \\ \vdots & \vdots & \ddots & \vdots \\ (\mathbf{A}_K + \mathbf{A}_1)/2 & (\mathbf{A}_K + \mathbf{A}_2)/2 & \cdots & \mathbf{A}_K \end{bmatrix},$$

• Unfolded adjacency spectral embedding (UASE, Jones and Rubin-Delanchy, 2020; Gallagher, Jones, and Rubin-Delanchy, 2021) – Obtain the embedding from the singular value the composition of the unfolded matrix $\tilde{\mathbf{A}} \in \{0,1\}^{n \times nK}$, defined as follows:

$$\tilde{\mathbf{A}} = [\mathbf{A}_1 \ \mathbf{A}_2 \ \cdots \ \mathbf{A}_K].$$

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- ullet In real-world applications, graphs evolve over time o dynamic multiplex graphs.
- Dynamic multiplex graph $\mathbb{G} = (V, \{E_{k,t}\}_{k=1,\dots,K,\ t=1,\dots,T})$ where:
 - V is the **shared node set** across layers and time points, with cardinality n = |V|.
 - $E_{k,t} \subseteq V \times V$ is the **edge set** for the k-th layer at the t-th time.
 - Denote the adjacency matrix for the k-th layer at the t-th time as $\mathbf{A}_{k,t}$.
- We propose a **dynamic multiplex RDPG (DMP-RDPG)** model where each node is represented by latent positions in $\mathcal{X}_k \subseteq \mathbb{R}^d$ and $\mathcal{Y}_t \subseteq \mathbb{R}^d$, $k = 1, \dots, K, \ t = 1, \dots, T$.
 - ullet Positions $oldsymbol{x}_{i,k} \in \mathcal{X}_k$ are shared across time but are different across layers.
 - ullet Positions $oldsymbol{y}_{j,t} \in \mathcal{Y}_t$ are shared across layers but vary over time.
 - \bullet The connectivity model for nodes i and j at time t in layer k is given by:

$$A_{i,j,k,t} \sim \text{Bernoulli}\left(\boldsymbol{x}_{i,k}^{\intercal} \boldsymbol{y}_{j,t}\right).$$

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DMP-RDPG: THE SETUP

Introduction

• For K and T fixed, let $\mathcal{X}_1, \ldots, \mathcal{X}_K, \mathcal{Y}_1, \ldots, \mathcal{Y}_T \subseteq \mathbb{R}^d$ for some shared $d \in \mathbb{N}$, such that $x^{\mathsf{T}}y \in [0,1]$ for any $x \in \mathcal{X}_k$ and $y \in \mathcal{Y}_t$, for all $k = 1, \ldots, K$ and $t = 1, \ldots, T$.

Vector autoregressive processes

- Let \mathcal{F} be a distribution over the product space $[\bigotimes_{k=1}^K \mathcal{X}_k^n] \bigotimes [\bigotimes_{t=1}^T \mathcal{Y}_t^n]$.
- Let $x_{1,1},\ldots,x_{n,1},\ldots,x_{n,K},y_{1,1},\ldots,y_{n,1},\ldots,y_{n,T}\sim\mathcal{F}$. By organising these positions into matrices $\mathbf{X}_k = [\boldsymbol{x}_{1.k}^\intercal, \cdots, \boldsymbol{x}_{n.k}^\intercal]^\intercal$ and $\mathbf{Y}_t = [\boldsymbol{y}_{1.t}^\intercal, \dots, \boldsymbol{y}_{n.t}^\intercal]^\intercal$, the stacked matrices $\mathbf{X} = [\mathbf{X}_1^\intercal \mid \cdots \mid \mathbf{X}_K^\intercal]^\intercal \in \mathbb{R}^{nK \times d}$ and $\mathbf{Y} = [\mathbf{Y}_1^\intercal \mid \cdots \mid \mathbf{Y}_T^\intercal]^\intercal \in \mathbb{R}^{nT \times d}$ can be constructed.
- We define the $n \times n$ probability matrices for each time point and layer as $\mathbf{P}_{k,t} = \mathbf{X}_k \mathbf{Y}_t^{\mathsf{T}}$.
- Consider the following $nK \times nT$ double unfolding of the matrices $\mathbf{P}_{k,t}$ and $\mathbf{A}_{k,t}$:

$$\mathbf{P} = \begin{bmatrix} \mathbf{P}_{1,1} & \dots & \mathbf{P}_{1,T} \\ \vdots & \ddots & \vdots \\ \mathbf{P}_{K,1} & \dots & \mathbf{P}_{K,T} \end{bmatrix} = \mathbf{X}\mathbf{Y}^{\mathsf{T}}, \qquad \mathbf{A} = \begin{bmatrix} \mathbf{A}_{1,1} & \dots & \mathbf{A}_{1,T} \\ \vdots & \ddots & \vdots \\ \mathbf{A}_{K,1} & \dots & \mathbf{A}_{K,T} \end{bmatrix}.$$

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DMP-RDPG: THE DEFINITION

This setup leads to the definition of the dynamic multiplex RDPG (DMP-RDPG) model.

Definition (DMP-RDPG: Dynamic Multiplex Random Dot Product Graph)

 $(\mathbf{A}, \mathbf{X}, \mathbf{Y}) \sim \mathrm{DMP\text{-}RDPG}(\mathcal{F})$ if, conditional on \mathbf{X}_k and \mathbf{Y}_t , the matrix $\mathbf{A}_{k,t}$ has entries $A_{i,j,k,t} \sim \mathrm{Bernoulli}(P_{i,j,k,t})$ for all $i,j \in V$, with $P_{i,j,k,t} = \boldsymbol{x}_{i,k}^\mathsf{T} \boldsymbol{y}_{j,t}$.

- Key features:
 - Positions $x_{i,k} \in \mathcal{X}_k$ are shared across time but are different across layers.
 - Positions $y_{j,t} \in \mathcal{Y}_t$ are shared across layers but vary over time.
- Given a realisation A from a DMP-RDPG, the inferential objective is to estimate X and
 Y. We propose a doubly unfolded adjacency spectral embedding estimator.

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Doubly unfolded adjacency spectral embedding (DUASE)

Definition (DUASE – Doubly unfolded adjacency spectral embedding)

For a given integer $d \in \{1, \dots, n\}$ and dynamic multiplex graph adjacency matrices $\mathbf{A}_{k,t} \in \{0,1\}^{n \times n}, \ k=1,\ldots,K, \ t=1,\ldots,T,$ consider the doubly unfolded matrix

Vector autoregressive processes

$$\mathbf{A} = egin{bmatrix} \mathbf{A}_{1,1} & \dots & \mathbf{A}_{1,T} \ dots & \ddots & dots \ \mathbf{A}_{K,1} & \dots & \mathbf{A}_{K,T} \end{bmatrix}.$$

The d-dimensional doubly unfolded adjacency spectral embedding (DUASE) of $\bf A$ is:

$$\hat{\mathbf{X}} = \mathbf{U}\mathbf{D}^{1/2} \in \mathbb{R}^{nK \times d}, \qquad \hat{\mathbf{Y}} = \mathbf{V}\mathbf{D}^{1/2} \in \mathbb{R}^{nT \times d},$$

where **D** is a $d \times d$ diagonal matrix containing the d largest singular values of **A**, and **U** and V are $nK \times d$ and $nT \times d$ matrices containing the corresponding singular vectors.

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RESULTS: TWO-TO-INFINITY NORM BOUND FOR DUASE

This theorem is an adaptation of Theorem 2 in Jones and Rubin-Delanchy, 2020.

Theorem $(2 \to \infty \text{ norm bound for DUASE})$

Let $(\mathbf{A}, \mathbf{X}, \mathbf{Y}) \sim \text{DMPRDPG}(F_o)$ with K_n layers and T_n time points. Then, for each $k \in [K_n]$ and $t \in [T_n]$, there exist sequences of matrices \mathbf{W}_X and $\mathbf{W}_Y \in \mathrm{GL}(d)$ such that

Vector autoregressive processes

$$\|\hat{\mathbf{X}}^{k}\mathbf{W}_{X}^{-1} - \mathbf{X}^{k}\|_{2\to\infty} = O_{\mathbb{P}} \left\{ \frac{\log^{1/2}(n)}{\rho_{n}^{1/2} n^{1/2} T_{n}^{1/2}} \right\},$$
$$\|\hat{\mathbf{Y}}^{t}\mathbf{W}_{Y}^{-1} - \mathbf{Y}^{t}\|_{2\to\infty} = O_{\mathbb{P}} \left\{ \frac{\log^{1/2}(n)}{\rho_{n}^{1/2} n^{1/2} K_{n}^{1/2}} \right\}.$$

• $\|\mathbf{M}\|_{2\to\infty} = \sup_{\|\mathbf{z}\|_2=1} \|\mathbf{M}\mathbf{z}\|_{\infty}$; $\|\mathbf{M}\| = O_{\mathbb{P}}(f)$ if for any $\alpha > 0$ there exist a constant C>0 and an integer n^* such that for all $n\geq n^*$, $\mathbb{P}\{\|\mathbf{M}\|\leq Cf(n)\}\geq 1-n^{-\alpha}$.

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Underlying assumptions and sparsity considerations

• A global sparsity factor $\rho_n \in (0,1)$ is used to control the asymptotic connection density of the network as the number of nodes in the network n tends to infinity.

Vector autoregressive processes

- We assume that the sequence ρ_n converges either to 0 or to some constant c.
- The desired sparsity regime is enforced by defining $x_{i,k} = \rho^{1/2} \xi_{i,k}, \xi_{i,k} \sim F_{X,k}$ for some distribution $F_{X,k}$ on \mathbb{R}^d and $y_{i,t} = \rho^{1/2} \nu_{i,t}$, $\nu_{i,t} \sim F_{Y,t}$. Therefore, the joint distribution \mathcal{F} factorises into the product of n identical marginals for each $k=1,\ldots,K,\ t=1,\ldots,T$.
- We adopt the notation \mathcal{F}_{ρ} to refer to this scaled distribution.
- For our main results to hold in the asymptotic regime where K_n and T_n tend to infinity, we further require the existence of the $d \times d$ matrices $\tilde{\Delta}_X = \lim_{n \to \infty} K_n^{-1} \sum_{k=1}^{K_n} \Delta_{X,k}$ and $\tilde{\Delta}_Y = \lim_{n \to \infty} T_n^{-1} \sum_{t=1}^{T_n} \Delta_{Y,t}$. where $\tilde{\Delta}_{X,k} = \mathbb{E}[\boldsymbol{\xi}_{i,k} \boldsymbol{\xi}_{i,k}^{\intercal}]$ and $\tilde{\Delta}_{Y,t} = \mathbb{E}[\boldsymbol{\nu}_{j,t} \boldsymbol{\nu}_{i,t}^{\intercal}]$ are the second moment matrices of F_{Xk} and F_{Yt} respectively.

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RESULTS: CENTRAL LIMIT THEOREM FOR DUASE

• This theorem mirrors Theorem 3 in Jones and Rubin-Delanchy, 2020.

Theorem (CLT for DUASE)

Introduction

Let $(\mathbf{A}, \mathbf{X}, \mathbf{Y}) \sim \text{DMPRDPG}(F_{\rho})$ with K_n layers and T_n time points. Given latent positions $x \in \mathcal{X}_k$ and $y \in \mathcal{Y}_t$, then for all $z \in \mathbb{R}^d$ and for any fixed $i \in [n]$, $k \in [K_n]$ and $t \in [T_n]$ there exist sequences of matrices \mathbf{W}_X , $\mathbf{W}_Y \in \mathrm{GL}(d)$ (dependent on n) such that, for $n \to \infty$:

Vector autoregressive processes

$$\mathbb{P}\left\{n^{1/2}T_n^{1/2}(\hat{\mathbf{X}}_k\mathbf{W}_X^{-1} - \mathbf{X}_k)_i^{\intercal} \leq \boldsymbol{z} \mid \boldsymbol{\xi}_{i,k} = \boldsymbol{x}\right\} \to \Phi\left\{\boldsymbol{z}, \tilde{\boldsymbol{\Delta}}_Y^{-1}\mathbf{V}_Y(\boldsymbol{x})\tilde{\boldsymbol{\Delta}}_Y^{-1}\right\},$$

$$\mathbb{P}\left\{n^{1/2}K_n^{1/2}(\hat{\mathbf{Y}}_t\mathbf{W}_Y^{-1} - \mathbf{Y}_t)_i^{\intercal} \leq \boldsymbol{z} \mid \boldsymbol{\nu}_{i,t} = \boldsymbol{y}\right\} \to \Phi\left\{\boldsymbol{z}, \tilde{\boldsymbol{\Delta}}_X^{-1}\mathbf{V}_X(\boldsymbol{y})\tilde{\boldsymbol{\Delta}}_X^{-1}\right\},$$

where $\Phi(z, \Sigma)$ is the CDF of a d-dimensional normal distribution centered at 0 (the identically zero vector of dimension d), with covariance matrix $\Sigma \in \mathbb{R}^{d \times d}$, evaluated at $z \in \mathbb{R}^d$. The form of V_Y and V_X is analytically available.

DYNAMIC MULTIPLEX STOCHASTIC BLOCKMODELS (DMP-SBM)

- DMP-RDPG to define a dynamic multiplex stochastic blockmodel (DMP-SBM).
- Assume $A_{i,j,k,t} \sim \text{Bernoulli}(B_{z_{i,k},z'_{i,t},k,t})$, where $z_{i,k} \in \{1,\ldots,G\}$ and $z'_{i,t} \in \{1,\ldots,G'\}$ are group labels for nodes i and j in the k-th layer and t-th time point respectively, and $\mathbf{B} \in [0,1]^{G \times G' \times K \times T}$ is a tensor of probabilities of connections between groups.

Vector autoregressive processes

• Under a DMP-RDPG representation, $B_{h,\ell,k,t}=m{\mu}_{h,k}^\intercal m{\lambda}_{\ell,t}$ for $m{\mu}_{h,k}, m{\lambda}_{\ell,t} \in \mathbb{R}^d$, which gives:

$$A_{i,j,k,t} \sim \operatorname{Bernoulli}(\boldsymbol{\mu}_{z_{i,k},k}^{\intercal} \boldsymbol{\lambda}_{z'_{j,t},t}).$$

• The indicators $z_{i,k}$ and $z'_{i,t}$ can be estimated via Gaussian mixture modelling on the output of **DUASE**, using the theoretical guarantees provided by our **DUASE-CLT**:

$$\mathbb{P}\left\{n^{1/2}T_n^{1/2}\left(\mathbf{W}_X^{-1}\hat{\boldsymbol{x}}_{i,k} - \boldsymbol{\mu}_{g,k}\right)^{\mathsf{T}} \leq \boldsymbol{q}\right\} \to \Phi\left\{\boldsymbol{q}, \boldsymbol{\Sigma}_{X,g}\right\}, \quad g \in [G], \; \boldsymbol{q} \in \mathbb{R}^d,$$

$$\mathbb{P}\left\{n^{1/2}K_n^{1/2}\left(\mathbf{W}_Y^{-1}\hat{\boldsymbol{y}}_{i,t} - \boldsymbol{\lambda}_{g,t}\right)^{\mathsf{T}} \leq \boldsymbol{q}\right\} \to \Phi\left\{\boldsymbol{q}, \boldsymbol{\Sigma}_{Y,g}\right\}, \quad g \in [G'], \; \boldsymbol{q} \in \mathbb{R}^d.$$

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References

DMP-SBM + DUASE: A SIMULATION

Introduction

• Simulate a DMP-SBM with G=3, G'=4 and the following connection probabilities:

$$\mathbf{B}_{1,1} = \begin{bmatrix} 0.08 & 0.02 & 0.18 & 0.10 \\ 0.02 & 0.20 & 0.04 & 0.10 \\ 0.18 & 0.04 & 0.02 & 0.02 \\ 0.10 & 0.10 & 0.02 & 0.06 \end{bmatrix}, \quad \mathbf{B}_{1,2} = \begin{bmatrix} 0.16 & 0.16 & 0.04 & 0.10 \\ 0.16 & 0.16 & 0.04 & 0.10 \\ 0.04 & 0.04 & 0.09 & 0.02 \\ 0.10 & 0.10 & 0.02 & 0.06 \end{bmatrix}, \quad \mathbf{B}_{1,3} = \begin{bmatrix} 0.08 & 0.02 & 0.18 & 0.10 \\ 0.02 & 0.20 & 0.04 & 0.10 \\ 0.18 & 0.04 & 0.02 & 0.02 \\ 0.10 & 0.10 & 0.02 & 0.06 \end{bmatrix}$$

Vector autoregressive processes

$$\mathbf{B}_{2,1} = \begin{bmatrix} 0.08 & 0.02 & 0.18 & 0.10 \\ 0.02 & 0.20 & 0.04 & 0.10 \\ 0.18 & 0.04 & 0.02 & 0.02 \\ 0.10 & 0.10 & 0.02 & 0.06 \end{bmatrix}, \quad \mathbf{B}_{2,2} = \begin{bmatrix} 0.16 & 0.16 & 0.04 & 0.10 \\ 0.16 & 0.16 & 0.04 & 0.10 \\ 0.04 & 0.04 & 0.09 & 0.02 \\ 0.10 & 0.10 & 0.02 & 0.06 \end{bmatrix}, \quad \mathbf{B}_{2,3} = \begin{bmatrix} 0.08 & 0.02 & 0.18 & 0.10 \\ 0.02 & 0.20 & 0.04 & 0.10 \\ 0.18 & 0.04 & 0.02 & 0.02 \\ 0.10 & 0.10 & 0.02 & 0.06 \end{bmatrix}$$

$$\mathbf{B}_{3,1} = \begin{bmatrix} 0.08 & 0.08 & 0.08 & 0.08 & 0.08 \\ 0.08 & 0.08 & 0.08 & 0.08 & 0.08 \\ 0.08 & 0.08 & 0.08 & 0.08 & 0.08 \end{bmatrix}, \quad \mathbf{B}_{3,2} = \begin{bmatrix} 0.08 & 0.08 & 0.08 & 0.08 & 0.08 \\ 0.08 & 0.08 & 0.08 & 0.08 & 0.08 \\ 0.08 & 0.08 & 0.08 & 0.08 \end{bmatrix}, \quad \mathbf{B}_{3,3} = \begin{bmatrix} 0.08 & 0.08 & 0.08 & 0.08 \\ 0.08 & 0.08 & 0.08 & 0.08 \\ 0.08 & 0.08 & 0.08 \end{bmatrix}, \quad \mathbf{B}_{3,3} = \begin{bmatrix} 0.08 & 0.08 & 0.08 & 0.08 \\ 0.08 & 0.08 & 0.08 & 0.08 \\ 0.08 & 0.08 & 0.08 \end{bmatrix}$$

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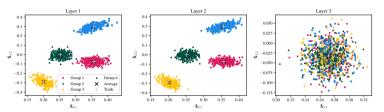


Figure 1. Left DUASE of a simulated DMP-SBM against the true latent positions (after orthogonal Procrustes rotation).

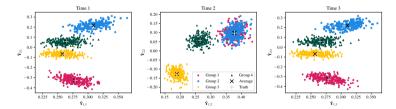


Figure 2. Right DUASE of a simulated DMP-SBM against the true latent positions (after orthogonal Procrustes rotation).

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References

Connections with literature on Euclidean mirrors

• Consider the right DUASE embedding $\hat{\mathbf{Y}} = [\hat{\mathbf{Y}}_1^\intercal \mid \cdots \mid \hat{\mathbf{Y}}_T^\intercal]^\intercal \in \mathbb{R}^{nT \times d}$ and extract the sequence of aligned time-specific embeddings $\hat{\mathbf{Y}}_1, \dots, \hat{\mathbf{Y}}_T$.

Vector autoregressive processes

• Athreya et al., 2024 calculates a dissimilarity (distance) matrix $\hat{\mathcal{D}}_{\phi}$ with entries:

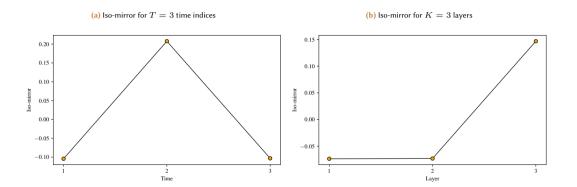
$$\hat{\mathcal{D}}_{\phi}(t,s) = \hat{d}_{\text{MV}}(\hat{\mathbf{Y}}_t, \hat{\mathbf{Y}}_s) = \min_{\mathbf{Q} \in \mathbb{O}(d)} \frac{1}{\sqrt{n}} ||\hat{\mathbf{Y}}_t - \hat{\mathbf{Y}}_s \mathbf{Q}||_2,$$

where $\mathbb{O}(d)$ is the orthogonal group with signature d.

- Since DUASE has **temporal stability**, one can simply set $\hat{\mathcal{D}}_{\phi}(t,s) = n^{-1/2} \|\hat{\mathbf{Y}}_t \hat{\mathbf{Y}}_s\|_2$.
- Apply CMDS to $\hat{\mathcal{D}}_{\phi}$ to estimate $\hat{\psi}(t) \in \mathbb{R}^c$ for $t = 1, \dots, T$, for some $c \in \mathbb{N}$.
- Apply ISOMAP to the points in $\text{CMDS}(\hat{\mathcal{D}}_{\phi}) = \{\hat{\psi}(t), t = 1, \dots, T\} \subset \mathbb{R}^c$ to obtain a 1-dimensional curve, which can be plotted against the time indices $t=1,\ldots,T$.
- This gives a joint (across layers) Euclidean mirror for dynamic multiplex networks.

Iso-MIRROR ON SBM WITH DUASE

Introduction



Vector autoregressive processes

Figure 3. Iso-mirrors calculated from DUASE on the simulated DMP-SBM in previous slides.

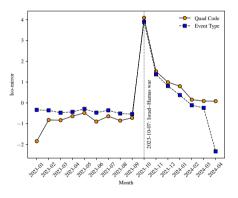
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APPLICATIONS ON REAL-WORLD MULTIPLEX NETWORKS

- PLOVER dataset (Halterman et al., 2023)
 - Consider the Political Language Ontology for Verifiable Event Records (PLOVER) dataset, consisting in 624,888 political interaction events between n = 104 countries.
 - We group the events across T=16 months ranging between January 2023 and April 2024. Each event is associated with one of K=16 event types based on the PLOVER categories.
 - Each of the event types is further grouped into $K^* = 4$ macro-groups called *quad categories*: material cooperation, verbal cooperation, verbal conflict, and material conflict.
- FinDKG dataset (Li and Sanna Passino, 2024)
 - The graph has a total of 241.948 edges between n=13.637 nodes, with K=15 different connection types related to financial concepts, such as "Raise", "Invests In" or "Produce".
 - Nodes represent financial institutions, politicians, businessmen, countries, financial concepts, and commodities.

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(a) Iso-mirror on right DUASE $\hat{\mathbf{Y}}$



(b) Iso-mirror on $\hat{\mathbf{X}}$, grouped by event type

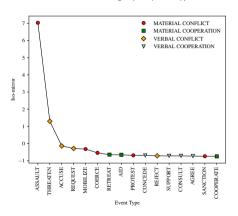
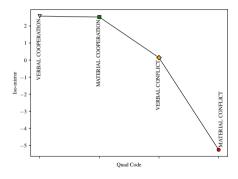


Figure 4. Iso-mirror across time and event types on the POLECAT data.

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PLOVER DATASET





(b) Iso-mirror on $\hat{\mathbf{X}}$ vs. PLOVER intensity

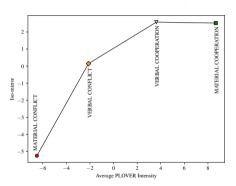
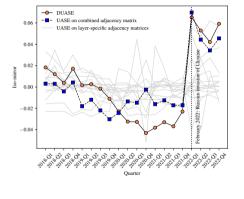


Figure 5. Iso-mirror across time and event types on the POLECAT data, grouped by quad-code.

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FINDKG DATASET





(b) Iso-mirror for left DUASE $\hat{\mathbf{X}}$

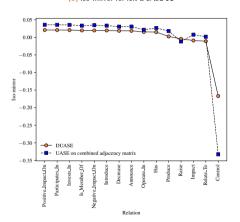


Figure 6. Iso-mirror across time and event types on the FinDKG data.

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- Panels of multivariate time series $\{\mathbf{X}_t = (X_{1,t},...,X_{N,t})^{\mathsf{T}}, \ X_{i,t} \in \mathbb{R}\}_{t \in \mathbb{Z}}$ exhibiting co-movement between components are central to many scientific disciplines such as environmental science, econometrics, and neuroscience.
- Often, $X_{i,t}$ depends not only on its own past values, but also on the **past values of a** subset of other panel components, $\{X_{j,s}: j \subseteq [N], s < t\}$, where $[N] = \{1, \ldots, N\}$.
- Vector autoregression (VAR) is a widely used model for multivariate time series:

$$\mathbf{X}_t = \mathbf{\Phi} \mathbf{X}_{t-1} + \boldsymbol{\varepsilon}_t, \ \boldsymbol{\varepsilon}_t \sim \mathbb{N}_N(\mathbf{0}_N, \boldsymbol{\Sigma}_{\varepsilon}),$$

where $\mathbf{\Phi} \in \mathbb{R}^{N imes N}$ is a matrix of **coefficients**, for some covariance structure $\mathbf{\Sigma}_{arepsilon}$.

• For large N, modelling via the VAR framework becomes **prohibitive** as the number of model parameters grows as $O(N^2)$ and can quickly exceed the number of observations.

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- Factor models The large panel of time series are modelled as stemming from a relatively small number of common latent factors (Stock and Watson, 2002).
- Factors with sparse regression Fan, Masini, and Medeiros (2023) combine the dimensionality reduction of factor modelling with the parsimony of sparse linear regression and give a novel test for covariance structure. Their proposed model is called the Factor Augmented Regression Model (FARM).
- Network VAR Knight et al. (2020) introduce generalise network autoregression (GNAR) which, given an observed network, fits a flexible network autoregressive model. Barigozzi, Cho, and Owens (2023) propose an L1-regularised Yule-Walker method for estimating a factor adjusted, idiosyncratic VAR model (FNETS).
- Community detection: Guðmundsson and Brownlees (2021) use estimated VAR coefficients to embed and cluster the panel components.

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NETWORK-INFORMED RESTRICTED VECTOR AUTOREGRESSIVE (NIRVAR) MODEL

Vector autoregressive processes

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Definition (NIRVAR - Network-informed restricted vector autoregressive model)

For some fixed $q \in [Q]$, let $\{\mathbf{X}_t^{(q)}\}_{t \in \mathbb{Z}}$ denote a zero mean, second order stationary stochastic process where $\mathbf{X}_t^{(q)} = (X_{1,t}^{(q)},\cdots,X_{N,t}^{(q)})^\intercal \in \mathbb{R}^N$ and $q \in [Q]$. The NIRVAR model for the *a*-th feature is

$$\mathbf{X}_t^{(q)} = \sum_{r=1}^{Q} (\mathbf{A}^{(r)} \odot \tilde{\mathbf{\Phi}}^{(r)}) \mathbf{X}_{t-1}^{(r)} + \boldsymbol{\varepsilon}_t^{(q)}, \qquad \boldsymbol{\varepsilon}_t^{(q)} \sim \mathbb{N}_N(\mathbf{0}_N, \sigma^2 \mathbf{I}_{N \times N}),$$

in which $(\mathbf{A}^{(r)}, \mathbf{Y}^{(r)}) \sim \mathrm{SBM}(\mathbf{B}^{(r)}, \pi^{(r)}), \ r \in [Q], \ \text{and} \ \tilde{\mathbf{\Phi}}^{(r)}, r \in [Q], \ \text{is an} \ N \times N$ matrix of fixed weights. Defining $\mathbf{\Phi}^{(r)} = \mathbf{A}^{(r)} \odot \tilde{\mathbf{\Phi}}^{(r)}$ and $\mathbf{\Phi} = (\mathbf{\Phi}^{(1)}| \cdots |\mathbf{\Phi}^{(Q)}), \ \text{we write}$ $\mathbf{X}_{t}^{(q)} \sim \text{NIRVAR}(\mathbf{\Phi}).$

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NIRVAR MODEL

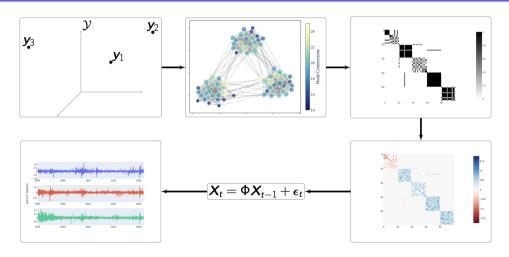
ullet Defining ${f \Phi}^{(r)}={f A}^{(r)}\odot ilde{f \Phi}^{(r)},$ we can write the NIRVAR model as

$$\mathbf{X}_t^{(q)} = \sum_{r=1}^Q \mathbf{\Phi}^{(r)} \mathbf{X}_{t-1}^{(r)} + \mathbf{arepsilon}_t^{(q)}, \qquad \mathbf{arepsilon}_t^{(q)} \sim \mathbb{N}_N(\mathbf{0}_N, \sigma^2 \mathbf{I}_{N imes N}),$$

which allows the interpretation of the NIRVAR model as a **restricted VAR whose restrictions** are determined by the graph. Note that the restrictions are static.

• Equivalently, we can define the coefficient matrices $\Phi^{(r)}$ as the adjacency matrix resulting from a **weighted SBM** (**WSBM**; Gallagher, Jones, and Rubin-Delanchy, 2021), for a suitable choice of the weight distribution. A **zero-inflated** component is added to the distribution to obtain sparsity.

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NIRVAR: ESTIMATION

- Recovering the **complete edge** set is difficult. Instead, we aim to **recover the community memberships** z_i , i = 1, ..., N for each node.
- The two step estimation approach is:
 - ① Estimate \hat{z}_i by performing clustering on an embedding $\hat{\mathbf{Y}} \in \mathbb{R}^{N \times d}$ of the panel components, and define the binary matrix $\hat{\mathbf{A}} = \{\hat{A}_{i,j}\} \in \{0,1\}^{N \times N}$ with the following entries:

$$\hat{A}_{i,j} = \mathbb{1} \left\{ \hat{z}_i = \hat{z}_j \right\},\,$$

where $\mathbb{1}\{\cdot\}$ is the indicator function.

- ② Set $\hat{\Phi}_{i,j} = 0$ if $\hat{A}_{i,j} = 0$ and estimate the remaining **unrestricted parameters** (corresponding to $\hat{A}_{i,j} = 1$) via ordinary least squares (OLS).
- If multiple features are used, the clustering can be done on **feature-specific embeddings** $\hat{\mathbf{Y}}^{(q)} \in \mathbb{R}^{N \times d}$, and the same procedure is followed to obtain $\hat{\mathbf{\Phi}}^{(q)}$.
- ullet The one-step-ahead prediction equation becomes: $\mathbf{X}_{t+1}^{(q)} = \sum_{r=1}^Q \hat{\mathbf{\Phi}}^{(r)} \mathbf{X}_t^{(r)}.$

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- Let $\mathfrak{X}^{(q)} = (\boldsymbol{x}_1^{(q)}, \dots, \boldsymbol{x}_T^{(q)})$ be the $N \times T$ design matrix of feature q, where $\boldsymbol{x}_t^{(q)} = (x_{1,t}^{(q)}, \dots, x_{N,t}^{(q)})'$ is a realisation of the random variable $\mathbf{X}_t^{(q)}$.
- To construct an embedding $\hat{y}_i^{(q)} \in \mathbb{R}^d$ we use **unfolded adjacency spectral embedding** (UASE, Jones and Rubin-Delanchy, 2020), which obtains embeddings $\hat{\mathbf{Y}}^{(q)} \in \mathbb{R}^{N \times d}$ by considering the SVD of $\tilde{\mathbf{A}} = (\mathbf{A}^{(1)}|\cdots|\mathbf{A}^{(Q)})$.
- UASE has key stability properties (Gallagher, Jones, and Rubin-Delanchy, 2021): it
 assigns the same position, up to noise, to vertices behaving similarly for a given feature
 (cross-sectional stability) and a constant position, up to noise, to a single vertex behaving
 similarly across different features (longitudinal stability).
- We consider the SVD of $\tilde{\mathbf{S}} = (\mathbf{S}^{(1)}|\cdots|\mathbf{S}^{(Q)}) \in \mathbb{R}^{N \times NQ}$ where $\mathbf{S}^{(q)} = \mathfrak{X}^{(q)}\mathfrak{X}^{(q)^{\mathsf{T}}}/T \in \mathbb{R}^{N \times N}$ is the sample covariance matrix for feature q.

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- Estimate the dimension of the latent space as the number of singular values of \tilde{S} that are greater than the Marčencko-Pastur distribution cutoff (Marčenko and Pastur, 1967).
- Let $\mathbf{D} \in \mathbb{R}^{\hat{d} \times \hat{d}}$ and $\mathbf{U} \in \mathbb{R}^{N \times \hat{d}}$ be the matrices containing the \hat{d} largest singular values and corresponding left singular vectors of $\tilde{\mathbf{S}}$. The corresponding **joint** embedding is:

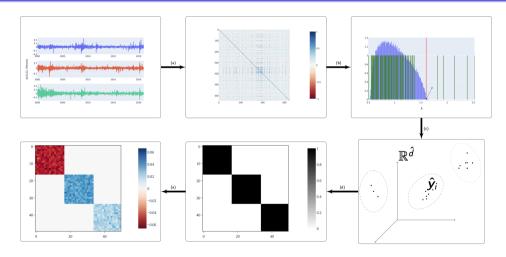
$$\hat{\mathbf{Y}} = \mathbf{U}\mathbf{D}^{1/2} \in \mathbb{R}^{N \times \hat{d}}.$$

- Cluster each $\hat{y}_i \in \mathbb{R}^{\hat{d}}$ into $K = \hat{d}$ groups using a Gaussian mixture model.
- Feature-specific embeddings $\hat{\mathbf{Y}}^{(q)} \in \mathbb{R}^{N \times \hat{d}}$ can be obtained by unstacking $\mathbf{V}\mathbf{D}^{1/2} \in \mathbb{R}^{NQ \times \hat{d}}$ into Q equal blocks $(\hat{\mathbf{Y}}^{(1)}; \dots; \hat{\mathbf{Y}}^{(Q)})$, where \mathbf{V} contains the right singular vectors of $\tilde{\mathbf{S}}$ corresponding to the \hat{d} largest singular values.

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NIRVAR: EMBEDDING ϕ ESTIMATION

Introduction



32/43 Imperial College London

Francesco Sanna Passino

WEIGHTED STOCHASTIC BLOCKMODELS

- How does $\hat{\boldsymbol{y}}_i^{(q)}$ relate to the ground truth positions $\boldsymbol{\theta}_{z_i}^{(q)}$? We derive a connection between the covariance $\boldsymbol{\Gamma}^{(q)} = \mathbb{E}\{(\mathbf{X}_t^{(q)})(\mathbf{X}_t^{(q)})^{\intercal}\}$ of a NIRVAR (Φ) process and $\boldsymbol{y}_i^{(q)}$ for Q=1.
- ullet If $oldsymbol{\Phi}$ is symmetric, the rank-d spectral embedding of $oldsymbol{\Gamma}$ and $oldsymbol{\Phi}$ are equivalent.

Proposition

Let $\mathbf{X}_t \sim \mathrm{NIRVAR}(\mathbf{\Phi})$ where $\mathbf{\Phi}$ is assumed to be symmetric. Consider the eigendecomposition $\mathbf{\Phi} = \mathbf{U}_{\mathbf{\Phi}} \mathbf{\Lambda}_{\mathbf{\Phi}} \mathbf{U}_{\mathbf{\Phi}}^{\mathsf{T}} + \mathbf{U}_{\mathbf{\Phi},\perp} \mathbf{\Lambda}_{\mathbf{\Phi},\perp} \mathbf{U}_{\mathbf{\Phi},\perp}^{\mathsf{T}}$, where $\mathbf{U}_{\mathbf{\Phi}} \in \mathbb{O}(N \times d)$ and $\mathbf{\Lambda}_{\mathbf{\Phi}}$ is a $d \times d$ diagonal matrix comprising the d largest eigenvalues in absolute value of $\mathbf{\Phi}$. Then the rank d truncated eigendecomposition of the covariance matrix $\mathbf{\Gamma} = \mathbb{E}(\mathbf{X}_t \mathbf{X}_t^{\mathsf{T}})$ is $\mathbf{\Gamma} = \mathbf{U}_{\mathbf{\Phi}} \mathbf{\Lambda}_{\mathbf{\Gamma}} \mathbf{U}_{\mathbf{\Phi}}^{\mathsf{T}}$ in which $\mathbf{\Lambda}_{\mathbf{\Gamma}}$ is a $d \times d$ diagonal matrix with diagonal elements $(\lambda_{\mathbf{\Gamma}})_i = 1/\{1 - (\lambda_{\mathbf{\Phi}})_i^2\}$ where $(\lambda_{\mathbf{\Phi}})_i$ is the corresponding diagonal entry of $\mathbf{\Lambda}_{\mathbf{\Phi}}$.

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MACROECONOMIC APPLICATIONS: FRED-MD DATASET

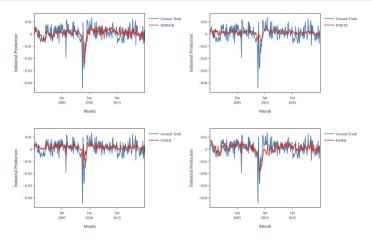
- FRED-MD is a publicly accessible database of monthly observations of macroeconomic variables.
- The prediction task is one-step ahead forecasts of the first order difference of the logarithm of the monthly industrial production (IP) index.
- We backtest NIRVAR, FARM, FNETS, and GNAR from January 2000 December 2019 using a rolling window framework with a lookback window of 480 observations.

| Metric | NIRVAR | FARM | FNETS | GNAR |
|-------------|--------|--------|--------|--------|
| Overall MSE | 0.0087 | 0.0089 | 0.0096 | 0.0101 |

Table 1. Overall MSE of each model for the task of forecasting US industrial production.

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Macroeconomic applications: FRED-MD dataset



Vector autoregressive processes

Figure 7. Predicted (log-differenced) IP against the realised (log-differenced) IP for each model.

- The previous close-to-close (pvCLCL) market excess returns of 648 financial assets from 03/01/2000 31/12/2020 were derived from databases provided by the CRSP.
- The task is to predict the sign of the next day pvCLCL market excess returns. A positive (negative) sign corresponds to a long (short) position in the asset.
- We backtest NIRVAR, FARM, FNETS, and GNAR using a rolling window from 01/01/2004 -31/12/2020 with a look-back window of four years.

| | NC1 | NC2 | NP1 | NP2 | FARM | FNETS | GNAR |
|---------------------|-------|-------|-------|-------|-------|-------|-------|
| Sharpe Ratio | 2.50 | 2.34 | 2.82 | 2.69 | 0.22 | 0.78 | 0.70 |
| Mean Absolute Error | 0.012 | 0.013 | 0.012 | 0.013 | 0.012 | 0.012 | 0.012 |

Table 2. Statistics on the financial returns predictive performance over the backtesting period.

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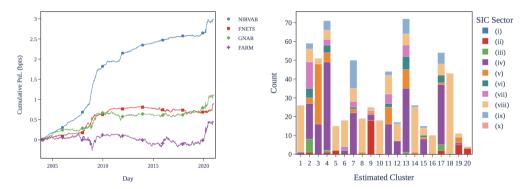


Figure 8. Left: the cumulative PnL in bpts over the backtesting period. Right: comparison of the NIRVAR estimated clusters with the SIC groups on 31/12/2020.

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- The first differences of the log daily number of bicycle rides from N=774 Santander Cycles stations in central London from 07/03/2018 until 10/03/2020 (T=735) were obtained using records from TfL Open Data (see https://cycling.data.tfl.gov.uk/).
- There were K = 7 clusters estimated by NIRVAR, differentiated by their mean number of bicycle rides as well as by the change in the number of bicycle rides on weekdays compared with weekends.

| Model | NIRVAR | FARM | FNETS | GNAR |
|-------|--------|-------|-------|-------|
| MSPE | 0.364 | 0.370 | 0.388 | 0.374 |

Table 3. Mean squared prediction errors (MSPE) for the Santander cycles data.

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SANTANDER CYCLES DATA

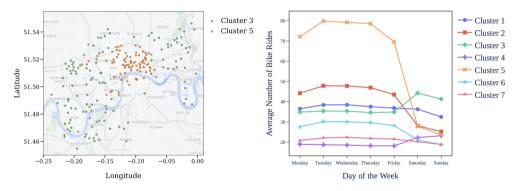


Figure 9. NIRVAR clusters on the Santander Cycles data.

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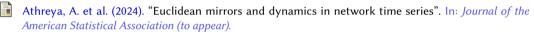
Conclusion

- This talk introduces two different models and related estimators:
 - OMP-RDPG + DUASE for dynamic multiplex graphs
 - Propose a DMP-RDPG model for dynamic multiplex graphs where $A_{i,j,t,k} \sim \text{Bernoulli}(\boldsymbol{x}_{i,k}^\intercal \boldsymbol{y}_{j,t})$.
 - Positions $x_{i,k} \in \mathcal{X}_k$ are shared across time but are different across layers.
 - Positions $y_{j,t} \in \mathcal{Y}_t$ are shared across layers but vary over time.
 - Estimates of the latent position are obtained via a double unfolding of the adjacency matrices.
 - NIRVAR model and estimator for panels of multivariate time series
 - We model the time series as a VAR process whose parameter matrix is a realisation of a WSBM.
 - We introduce an estimation framework that firstly determines the restrictions to be placed on the VAR parameters and secondly estimates the remaining unrestricted parameters via OLS.
 - The method can be used to estimate VAR parameters when the underlying network is **unobserved**.
 - More details can be found in Martin et al. (2024) (preprint available on arXiv).
 - Martin, B. et al. (2024). "NIRVAR: Network Informed Restricted Vector Autoregression". In: *arXiv e-prints*. arXiv: 2407.13314.

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Francesco Sanna Passino Imperial College London Low-rank models for dynamic multiplex graphs and vector autoregressive processes

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