

# Shrinkage in Space

Spillovers and Networks in a Hierarchical Model

Nikolas Kuschnig

Causal Panel Data Conference, Stanford GSB

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Vienna University of Economics and Business

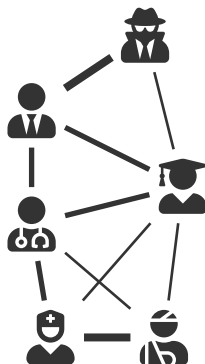
[nikolas.kuschnig@wu.ac.at](mailto:nikolas.kuschnig@wu.ac.at)

# Motivation

Economic activities rarely occur in isolation — agents are **embedded in networks** and **experience spillovers**.<sup>a</sup>

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<sup>a</sup>See, e.g., Akcigit et al., 2021; Alfaro-Ureña et al., 2022; Ambrus and Elliott, 2021; Canen et al., 2023; Chetty et al., 2022; Dhyne et al., 2021; Giovanni et al., 2022; Vom Lehn and Winberry, 2022; Weidmann and Deming, 2021.

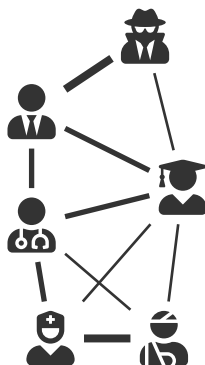


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## The issue

We **rarely observe** the networks behind spillovers, and models suffer from the **curse of dimensionality**.

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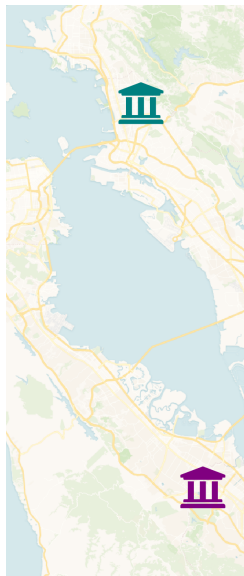
# Overview

With networks unknown, models rely on **assumptions** and **approximate information**.

*How far is **Berkeley** from **Stanford**?*

*Who are your five best friends?*

*Who do you ask for advice?*



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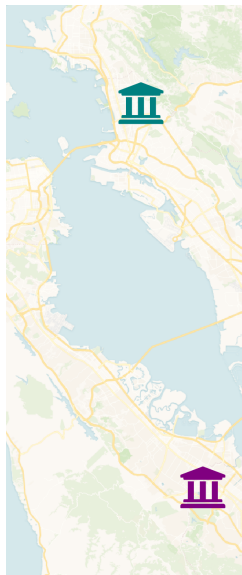
*How far is Berkeley from Stanford?*

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Today, I will show

- that these restrictions **distort inference**, and
- how to address this with a **Bayesian approach**.



# Contributions and literature

Today, I'll focus on the *main contribution* to a growing literature<sup>a</sup> — a **Bayesian hierarchical approach** to model **spillovers** and **latent networks** behind them.

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Compared to the literature, my approach

- **flexibly leverages information** of all kinds,
- naturally conveys uncertainty via full posteriors,
- is generally applicable.

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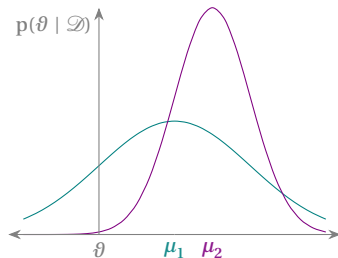
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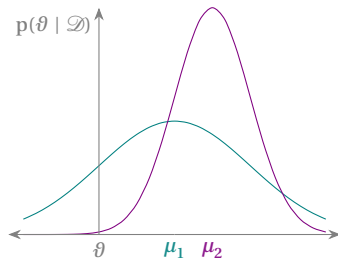
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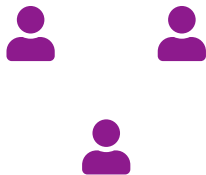
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# Setting

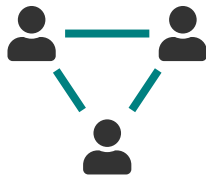
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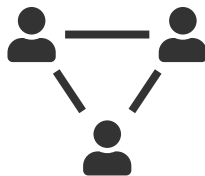
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We want to **learn about the relationship**

$$Y = f(X, \mathcal{G}) + \varepsilon,$$

and will need to *impose some structure* on  $f$  and  $\mathcal{G}$ .

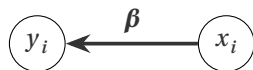


# A model for $f$

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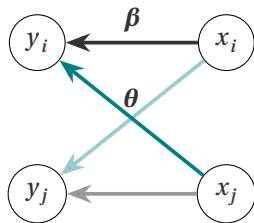
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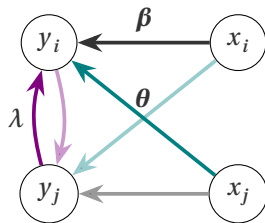
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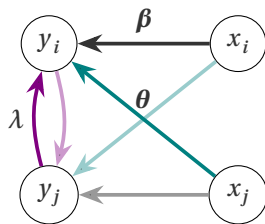
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## Linear network model

The network is represented by  $\mathbf{W}$ . Special cases are the *linear-in-means* and *spatial Durbin* models, which constrain  $\mathbf{W}$  and treat it as given.



# Does the network matter?

Consider *network effects*<sup>a</sup> based on

1. **contiguity of US states**, proxied with
2. averages of contiguous states, and
3. distance-decay between centers.



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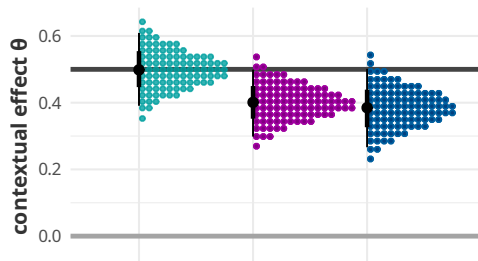
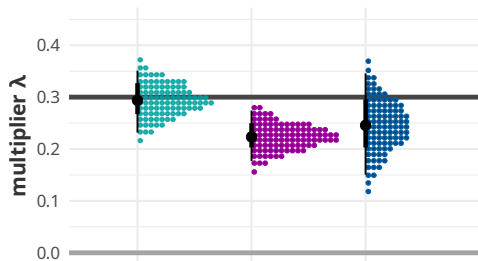
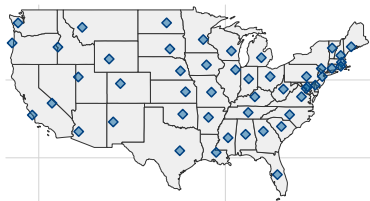
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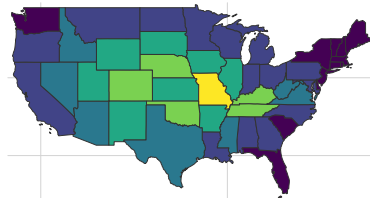
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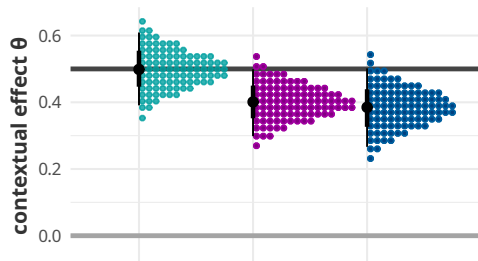
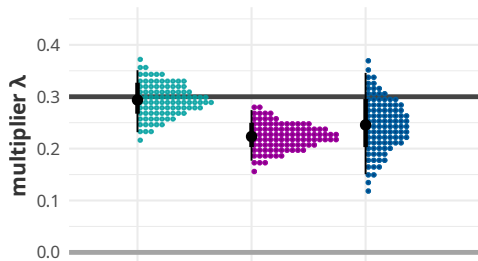
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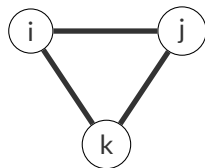


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 $\mathcal{G} = \{\mathcal{A}, \mathcal{E}\}$ , which we allow to be



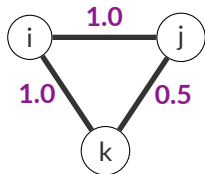
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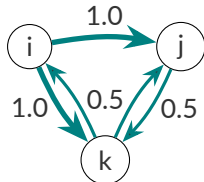
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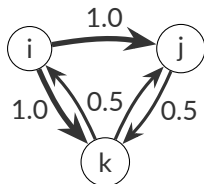
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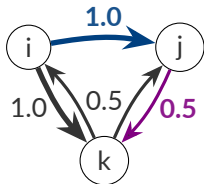
## Adjacency matrix

The graph corresponds to the matrix  $\mathbf{G}$  with entries given by  $g_{ij} = g(i, j)$ .

$$\mathbf{G} = \begin{bmatrix} 0 & g_{12} & \cdots & g_{1n} \\ g_{21} & 0 & \cdots & g_{2n} \\ \vdots & \vdots & \ddots & \vdots \\ g_{n1} & g_{n2} & \cdots & 0 \end{bmatrix}.$$

# The normalized adjacency matrix

In practice, a **normalized adjacency matrix  $\mathbf{W}$**  is used, such that  $\lambda$  and  $\theta$  are identified.



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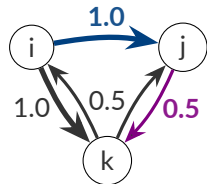


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The standard is to transform  $\tilde{\mathbf{W}}$  to be **row-stochastic**, such that  $\sum_j w_{ij} = 1 \ \forall i$ .



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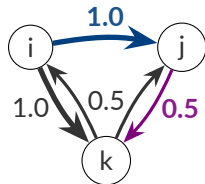
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## Scalar normalization

We will use **scalar normalization**, such that  $w_{ij} = g_{ij} \times \varsigma \ \forall i, j$ , in order to **preserve the network structure**. [► See more](#)



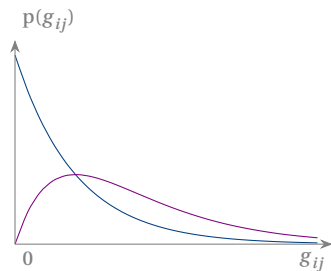
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We want to model links, and could do so directly

$$g_{ij} \sim f(\cdot) \quad \forall i \neq j.$$



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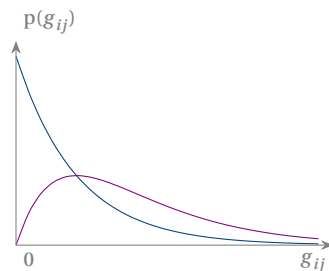
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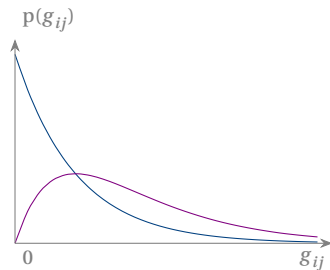
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We want to **constrain the dimensionality** by **imposing some structure**<sup>a</sup> on  $\mathcal{G}$ , allowing for more nuance where it is needed.

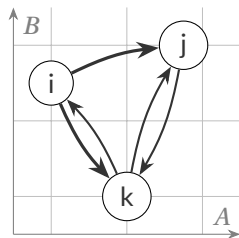
<sup>a</sup>Lewbel et al. (2023), e.g., constrain links to *sub-networks*. [► Illustration](#)

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# Imposing structure — A (Metric) Space Odyssey

Alternatively, assume that we can **locate our agents** in some (generalized) **metric space**  $(\mathcal{P}, d)$ .<sup>a</sup>



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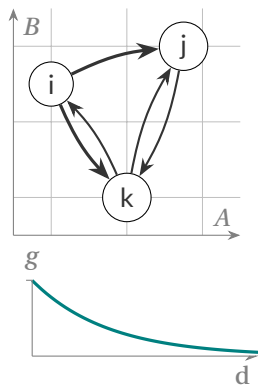
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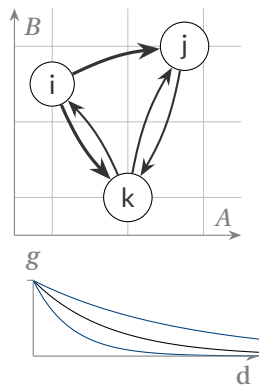
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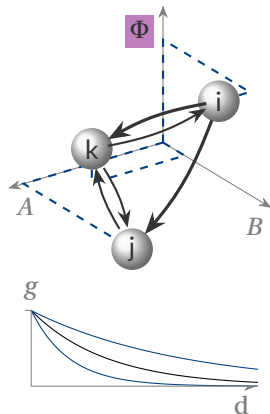
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- We may also consider, e.g., the **speed of decay**,
- or *asymmetries* via **popularity** or gravity.



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## Estimation or: How I Learned to Stop Worrying and Love MCMC

**Adaptive MCMC** facilitates full posterior inference, nuanced **weakly informative priors** improve convergence, and a **Gaussian process approximation** for costly Jacobians improves speed. [► See more](#)

# The approach in practice

We'll simulate repeatedly from

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## Setup

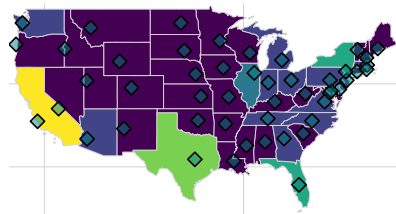
I'll model the *locations*, speed of *distance-decay*, and *popularity*. The **priors will be too flat**, and the *sampler's initialised* first at the true values, then at draws from the prior.

# Simulation results — US population

The first network is determined as

$$g_{ij} = \exp \left\{ -\delta_i \times d \left( \mathbf{p}_i, \mathbf{p}_j \right) \right\},$$

where **popularity**  $\delta_i$  is driven by a state's population, and  $\mathbf{p}_i$  its **population center**.



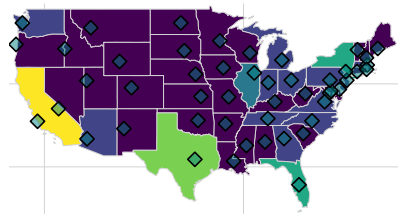


# Simulation results — US population

The first network is determined as

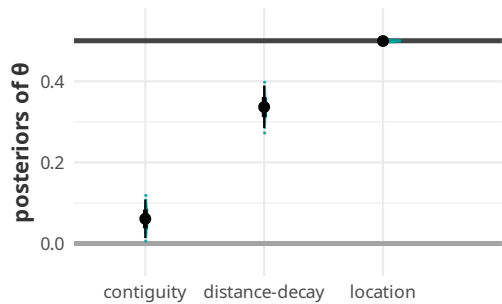
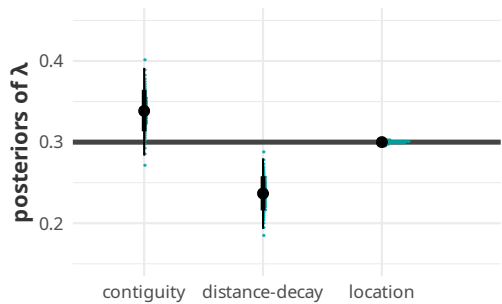
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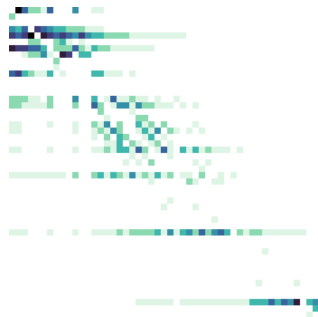
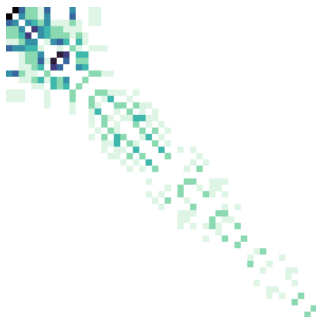
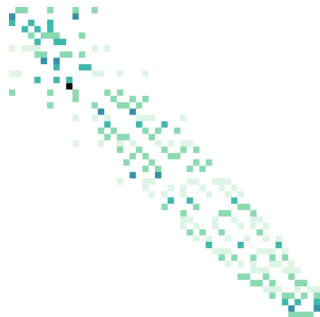
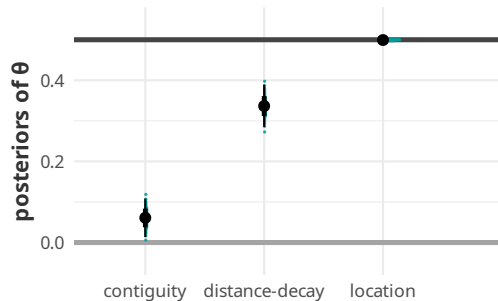
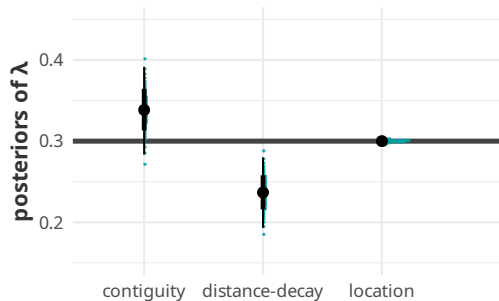
where **popularity**  $\delta_i$  is driven by a state's population, and  $\mathbf{p}_i$  its **population center**.



I'll start by estimating  $\lambda, \theta$  in a long panel ( $T = 50$ ) using

- **contiguity** between states,
- **distance-decay** between centers, and
- the *true model* of **location & popularity**.





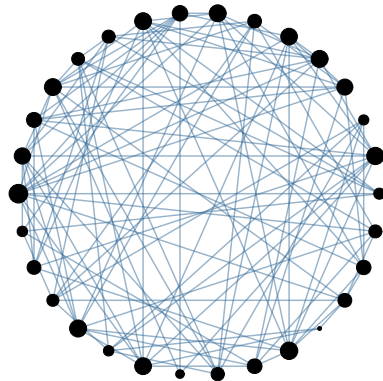
$W$  based on **contiguity**, **distance-decay**, and the **true network**.

# Simulation results — the Erdős–Rényi graph

The graph ( $N \in \{30, 50\}$ ) is determined as

$$g_{ij} = \begin{cases} 1 & \text{with probability 0.25,} \\ 0 & \text{otherwise.} \end{cases}$$

Our model uses the same **distance-decay** specification as before.



A realization of  $\mathcal{G}(30, 0.25)$ .

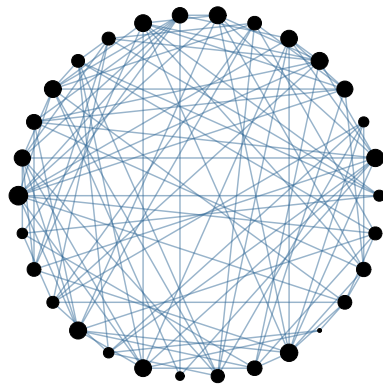
# Simulation results — the Erdős–Rényi graph

The graph ( $N \in \{30, 50\}$ ) is determined as

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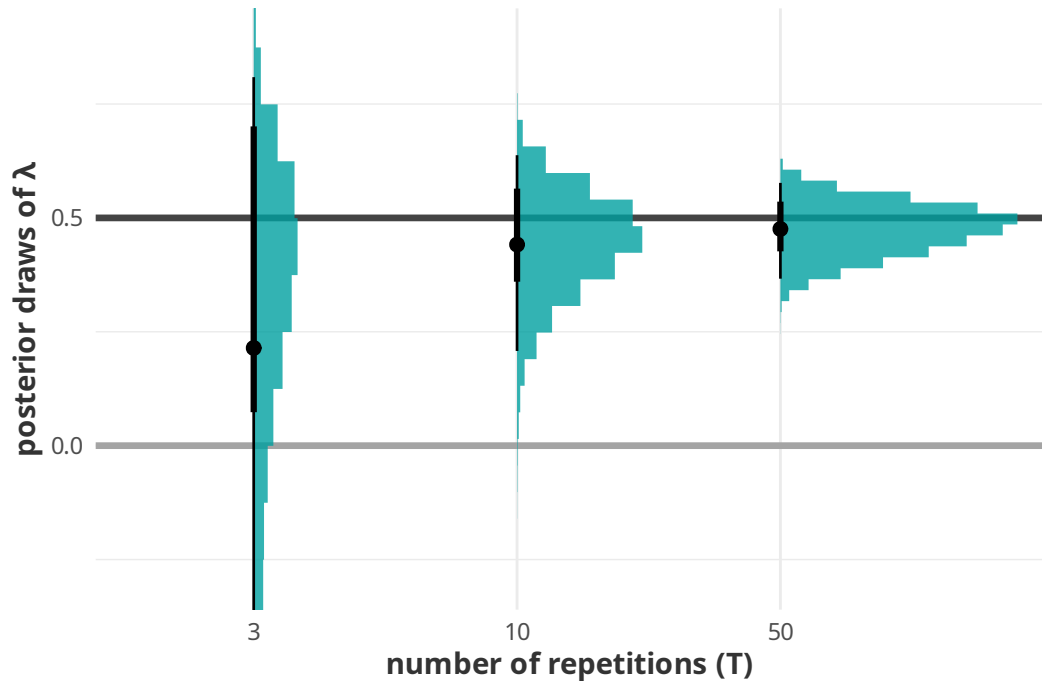
Our model uses the same **distance-decay** specification as before.

To **highlight convergence**, we'll have a look at **posteriors of  $\lambda$**  for multiple simulations after a **short burn-in** of 1,000 draws.

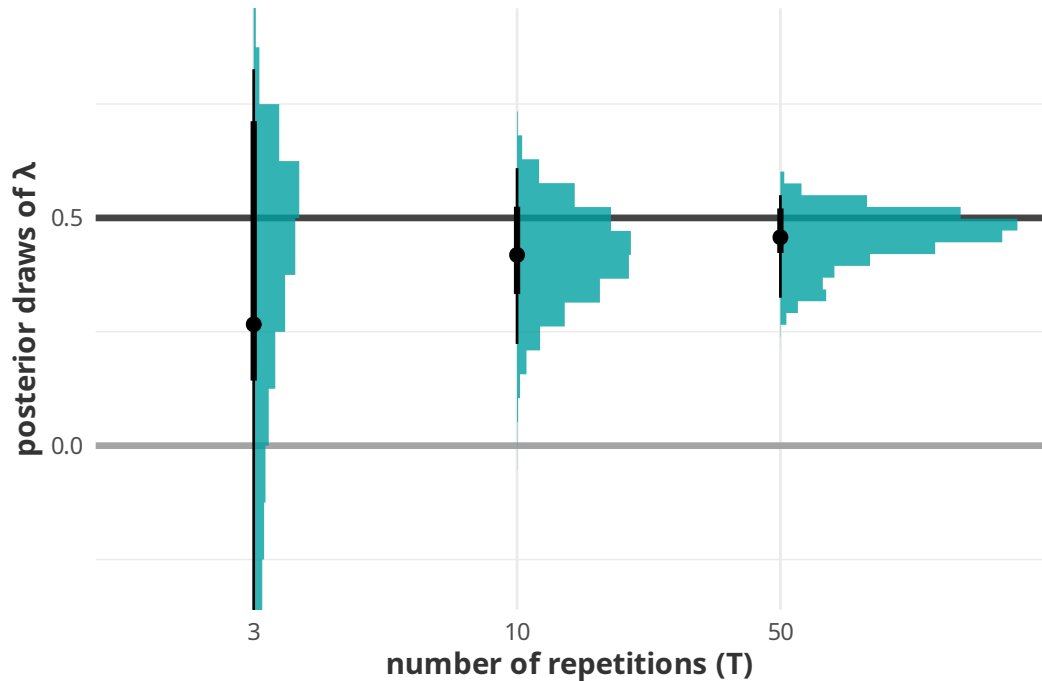


A realization of  $\mathcal{G}(30, 0.25)$ .

## Connectivity strength $\lambda$ (N = 30)



## Connectivity strength $\lambda$ (N = 50)



# Conclusion

- I developed a **framework for jointly modelling**  $f$  and  $\mathcal{G}$ ,
- that flexibly leverages **data, structure, and shrinkage**.
- It's widely applicable to **network and spatial settings**
  - *with no, limited, or uncertain information on diverse links,*
- allowing us to gain **deeper insights into spillovers**,
  - *at moderate to pronounced computational costs.*



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For **more details and info** on identification, priors, sampling, and applications have a look at *the appendix*, or — coming soon™ to a repository near you — a draft.



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# Model

We are interested in

$$p(\Theta, \mathcal{E} \mid \mathcal{D}) \propto p(\Theta, \mid \mathcal{E}, \mathcal{D}) \times p(\mathcal{E} \mid \Theta, \mathcal{D}),$$

or, to be more concrete, in

$$\mathbf{y} = (\mathbf{I} - \lambda \mathbf{W})^{-1} \mathbf{z},$$
$$\mathbf{z} = \mathbf{X}\boldsymbol{\beta} + \dot{\mathbf{W}}\mathbf{X}\boldsymbol{\theta} + \boldsymbol{\varepsilon},$$

where  $\mathbf{W} = g(\cdot)\zeta$ ,  $\dot{\mathbf{W}} = g(\cdot)\dot{\zeta}$ , and  $g$  is based on a network model of choice. Options include the ones described, many others, or a combination thereof.

## Identification

We can identify the parameters  $\lambda, \boldsymbol{\theta}$  with mild constraints. Network parameters are generally only weakly identified. We can alleviate this by imposing constraints from the literature or prior information.

# Normalization — multiplier effect

Consider a **network autoregression**

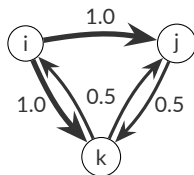
$$\mathbf{y} = \lambda \mathbf{W} \mathbf{y} + \mathbf{e}.$$

The following result guarantees stability.

## Theorem

Let  $\mathbf{I}$  denote the identity matrix, and  $\alpha$  be a real scalar. Then  $\mathbf{I} - \alpha \mathbf{A}$  is stationary for  $\alpha \in (-\rho_{\mathbf{A}}, \rho_{\mathbf{A}})$ , where  $\rho_{\mathbf{A}}$  denotes the spectral radius of  $\mathbf{A}$ .

By normalizing with the *spectral radius* (i.e.  $\varsigma = \rho_{\mathbf{G}}^{-1}$ ), we can let  $\lambda \in (-1, 1)$ . This relates  $\lambda$  to the *dominant eigenvector* of the network, which generally *does not* coincide with the *average partial effect* of  $\mathbf{W}$ .



Row-normalization distorts, e.g., the *eigenvector centrality*  $c$ .

	$i$	$j$	$k$
$c_{\mathbf{G}}$	$2^{-1}$	$6^{-1}$	$3^{-1}$
$c_{\bar{\mathbf{W}}}$	$3^{-1}$	$3^{-1}$	$3^{-1}$

In fact,  $\lambda$  is *at least* the average partial effect (by the spectral radius being the infimum norm).

# Normalization — contextual effect

Consider a contextual model

$$\mathbf{y} = \lambda \mathbf{W} \mathbf{X} \boldsymbol{\theta} + \mathbf{e}.$$

$$\frac{\partial \mathbf{z}}{\partial \mathbf{x}_k} = \mathbf{I} \beta_k + \mathbf{W} \theta_k.$$

In this case, we have fewer requirements of the normalization. One sensible option is to fix  $\boldsymbol{\theta}$  at the *average partial effect* of the network characteristics.<sup>a</sup>

We can achieve that by setting  $\mathbf{W}$  such that it sums to  $N$  — we scale with  $\zeta = \frac{N}{\sum_i \sum_j g_{ij}}$ .

This applies similarly to the nested linear model when a network multiplier is present. [► Go back](#)

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<sup>a</sup>An alternative when some agents are not linked in the network, is the average partial effect for all agents that are linked within the network.

# Estimation

For full posterior inference, we extend existing *MCMC methods* – the central term (suppressing contextual effects) is given by

$$|\mathbf{S}(\lambda, \cdot)| \exp \left\{ -\frac{1}{2\sigma^2} (\mathbf{S}(\lambda, \cdot)\mathbf{y} - \mathbf{X}\boldsymbol{\beta})' (\mathbf{S}(\lambda, \cdot)\mathbf{y} - \mathbf{X}\boldsymbol{\beta}) \right\},$$

where  $\mathbf{S}(\lambda, \cdot) = (\mathbf{I} - \lambda\mathbf{W})$  is a spatial filter. The main concerns are essentially computational – we need

1. convergence of parameters,
2. to evaluate the  $N \times N$  **Jacobian determinant**.

We use *Gibbs* and *Metropolis-Hastings* steps, as well as a *rejection sampler* to draw from the posterior.

# Weakly informative priors

We use weakly informative priors to (1) help improve convergence of the MCMC samples, and (2) build a more credible, realistic model.

Two central parameters are  $\lambda$  and  $\theta$ . For the former, I propose the *hierarchical prior*

$$\lambda \sim \text{Beta}(1 + \tau, 1 + \tau), \quad \tau \sim \text{Gamma},$$

which adds barely any computational overhead,<sup>a</sup> but facilitates much better shrinkage towards, e.g., zero, while providing wide support.

For  $\theta$ , which resembles a standard coefficient, standard global-local shrinkage priors are applicable.

Parameters for the network structure also greatly benefit from weakly informative priors, and even more from actually informative ones. Options for  $\theta$  include the Horseshoe, Dirichlet-Laplace, and the Normal-Gamma shrinkage priors.

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<sup>a</sup>Thanks to a *rejection sampler* based on a Gamma proposal density.

# A Beta prior ...

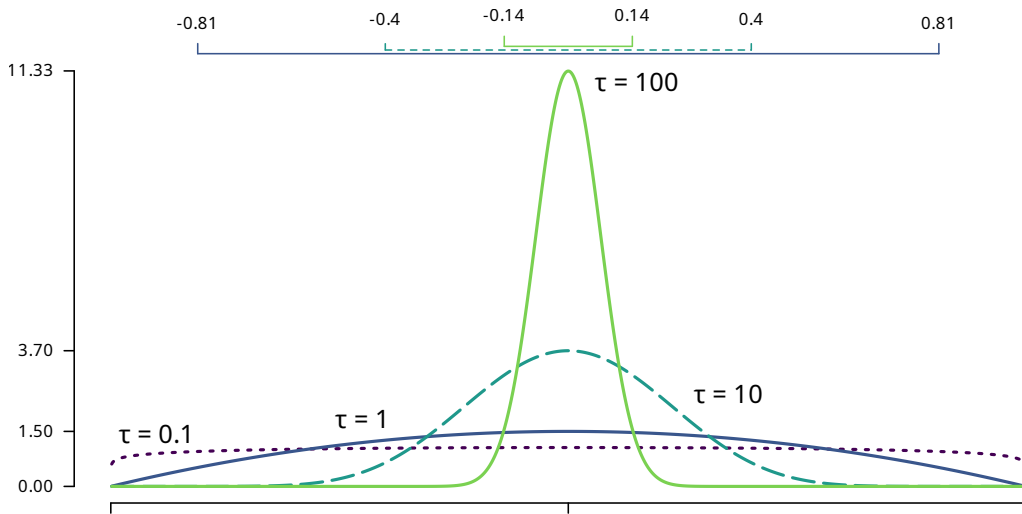


Figure 1: Scaled Beta(1 +  $\tau$ , 1 +  $\tau$ ) densities with increasing weight,  $\tau$ .

...with a Gamma mixing distribution

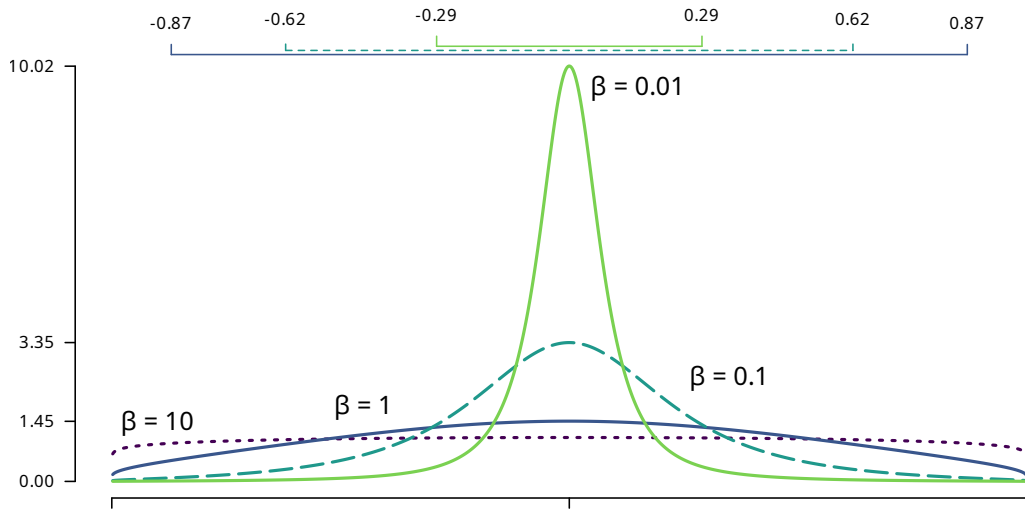


Figure 2: Scaled Beta(1 +  $\tau$ , 1 +  $\tau$ ),  $\tau \sim \text{Exp}(\beta)$  with increasing weight,  $\mathbb{E}[\tau] = 1/\beta$ . [Go back](#)



# Jacobian determinant

In standard models, we'd compute the **spectral decomposition** of **W** **once** and compute the determinant using **W**'s eigenvalues ( $\eta_i$ ), as

$$\ln |\mathbf{I} - \lambda \mathbf{W}| = \sum_{i=1}^N \ln(1 - \lambda \eta_i).$$

However, our **W** is **mutable**, and computing eigenvalues for every draw of  $\delta$  is prohibitive at  $\mathcal{O}(N^3)$  complexity.

For models with limited parameters for the network structure, I propose a **Gaussian process approximation** instead —

$$|\mathbf{S}(\lambda, \cdot)| \approx \text{GP}(\mu(\lambda, \cdot), \Sigma(\lambda, \cdot)).$$

# Gaussian process approximation

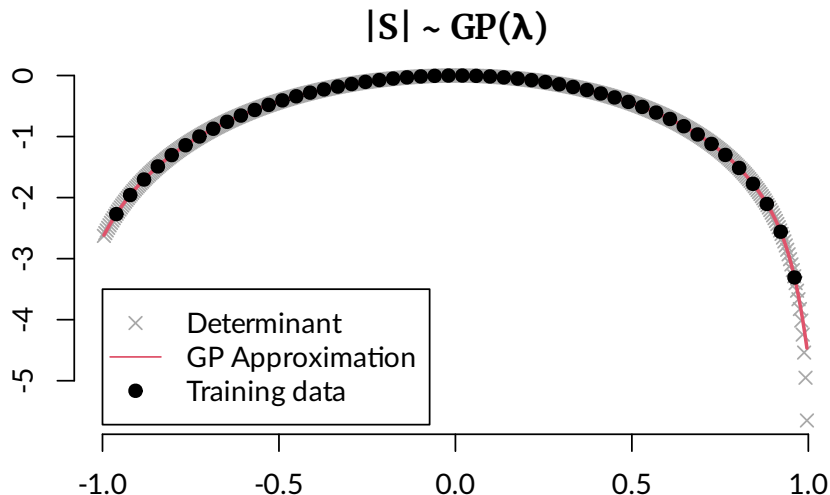


Figure 3: GP approximation to  $|S(\lambda)|$  using 50 training samples. Distances are between  $N = 100$  locations with Uniform random coordinates.

$|\mathbf{S}| \sim \text{GP}(\boldsymbol{\lambda}, \boldsymbol{\delta}), \text{ absolute error}$

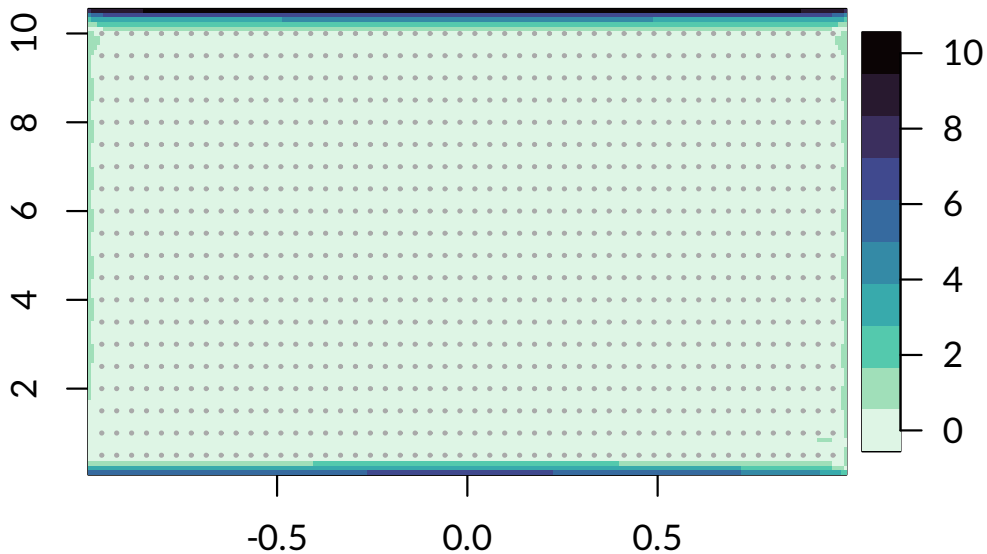
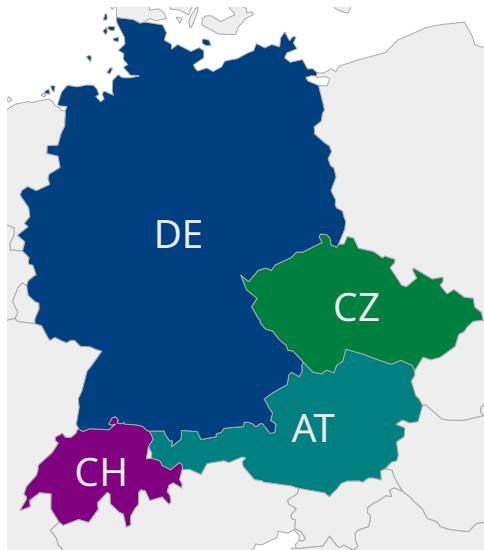


Figure 4: GP approximation to  $|\mathbf{S}(\boldsymbol{\lambda}, \boldsymbol{\delta})|$  using  $50 \times 20$  training samples.

# Imposing structure — Links Widely Shut

One way to reduce the dimensionality is by **constraining links to only occur within groups.**<sup>a</sup>

► Go back



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<sup>a</sup>This is essentially the approach of Lewbel et al., 2023.

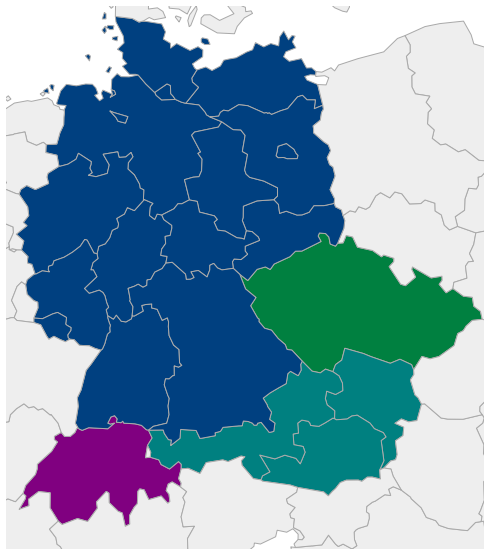
# Imposing structure — Links Widely Shut

One way to reduce the dimensionality is by **constraining links to only occur within groups.**<sup>a</sup> [▶ Go back](#)

Country	NUTS 1
Austria	3
Czechia	1
South Germany <sup>b</sup>	2 (16)
Switzerland	1
Total ( $N^2 - N$ )	42
Grouped ( $\sum_i N_i^2 - N_i$ )	10

<sup>a</sup>This is essentially the approach of Lewbel et al., [2023](#).

<sup>b</sup>Thanks for nothing, re-unification.



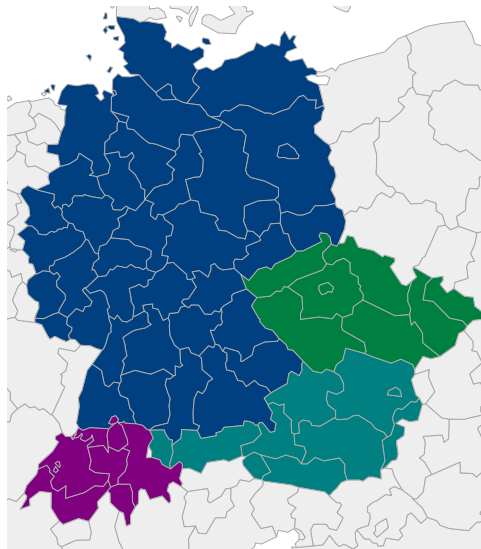
# Imposing structure — Links Widely Shut

One way to reduce the dimensionality is by **constraining links to only occur within groups.**<sup>a</sup> [▶ Go back](#)

Country	NUTS 1	NUTS 2
<b>Austria</b>	3	9
<b>Czechia</b>	1	8
South <b>Germany</b> <sup>b</sup>	2 (16)	11 (38)
<b>Switzerland</b>	1	7
Total ( $N^2 - N$ )	42	1190
Grouped ( $\sum_i N_i^2 - N_i$ )	10	280

<sup>a</sup>This is essentially the approach of Lewbel et al., [2023](#).

<sup>b</sup>Thanks for nothing, re-unification.



# Simulation results — US centroids

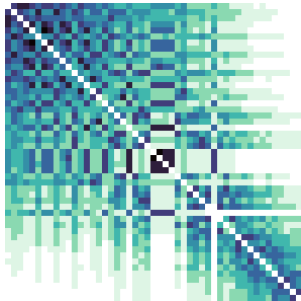
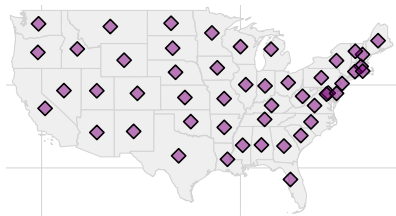
The second network is determined as

$$g_{ij} = \exp \left\{ -\delta \times d \left( \mathbf{p}_i, \mathbf{p}_j \right) \right\},$$

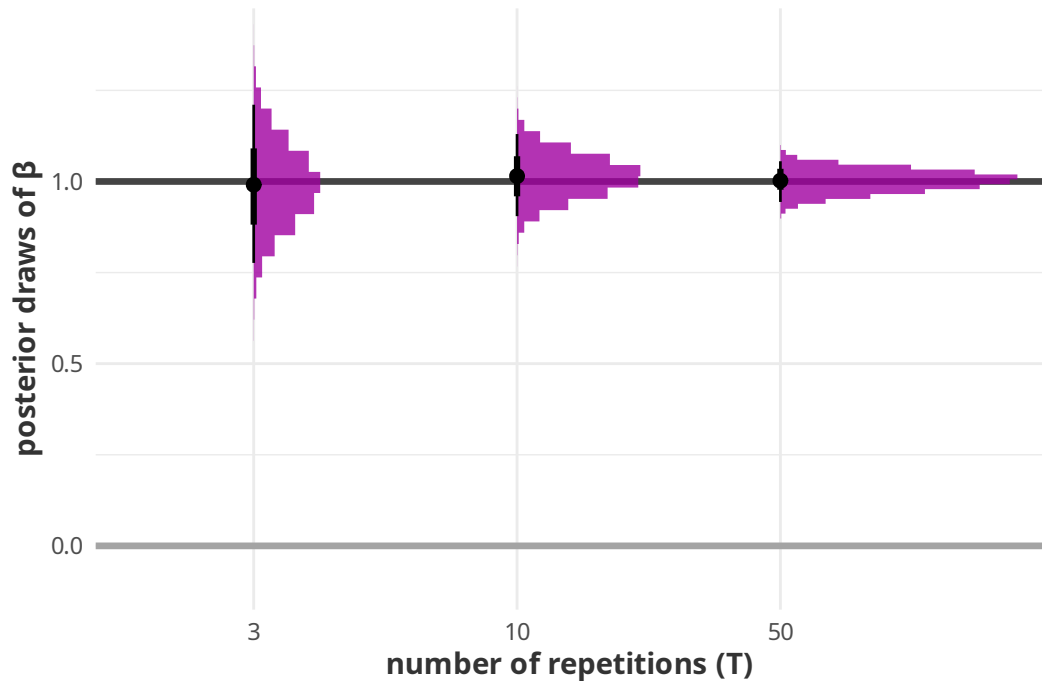
where  $\mathbf{p}_i$  are **centroids**, and

- the network is rather **dense**,
- I initialize the sampler at **random locations** to illustrate convergence.

We'll have a look at **posteriors of  $\lambda$**  for multiple simulations after a **short burn-in** of 1,000 draws. [▶ Go back](#)

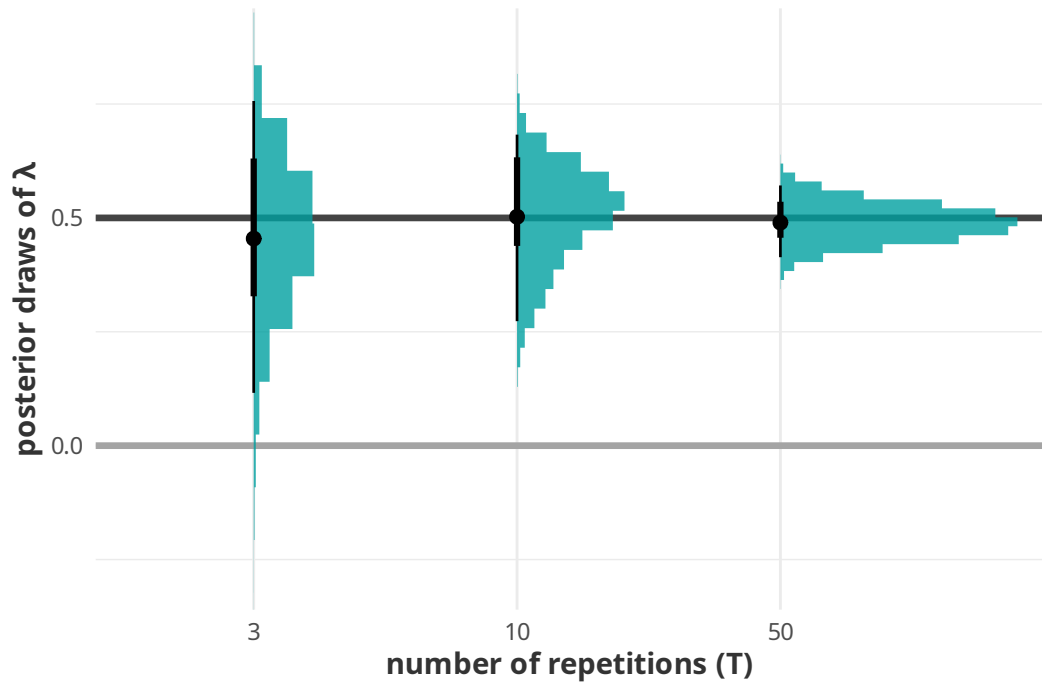


# Coefficient $\beta$ (N = 49)





# Connectivity strength $\lambda$ (N = 49)



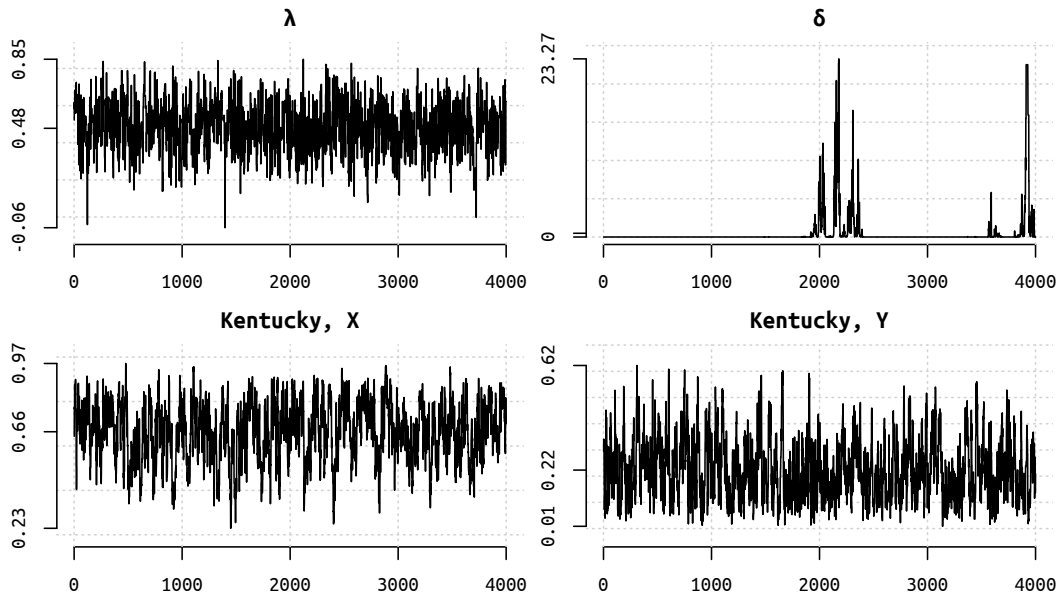


Figure 5: Traceplots for the posterior draws of  $\lambda$ ,  $\delta$  (note the poor mixing), and the (scaled) coordinates of Kentucky based on  $T = 3$  network observations.

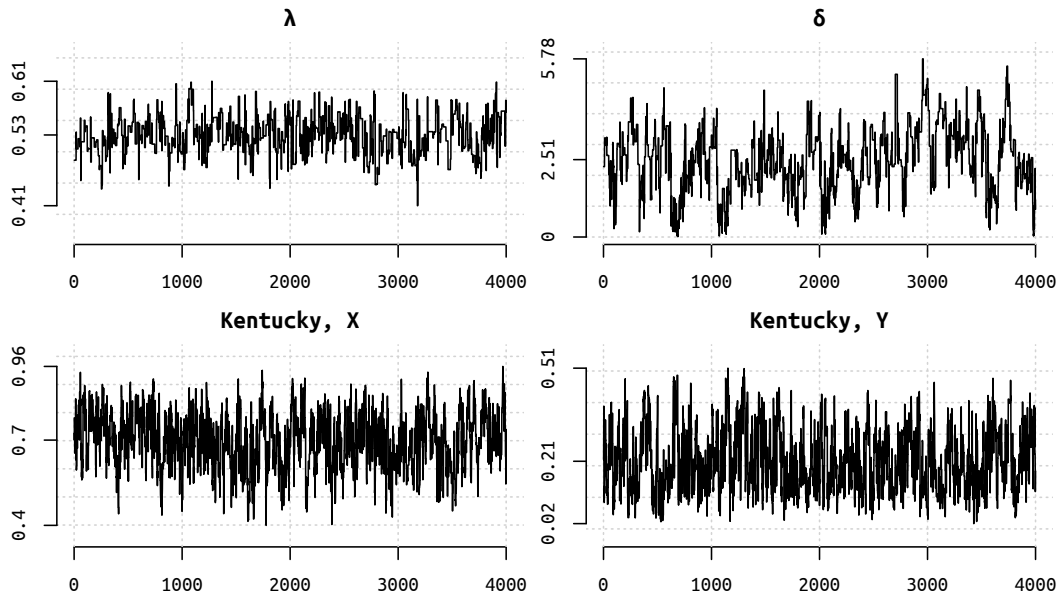


Figure 6: Traceplots for the posterior draws based on  $T = 50$  network observations. Note the improved mixing behavior of  $\delta$ . [► Go back](#)