# Almost-Matching-Exactly for Treatment Effect Estimation under Network Interference

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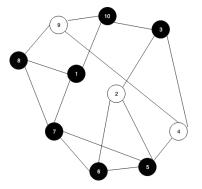
# Setting - Causal Inference

- ⊚ We have i = 1, ..., n experimental units
- ⊚ Treatment  $t_i \in \{0,1\}$  with  $\mathbf{t} \in \{0,1\}^n$  is a binary vector with the treatment level of every unit.
- © Potential outcomes  $Y_i(t_i, \mathbf{t})$  are random variables and depend on both treatment of unit i (1st argument), and treatment of **all other units** (2nd argument).
- ⊚ Observed treatment  $\mathbf{T} \in \{0,1\}^n$  is assigned **uniformly at** random.
- ⊚ Observed outcome:  $Y_i = T_i Y(1, \mathbf{T}) + (1 T_i) Y_i(0, \mathbf{T})$ 
  - Since treatment is randomized:  $\mathbb{E}[Y_i|\mathbf{T}=\mathbf{t},T_i=t]=\mathbb{E}[Y_i(t,\mathbf{T})]$  (Ignorability).
- ⊚ Units are connected in a network, G, in which unit i's treated neighborhood subgraph is  $G_{N_i^t}$ .

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#### The Problem: No SUTVA!

Usually we assume SUTVA: that *units' treatments don't influence other units' outcomes*, but we can't do that here because our units are connected in a network:



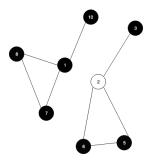
It could be that the treatment assigned to j influences the outcome of i through their connection in the network.

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# Similar Graphs Carry Similar Interference

#### Idea

What if the amount of interference experienced by a unit depended on the *shape* of its treated neighborhood subgraph?



Then, in expectation, two units with the same treated neighborhood graph will respond similarly to the treatment. We can use this idea to do matching to reduce interference.

## Assumptions

- 1. Outcome model:  $Y_i = \alpha + t_i \beta_i + f(G_{N_i^t}) + \epsilon_i$ 
  - f is some interference function dependent on  $G_{N_i^t}$ , the **treated neighborhood subgraph** of unit i.
- 2.  $\mathbb{E}[\epsilon_i] = 0$ 
  - Ignorability
- 3.  $|\mathbb{E}[f(g)] \mathbb{E}[f(h)]| \le K_1||g h||$ 
  - The more similar the neighborhood graphs of *i* and *j*, the more similar the amount of interference they receive.
  - Together with (1), this assumption encodes a version of SANASIA (Airoldi and Sussman, 2018) conditional on unit's neighborhood subgraphs.

#### Problem

How do we represent similarity between neighborhood graphs?

# Subgraph Counts

- We'll say that neighborhood graphs are similar if they contain similar counts of subgraphs
- What subgraphs? How similar must the counts be? How similar does this make the graphs?
- Use FLAME to decide

#### FLAME: An Overview

- FLAME (Fast Large-Scale Almost Matching Exactly) is a method for creating interpretable matches between units with discrete covariates
  - 1. Match units exactly on as many covariates as possible
  - 2. Drop a covariate
  - 3. Repeat
- At each step, drop the covariate maximizing match quality:

$$MQ = C \cdot BF - PE$$

- ⊚ BF = prop. controls matched + prop. treated matched
- Tradeoff between making matches and accurate prediction

#### ·AME-Networks

#### Problem #2

How do we choose which subgraphs we should use to represent the treated neighborhood graphs of our units?

#### Use FLAME to choose which subgraphs to count!

- 1. Enumerate (up to isomorphism) all p subgraphs  $S_1, \ldots, S_p$  seen across all the  $\mathcal{N}_i$ ,  $i = 1, \ldots, n$
- 2. For each unit i, define the p-dimensional vector  $S(G_{N_i^t})$  as the vector whose j'th entry is the number of  $S_j$  in  $\mathcal{N}_i$
- 3. These are likely a lot (maximum on the order of  $|\mathcal{N}_i|^2$ ) and it's unlikely that many units will have identical counts
- 4. Use FLAME to make almost-exact matches on subgraph counts

## A Small Change to FLAME

We know from our theoretical setup that network statistics should do two things well:

- 1. Predict the **outcomes**
- 2. Predict the **network**

To measure how well the our network statistics (subgraphs) are predicting the network, we model the edges of the network as independent conditional on the observed statistics

$$E_{ij}|x_i, x_j \stackrel{iid}{\sim} \text{Bern}(\text{logit}(\beta_1'x_i + \beta_2'x_j))$$

and consider the AIC of the resulting model, AIC<sub>network</sub>.

#### The Modified PE Function

To ensure FLAME strikes a balance between predicting both the **outcomes** and the **network**, we modify the PE function:

$$PE = \sum_{t=0}^{1} \underset{f \in \mathcal{F}}{\operatorname{arg \, min}} \frac{1}{n} \sum_{i=1}^{n} (Y_i - f(S(G_{\mathcal{N}_i^t})))^2$$
$$- \underbrace{D \cdot AIC_{\text{network}}}_{\text{new component}}$$

# Simulation Setup

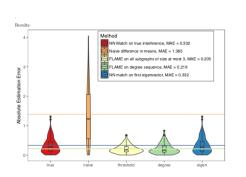
- © Recall the outcome model:  $Y_i = \alpha + \beta_i t_i + f(G_{N_i^t}) + \epsilon_i$
- The graph G is generated according to Erdos-Renyi or Stochastic Block models
- $\odot$  We consider various forms of f based off different features:
  - $\circ$   $d_i$ : the degree of unit i
  - $\Delta_i$ : the number of triangles in  $\mathcal{N}_i$
  - $\dagger_i^k$ : the number of units in  $\mathcal{N}_i$  with degree  $\geq k$
  - $\bigstar_i^k$ : the number of *k*-stars in  $\mathcal{N}_i$
  - Betweenness *i*: the vertex betweenness of unit *i*
  - Closeness<sub>i</sub>: the closeness centrality of unit i

#### **Estimators**

- ⊚ True: nearest neighbor (NN) on true interference
- Naive: naive difference in means
- © Eigen All: NN on eigenvalues of adjacency matrix *A*
- Eigen All: NN on largest eigenvalue of A
- Stratified Naive: Stratified degree estimator
- SANIA: MIVLUE under SANIA
- FLAME: Our approach

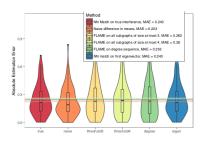
## (OLD) Simulation Results





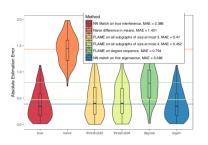
## (OLD) Simulation Results





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That the Interference that is not related to subgraph counts (denser graph) interference: f_r(\mathbf{z}) = \delta * Betweemass_r(\mathbf{z}) + \lambda * Closeness_r(\mathbf{z}) Outcomes: Y_i = a_i + \beta_i z_i + f_i(\mathbf{z}). Parameters: N = 100, N \sin = 100, \ a_i \sim N(0,1), \ \dot{\beta}_i \sim N(5,1), \ \dot{\delta} = 40, \ \lambda = 30, \ G \sim ER(0,1), \ Z_i \sim Ber(0,5) . Results:
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#### Bias Bound for oracle AME

As a preliminary result we can say that, under all the assumptions made before, with the true value of  $\phi$  and S known, and if we choose a match for unit i with treated neighborhood graph g such that:

$$j \in \mathrm{MG}(\mathbf{g}) \text{ if } j \in \underset{j=1,\ldots,n,T_j=0}{\arg\min} | \boldsymbol{\phi}^T S(g) - \boldsymbol{\phi}^T S(G_{\mathcal{N}_j^{\mathbf{t}}})|,$$

then the bias for the CATT of i can be upper bounded by:

$$|\mathbb{E}[Y_i - Y_j] - \tau_i| \leq K_1 \sum_{h \in \mathcal{G}} |\phi^T S(g) - \phi^T S(h)| \frac{exp(\phi^T S(h))}{\sum_{\ell \in \mathcal{G}} exp(\phi^T S(\ell))} \times \left[ \sum_{d=S(g)-|S(g)-S(h)|}^{S(g)+|S(g)-S(h)|} \frac{D_{\phi,S}(d) exp(d)}{\sum_{\ell \in \mathcal{G}} exp(\phi^T S(\ell))} \right]^{n-1}$$

### Plan: We Want to Hit the October 8 AISTATS Deadline

- 1. We need to implement the revised PE function
- 2. We need to redo the simulations
- 3. More theory? Statements on how well the subgraph count can "encode" a given graph would be nice
- 4. Find a good application
- 5. Write!