Randomized Graph Cluster Randomization With R Package: RGCR

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- ▶ Ugander, J. and Yin, H., 2023. Randomized graph cluster randomization. *Journal of Causal Inference*, 11(1), p.20220014.
- ► You can download 'RGCR' from https://github.com/RUC-ChangHao/RGCR.

Classic causal inference and randomization

- ▶ Binary treatment $Z_i \in \{0,1\}$ ($Z \sim P(z)$, CR, Bernoulli design, etc)
- ▶ Potential outcomes $Y_i(1)$ and $Y_i(0)$ (SUTVA assumption)
- Causal effects of interest: average causal effect (ACE)

$$ACE \stackrel{\text{def}}{=} \sum_{i=1}^{n} \{ Y_i(1) - Y_i(0) \}$$

Estimators: ht (ipw), hajek, aipw, etc

$$\hat{\tau}_{ht} = \frac{1}{n} \sum_{i=1}^{n} \frac{Y_i(1)Z_i}{P(Z_i=1)} - \frac{1}{n} \sum_{i=1}^{n} \frac{Y_i(0)(1-Z_i)}{P(Z_i=0)}$$

Technically, a missing data problem

	1	2	3		n
$Y_i(1)$	36.5	?	?		38.0
$Y_i(0)$?	38.0	37.0	• • •	?

▶ Question: To get better estimation of ATE, how do we assign treatment Z? (i.e., seek the 'optimal' experiment design $Z \sim P(z)$)

Causal inference under interference

- Violation of SUTVA
- Common in advertising, epidemiology and educational studies
- ▶ Potential outcomes $Y_i(\mathbf{z})$, where $\mathbf{z} \in \{0,1\}^n$
- ► Causal effects of interest: total treatment effect (TTE)

$$\text{TTE} \stackrel{\text{def}}{=} \frac{1}{n} \sum_{i=1}^{n} \{ Y_i(\mathbf{1}) - Y_i(\mathbf{0}) \}$$

Classic randomization does not work here (why?)

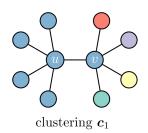
Table: Bias (SD) of HT Estimators

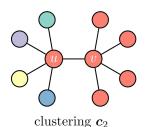
sample size	16 ²	24 ²	32 ²	48 ²	64 ²
SUTVA	0.00 (0.22)	-0.04 (0.16)	0.02 (0.13)	0.00 (0.08)	0.01 (0.05)
interference	0.09 (7.91)	-1.00 (3.74)	-0.47 (8.47)	-0.39 (2.33)	0.22 (4.61)

General framework for interference

- ightharpoonup A social network G = (V, E)
 - through which individuals interfere each other
 - observable and correctly measured
- An exposure mapping
 - determines the extent and intensity of the interference
 - technically reduces the number of potential outcomes
 - canonical examples (minor notation abuse)
 - ightharpoonup (no interference) $Y_i(z) = Y_i(z_i)$
 - (neighborhood interference) $Y_i(z) = Y_i(z_{N_i})$
 - ightharpoonup (arbitrary interference) $Y_i(z) = Y_i(z)$
 - ("individualized" interference) $Y_i(z) = Y_i(?)$
- Estimators: ht, hajek, difference-in-means, etc
- Experimental designs $Z \sim P(z)$: complete randomization, bernoulli randomization, cluster randomization, etc

A toy example





Under full neighborhood interference

$$\text{TTE} \stackrel{\mathsf{def}}{=} \tfrac{1}{n} \textstyle \sum_{i=1}^n \{ Y_i(\mathbf{1}) - Y_i(\mathbf{0}) \} = \tfrac{1}{n} \textstyle \sum_{i=1}^n \{ Y_i(\mathbf{1}_{\mathcal{N}_i}) - Y_i(\mathbf{0}_{\mathcal{N}_i}) \}$$

- what really matters is valid sample size
- what's a good design? intuitively (under proper social networks)
 - gives valid sample size a lower bound (increasing with sample size)
 - gives every node a positive exposure probability (due to ht estimators)

Randomized graph cluster randomization

- Randomized graph clustering
 - a weight indicating which nodes are more important
 - related to the variance of $\hat{\tau}_{ht}$, roughly speaking, related to the inverse exposure probabilities $1/\pi_i^1$ and $1/\pi_i^0$
 - types: uniform, degree, spectral weighting
 - ► a randomized clustering design deduced by the weight
 - types: 3-net and 1-hop-max (heuristically)
 - a clustering results generated from the clustering design
- Randomization at the level of clusters
 - types: complete design and bernoulli design
- Remark: no literature really solves this problem, all algorithms are heuristic in some sense

$$d^*, \hat{\tau}^* = \arg\min_{d \in \mathcal{D}, \hat{\tau} \in \mathcal{M}} \max_{\{Y_i(z)\} \subset \mathcal{Y}} E_{\{Y_i(z)\}, d, \hat{\tau}} \left[\operatorname{dist} \left(\hat{\tau}_d(Y, Z), \tau(\{Y_i(z)\}) \right) \right]$$

Weighted 3-net clustering

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Algorithm 4: Weighted 3-net clustering.
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Input: Graph G = (V, E), node weights \boldsymbol{w} \in \mathbb{R}^n_+.
    Output: Graph clustering c \in \mathbb{R}^n
 1 for i \in V do
 \mathbf{z} \mid X_i \leftarrow \beta(w_i, 1)
 \mathbf{3} \ \pi \leftarrow \operatorname{arg\,sort}([X_i]_{i \in V}, descend)
 4 S \leftarrow \emptyset, unmark all nodes
 5 for i \in \pi do
         if i is unmarked then
             S \leftarrow S \cup \{i\}
             for i \in B_2(i) do
               mark node j if it is unmarked yet
 9
10 for i \in V do
        c_i \leftarrow \arg\min\{j \in S, j \to \mathbf{dist}(i, j)\}\, i.e., the id of the node in S
          with shortest graph distance to i (arbitrary tie breaking)
12 return c
```

Weighted 1-hop-max clustering

Algorithm 3: Weighted 1-hop-max clustering.

Input: Graph G = (V, E), node weights $\boldsymbol{w} \in \mathbb{R}^n_+$.

Output: Graph clustering $c \in \mathbb{R}^n$

- 1 for $i \in V$ do
- $\mathbf{z} \mid X_i \leftarrow \beta(w_i, 1)$
- $\mathbf{3}$ for $i \in V$ do
- 4 $\lfloor c_i \leftarrow \max([X_j \text{ for } j \in B_1(i)])$
- $_{5}$ return $_{c}$

R package 'RGCR'

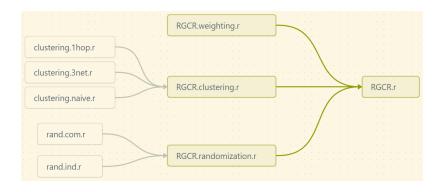
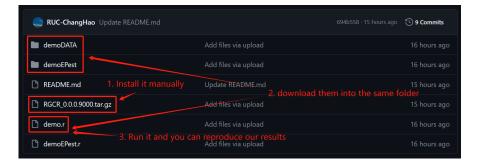


Figure: The Structure of functions

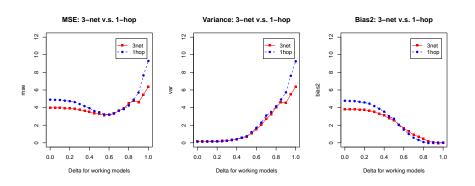
Source: https://github.com/RUC-ChangHao/RGCR

A short guidance to use 'RGCR'



Application of 'RGCR'

► The real data analysis of experimental design is actually a simulation



► Conclusion: the design 'spectral weighting + 3-net clustering + complete randomization' is recommended (after numerous simulations...)

Thank you!