Social Network Analysis Assignment Time-Aware Network Centrality Measures & Link Prediction

Dr. Dionisios N. Sotiropoulos

June 8, 2022

This assignment focuses on the algorithmic manipulation of the **Stack Over**flow Temporal Network, whose description and related dataset may be found at ^{[1](#page-0-0)}. Each edge of the underlying graph is associated with a timestamp, indicating the exact time instance where the edge was created. The complete set of directional edges for the aforementioned network, along with the associated timestamps, is row-wise stored in the file sx-stackoverflow.txt as consecutive triplets of the form (source_id,target_id,timestamp). Thus, the dataset is actually a collection of timestamped edges of the following form

$$
E = \{e_{ij}(t) = (v_i, v_j, t) : t_{min} \le t \le t_{max}\}\
$$
 (1)

where t_{min} and t_{max} identify the oldest and latest time instances respectively.

Your analysis should be conducted on a sequence of N non-overlapping time periods $\{T_1, \ldots T_N\}$, of equal duration δt , that span the entire time interval

$$
T = [t_{min}, t_{max}] \tag{2}
$$

The set of N time periods may be defined by considering a sequence of $N + 1$ time-instances $\{t_0, \ldots, t_N\}$ such that:

$$
t_j = t_{min} + j * \delta t, \ 0 \le j \le N \tag{3}
$$

where $\delta t = \frac{\Delta T}{N}$ and $\Delta T = t_{max} - t_{min}$. It is easy to deduce that the j-th time period may be defined as:

$$
T_j = \begin{cases} \begin{bmatrix} t_{j-1}, t_j \end{bmatrix}, & 1 \le j \le N-1; \\ \begin{bmatrix} t_{j-1}, t_j \end{bmatrix}, & j = N. \end{cases}
$$
\n
$$
(4)
$$

For each time period T_j , with $1 \leq j \leq N$, we may consider the corresponding undirected subgraphs of the network, denoted as:

$$
G[t_{j-1}, t_j] = (V[t_{j-1}, t_j], E[t_{j-1}, t_j])
$$
\n(5)

where

$$
E[t_{j-1}, t_j] = \{e_{ij}(t) : t \in T_j\}
$$
\n(6)

The set $V[t_{j-1}, t_j]$ of vertices for each period may be implicitly defined as the set of nodes that appear at the end points of edges pertaining to the set $E[t_{i-1}, t_i]$.

Part I Questions (Grade Percentage 35%):

¹https://snap.stanford.edu/data/sx-stackoverflow.html

- 1. Partition the complete time period $T = [t_{min}, t_{max}]$ into a set of nonoverlapping time periods $\{T_1, \ldots, T_N\}$ by computing the corresponding set of time instances $\{t_0, \ldots, t_N\}$ where $t_0 = t_{min}$ and $t_N = t_{max}$. Mind that N is a user defined parameter.
- 2. Choose an appropriate representation for each subgraph $G[t_{j-1}, t_j]$ of the network for each time period T_j where $1 \leq j \leq N$.
- 3. Provide a graph depicting the time evolution of the quantities $|V[t_{i-1}, t_i]|$ and $|E[t_{j-1}, t_j]|$ for each time period T_j where $1 \leq j \leq N$.
- 4. For each subgraph $G[t_{j-1}, t_j]$ compute and graphically represent the probability density functions (i.e. histograms of relative frequencies) for the following centrality measures:
	- (a) Degree Centrality
	- (b) Closeness Centrality
	- (c) Betweenness Centrality
	- (d) Eigenvector Centrality
	- (e) Katz Centrality

Acquiring a more accurate description for the evolution of the network between successive time periods can be facilitated by considering the set of nodes that persist during the transition from T_j to T_{j+1} , formulated as:

$$
V^*[t_{j-1}, t_{j+1}] = V[t_{j-1}, t_j] \cap V[t_j, t_{j+1}], \ 1 \le j \le N \tag{7}
$$

In this setting, we are particularly interested in restricting the sets $E[t_{i-1}, t_i]$ and $E[t_j, t_{j+1}]$ within the common set of nodes $V^*[t_{j-1}, t_{j+1}]$ as:

$$
E^*[t_{j-1}, t_j] = \{(u, v) \in E[t_{j-1}, t_j] : u \in V^*[t_{j-1}, t_{j+1}] \land v \in V^*[t_{j-1}, t_{j+1}] \} (8)
$$

$$
E^*[t_j, t_{j+1}] = \{(u, v) \in E[t_j, t_{j+1}] : u \in V^*[t_{j-1}, t_{j+1}] \land v \in V^*[t_{j-1}, t_{j+1}] \} \tag{9}
$$

Part II Questions (Grade Percentage 35%):

- 1. For each pair of successive network instances $(G[t_{j-1}, t_j], G[t_j, t_{j+1}])$, where $1 \leq j \leq N-1$, compute the following sets
	- (a) $V^*[t_{j-1}, t_{j+1}]$
	- (b) $E^*[t_{j-1}, t_j]$
	- (c) $E^*[t_j, t_{j+1}]$

and graphically represent their volumes $|V^*[t_{j-1}, t_{j+1}]|, |E^*[t_{j-1}, t_j]|$ and $|E^*[t_j, t_{j+1}]|$ as functions of the coupled time periods (T_j, T_{j+1}) .

2. For each pair of nodes $(u, v) \in V^*[t_{j-1}, t_{j+1}]$ and for every set of common vertices $V^*[t_{j-1}, t_{j+1}]$, where $1 \leq j \leq N-1$, compute the following similarity matrices:

(a)
$$
S_{GD}: S_{GD}(u, v) = -d_{geodesic}(u, v)
$$
 [Graph Distance]

(b) $\mathbf{S_{CN}}$: $S_{CN}(u, v) = |\Gamma(u) \cap \Gamma(v)|$ [Common Neighbors]^{[2](#page-2-0)}

\n- (c)
$$
\mathbf{S_{JC}} : S_{JC}(u, v) = \frac{|\Gamma(u) \cap \Gamma(v)|}{|\Gamma(u) \cup \Gamma(v)|} \left[\text{Jaccard's Coefficient} \right]
$$
\n- (d) $\mathbf{S_A} : S_A(u, v) = \sum_{z \in \Gamma(u) \cap \Gamma(v)} \frac{1}{\log(|\Gamma(z)|)} \left[\text{Adamic } / \text{ Adar} \right]$
\n- (e) $\mathbf{S_{PA}} : S_{PA}(u, v) = |\Gamma(u)| * |\Gamma(c)| \left[\text{Preferential Attachment} \right]$
\n

According to the previous discussion, it is easy to deduce that during the successive time periods T_j and T_{j+1} , the set of all possible edges between vertices in the common set of nodes $V^*[t_{j-1}, t_{j+1}]$ may be given as:

$$
E^{0}[t_{j-1}, t_{j+1}] = V^{*}[t_{j-1}, t_{j+1}] \times V^{*}[t_{j-1}, t_{j+1}], \ 1 \leq j \leq N - 1 \qquad (10)
$$

However, the subset of edges that are actually realized corresponds to the set

$$
E^*[t_{j-1}, t_{j+1}] = E^*[t_{j-1}, t_j] \cup E^*[t_j, t_{j+1}] \tag{11}
$$

In the context of the link prediction task, $E^*[t_{j-1}, t_j]$ will serve as the training set, whereas $E^*[t_j, t_{j+1}]$ will be used for testing. Each one of the previously defined similarity metrics

$$
\mathbf{S}_{\mathbf{X}}: X \in \{GD, CN, JC, A, PA\} \tag{12}
$$

can be employed in order to implement a simple classification mechanism that provides an estimation for the actual set of edges $E^*[t_{j-1}, t_{j+1}]$ according to the following equation:

$$
\hat{E}_X^*[t_{j-1}, t_{j+1}] = \{(u, v) \in E^0[t_{j-1}, t_{j+1}] : S_X(u, v) \in R_X\}
$$
\n(13)

where R_X indicates a range of values for the similarity score S_X . The prediction accuracy of each classification rule defined by Eq. [13](#page-2-1) can be assessed with respect to a ground truth set of edges E through the utilization of the quantity given below:

$$
ACC(R_X, E) = \lambda * TPR(R_X, E) + (1 - \lambda) * TNR(R_X, E)
$$
(14)

where

$$
TPR(R_X, E) = \frac{|\hat{E}_X^*[t_{j-1}, t_{j+1}] \cap E|}{|E|} \tag{15}
$$

$$
TNR(R_X, E) = 1 - \frac{|\hat{E}_X^*[t_{j-1}, t_{j+1}]| - |\hat{E}_X^*[t_{j-1}, t_{j+1}] \cap E|}{|E^0[t_{j-1}, t_{j+1}]| - |E|}
$$
(16)

$$
\lambda = \frac{|E|}{|E^0[t_{j-1}, t_{j+1}]|} \tag{17}
$$

The simplest way to define R_X is as a continuous interval of the following form:

$$
R_X = [S_X^L, S_X^U] \tag{18}
$$

where S_X^L and S_X^U are the lower and upper bounds respectively. However, more accurate classification results can be obtained by considering a more composite

²For a graph $G = (V, E)$, function $\Gamma : V \to P(V)$, evaluated on a particular node $z \in V$ provides the subset $\Gamma(z) \subset V$ of nodes neighboring with z

form for the set R_X , composed by the union of n_X non-overlapping intervals, formulated as:

$$
R_X = \bigcup_{k=1}^{k=n_X} [S_X^{L_k}, S_X^{U_k}]
$$
\n(19)

Part III Questions (Grade Percentage 30%):

1. Describe and implement a training algorithm which determines the optimal range sets R_X^* , defined by Eq. [19,](#page-3-0) for each similarity measure. The goal of the training algorithm should be the maximization of accuracy given by Eq. [14](#page-2-2) within the training set. Therefore, the training algorithm reduces to solving the following maximization problem:

$$
R_X^* = \arg\max_{R_X} ACC(R_X, E^*[t_{j-1}, t_j])
$$
\n(20)

- 2. Having determined the optimal range sets R_X^* for each similarity measure, evaluate and rank the corresponding training accuracy measurements $ACC(R_X^*, E^*[t_{j-1}, t_j]).$
- 3. Evaluate and rank the testing accuracy measurements $ACC(R_X^*, E^*[t_j, t_{j+1}])$.

For this assignment you can work in groups of no more than 3 students. Your implementation can be in any programming language. The final deliverable should contain:

- 1. Well documented code of your implementation.
- 2. A concise report explaining your assumptions and implementation decisions.
- 3. Example runs of your code providing the required graphical representations and classification measurements.