DyREP: Learning Representations over Dynamic Graphs

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Paper under double-blind review

Abstract

Representation Learning over graph structured data has received significant attention recently due to its ubiquitous applicability. However, most advancements have been made in static graph settings while efforts for jointly learning dynamic of the graph and dynamic on the graph are still in an infant stage. Two fundamental questions arise in learning over dynamic graphs: (i) How to elegantly model dynamical processes over graphs? (ii) How to leverage such a model to effectively encode evolving graph information into low-dimensional representations? We present **DyRep** - a novel modeling framework for dynamic graphs that posits representation learning as a *latent mediation process* bridging two observed processes namely – dynamics *of* the network (realized as topological evolution) and dynamics *on* the network (realized as activities between nodes). Concretely, we propose a two-time scale deep temporal point process model that captures the interleaved dynamics of the observed processes. This model is further parameterized by a temporal-attentive representation network that encodes temporally evolving structural information into node representations which in turn drives the nonlinear evolution of the observed graph dynamics. Our unified framework has inductive capability to generalize over unseen nodes and we design an efficient unsupervised procedure for end-to-end training. We demonstrate that DyRep outperforms state-of-art baselines in quantitative analysis using dynamic link prediction and time prediction tasks. We further present extensive qualitative insights into our framework to discern indispensable role of various components of our framework.

1 Introduction

Representation learning over graph structured data has emerged as a keystone machine learning task due to its ubiquitous applicability in variety of domains such as social networks, bioinformatics, natural language processing, and relational knowledge bases. Learning node representations to effectively encode high-dimensional and non-Euclidean graph information is a challenging problem but recent advances in deep learning has helped important progress towards addressing it (Cao et al., 2015; Grover & Leskovec, 2016; Perozzi et al., 2014; Tang et al., 2015; Wang et al., 2016a; 2017; Xu et al., 2017), with majority of the approaches focusing on advancing the state-of-art in static graph setting. However, several domains now present highly dynamic data that exhibit complex temporal properties in addition to earlier cited challenges. For instance, social network communications, financial transaction graphs or longitudinal citation data contain fine-grained temporal information on nodes and edges that characterize the dynamic evolution of a graph and its properties over time.

These recent developments have created a conspicuous need for principled approaches to advance graph embedding techniques for dynamic graphs (Hamilton et al., 2017b). We focus on two pertinent questions fundamental to representation learning over dynamic graphs: (i) *What can serve as an elegant model for dynamic processes over graphs?* — A key modeling choice in existing representation learning techniques for dynamic graphs (Goyal et al., 2017; Zhou et al., 2018; Trivedi et al., 2017; Ngyuyen et al., 2018; Yu et al., 2018) assume that graph dynamics evolve as a single time scale process. In contrast to these approaches, we observe that most real-world graphs exhibit at least two distinct dynamic processes that evolve at different time scales — *Topological Evolution*: where the number of nodes and edges are expected to grow (or shrink) over time leading to structural changes in the graph; and *Node Interactions*: which relates to activities between nodes that may or may not be structurally connected. Modeling interleaved dependencies between these non-linearly evolving dynamic processes is a crucial next step for advancing the formal models of dynamic graphs.
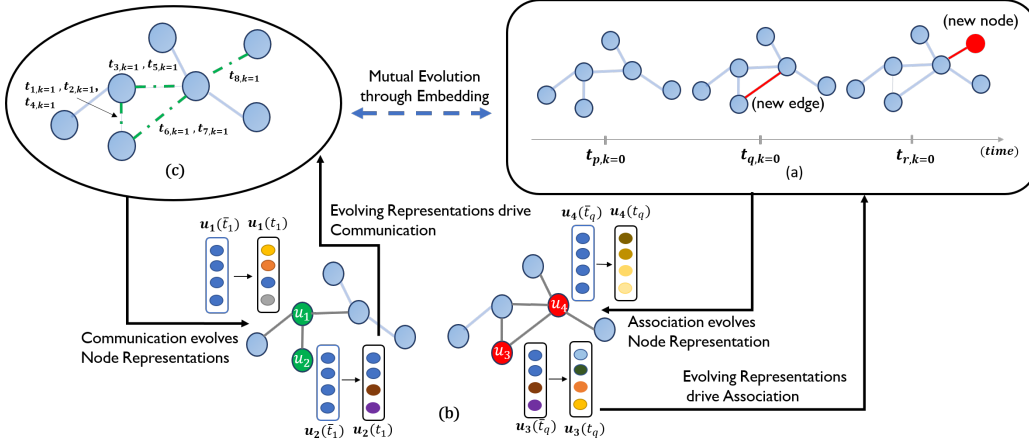
Figure 1: Evolution Through Mediation. **(a)** Association events (k=0) where the node or edge grows. **(c)** Communication Events (k=1) where nodes interact with each other. For both these processes, $t_{p,k=0} < (t_1, t_2, t_3, t_4, t_5)_{k=1} < t_{q,k=0} < (t_6, t_7)_{k=1} < t_{r,k=0}$. **(b)** Evolving Representations.

(ii) *How can one leverage such a model to learn dynamic node representations that are effectively able to capture evolving graph information over time?* — Existing techniques in this direction can be divided into two approaches: a.) Discrete-Time Approach, where the evolution of a dynamic graph is observed as collection of static graph snapshots over time (Zhu et al., 2016; Goyal et al., 2017; Zhou et al., 2018). These approaches tend to preserve (encode) very limited structural information and capture temporal information at a very coarse level which leads to loss of information between snapshots and lack of ability to capture fine-grained temporal dynamics. Another challenge in such approaches is the selection of appropriate aggregation granularity which is often misspecified. b.) Continuous-Time Approach, where evolution is modeled at finer time granularity in order to address the above challenges. While existing approaches have demonstrated to be very effective in specific settings, they either model simple structural and complex temporal properties in a decoupled fashion (Trivedi et al., 2017) or use simple temporal models (exponential family in (Ngyuyen et al., 2018)). But several domains exhibit highly nonlinear evolution of structural properties coupled with complex temporal dynamics and it remains an open problem to effectively model and learn informative representations capturing various dynamical properties of such complex systems.

As noted in (Chazelle, 2012), an important requirement to effectively learn over such dynamical systems is the ability to express the dynamical processes at different scales. We propose that any dynamic graph must be minimally expressed as a result of two fundamental processes evolving at different time scales: **Association Process** (dynamics *of* the network), that brings change in the graph structure and leads to long lasting information exchange between nodes; and **Communication Process** (dynamics *on* the network), that relates to activities between (not necessarily connected) nodes which leads to temporary information flow between them (Farine, 2017; Artime et al., 2017). We, then, posit our goal of learning node representations as modeling a *latent mediation process* that bridges the above two observed processes such that learned embeddings drive the complex temporal dynamics of both processes and these processes subsequently lead to the nonlinear evolution of node representations. Further, the information propagated across the graph is governed by the temporal dynamics of communication and association histories of nodes with its neighborhood. For instance, in a social network, when a node's neighborhood grows, it changes that node's representation which in turn affects her social interactions (*association* → **embedding** → *communication*). Similarly, when node's interaction behavior changes, it affects the representation of her neighbors and herself which in turn changes the structure and strength of her connections due to link addition or deletion (*communication* → **embedding** → *association*). We call this phenomenon — *evolution through mediation* and illustrate it graphically in Figure 1.

In this work, we propose a novel representation learning framework for dynamic graphs, **DyRep**, to model interleaved evolution of two observed processes through latent mediation process expressed above and effectively learn richer node representations over time. Our framework ingests dynamic graph information in the form of association and communication events over time and updates the node representations as they appear in these events. We build a two-time scale deep temporal point

process approach to capture the continuous-time fine-grained temporal dynamics of the two observed processes. We further parameterize the conditional intensity function of the temporal point process with a deep inductive representation network that learns functions to compute node representations. Finally, we couple the structural and temporal components of our framework by designing a novel *Temporal Attention Mechanism*, which induces temporal attentiveness over neighborhood nodes using the learned conditional intensity function. This allows to capture highly interleaved and nonlinear dynamics governing node representations over time. We design an efficient unsupervised training procedure for end-to-end training of our framework. We demonstrate consistent and significant improvement over state-of-art representative baselines on two real-world dynamic graphs for the tasks of dynamic link prediction and time prediction. We also present an extensive qualitative analysis through embedding visualization and ablation studies to discern the effectiveness of our framework.

2 Background and Preliminaries

2.1 Related Work

Representation Learning approaches for static graphs either perform node embedding (Cao et al., 2015; Grover & Leskovec, 2016; Perozzi et al., 2014; Tang et al., 2015; Wang et al., 2016a; 2017; Xu et al., 2017) or sub-graph embedding (Scarselli et al., 2009; Li et al., 2016; Dai et al., 2016) which can also utilize convolutional neural networks (Kipf & Welling, 2017; 2016; Bruna et al., 2014). Among them, GraphSage (Hamilton et al., 2017a) is an inductive method for learning functions to compute node representations that can be generalized to unseen nodes. Most of these approaches only work with static graphs or can model evolving graphs without temporal information. Dynamic network embedding is pursued through various techniques such as matrix factorization (Zhu et al., 2016), structural properties (Zhou et al., 2018), CNN-based approaches (Seo et al., 2016), deep recurrent models (Trivedi et al., 2017), and random walks (Ngyuyen et al., 2018). There exists a rich body of literature on temporal modeling of dynamic networks (Kim et al., 2017), that focus on link prediction tasks but their goal is orthogonal to our work as they build task specific methods and do not focus on representation learning. Authors in (Yang et al., 2017; Sarkar et al., 2007) proposed models of learning dynamic embeddings but none of them consider time at finer level and do not capture both topological evolution and interactions simultaneously. In parallel, research on deep point process models include parametric approaches to learn intensity (Du et al., 2016; Mei & Eisner, 2017) using recurrent neural networks and GAN based approaches to learn intensity functions (Xiao et al., 2017). More detailed related works are provided in **Appendix D**.

2.2 Temporal Point Processes

Stochastic point processes (Daley & Vere-Jones, 2007) are random processes whose realization comprises of discrete events in time, $t_1, t_2, \ldots$. A temporal point process is one such stochastic process that can be equivalently represented as a counting process, $N(t)$, which contains the number of events up to time t. The common way to characterize temporal point processes is via the conditional intensity function $\lambda(t)$, a stochastic model of rate of happening events given the previous events. Formally, $\lambda(t)\mathrm{d}t$ is the conditional probability of observing an event in the tiny window $[t, t + \mathrm{d}t)$, $\lambda(t)\mathrm{d}t := \mathbb{P}\{\text{event in } [t, t + \mathrm{d}t)|\mathcal{T}(t)\} = \mathbb{E}[\mathrm{d}N(t)|\mathcal{T}(t)]$, where $\mathcal{T}(t) = t_k|t_k < t$ is history until t. Similarly, for $t > t_n$ and given history $\mathcal{T} = t_1, \ldots, t_n$, we characterize the conditional probability that no event happens during $[t_n, t)$ as $S(t|\mathcal{T}) = \exp\left(-\int_{t_n}^t \lambda(\tau)\,\mathrm{d}\tau\right)$, which is called survival function of the process (Aalen et al., 2008). Moreover, the conditional density that an event occurs at time t is defined as $f(t) = \lambda(t)\,S(t)$. The intensity $\lambda(t)$ is often designed to capture phenomena of interests – common forms include Poisson Process, Hawkes processes (Farajtabar et al., 2014; Hawkes, 1971; Wang et al., 2016b; Tabibian et al., 2017), Self-Correcting Process (Isham & Westcott, 1979). Temporal Point Processes have previously been used to model both – dynamics on the network (Farajtabar et al., 2016; Zarezade et al., 2017; Farajtabar et al., 2017) and dynamics of the network (Tran et al., 2015; Farajtabar et al., 2015).

2.3 Notations and Dynamic Graph Setting

Notations. Let $\mathcal{G}_t = (\mathcal{V}_t, \mathcal{E}_t)$ denote graph $\mathcal{G}$ at time t, where $\mathcal{V}_t$ is the set of nodes and $\mathcal{E}_t$ is the set of edges in $\mathcal{G}_t$. ***Event Observation*** **–** Both communication and association processes are realized in the form of dyadic events observed between nodes on graph $\mathcal{G}$ over a temporal window $[t_0, T]$ and ordered by time. We use the following canonical tuple representation for any type of event at time t of the form $e = (u, v, t, l, k)$, where u, v are the two nodes involved in an event. t represents

time of the event. $l \in \{0, 1\}$ represent link status – $l = 1$ signify a structural edge and $l = 0$ signify the two nodes are not directly connected. $k \in \{0, 1\}$ and $k = 0$ when the current event is an association and $k = 1$ when the current event is a communication. We then represent complete set of P observed events ordered by time in window $[0, T]$ as $\mathcal{O} = \{(u, v, t, l, k)_p\}_{p=1}^P$. Here, $t_p \in \mathbb{R}^+$, $0 \leq t_p \leq T$. **Node Representation—** Let $\mathbf{z}^v \in \mathbb{R}^d$ represent d-dimensional representation of node v. As the representations are evolving, we qualify them as function of time and denote it as $\mathbf{z}^v(t)$ which signifies the representation of node v is updated after an event involving v at time t. We use $\mathbf{z}^v(\bar{t})$ for most recently updated embedding of node v just before t.

Dynamic Graph Setting. Let $\mathcal{G}_{t_0} = (\mathcal{V}_{t_0}, \mathcal{E}_{t_0})$ be the initial snapshot of a graph at time t_0. Please note that $\mathcal{G}_{t_0}$ may be empty or it may contain an initial structure (association edges) but it will not have any communication history. Our framework observes evolution of graph as a stream of events $\mathcal{O}$ and hence any new node will always be observed as a part of such an event. This will induce a natural ordering over nodes as available from the data. As our method is inductive, we never learn node-specific representations and rather learn functions to compute node representations. In this work, we only support growth of network i.e. we only model addition of nodes and structural edges and leave deletion as future work. Further, for general description of the model, we will assume that an edge in the graph do not have types and nodes do not have attributes but we will discuss how to extend our model to accommodate these features. Finally, one can use $k > 1$ to signify more temporal scales for communication events (e.g. if there are different types of events in dynamic heterogeneous graphs and one wants to assign different scale corresponding to each type).

3 PROPOSED METHOD: DYREP

The key idea of DyRep is to build a unified architecture that can ingest evolving information over graphs and effectively model the *evolution through mediation* phenomenon described in Section 1. To achieve this, we design a two-time scale temporal point process model of observed processes and parameterize it with an inductive representation network which subsequently models the latent mediation process of learning node representations. The rationale behind our framework is that the observed set of events are the realizations of the nonlinear dynamic processes governing the changes in topological structure of graph and interactions between the nodes in the graph. Now, when an event is observed between two nodes, information flows from the neighborhood of one node to the other and affects the representations of the nodes accordingly. While a communication event (interaction) only propagates local information across two nodes, an association event changes the topology and thereby has more global effect. The goal is to learn node representations that encode information evolving due to such local and global effects and further drive the dynamics of the observed events.

3.1 MODELING TWO-TIME SCALE OBSERVED GRAPH DYNAMICS

The observations over dynamic graph contain temporal point patterns of two interleaved complex processes in the form of communication and association events respectively. At any time t, the occurrence of an event, from either of these processes, is dependent on the most recent state of the graph, *i.e.*, two nodes will participate in any event based on their most current representations. Given an observed event $p = (u, v, t, l, k)$, we define a continuous-time deep model of temporal point process using the conditional intensity function $\lambda_k^{u,v}(t)$ that models the occurrence of event p of dynamics k between nodes u and v at time t:

$$\lambda_k^{u,v}(t) = f_k(g_k(\bar{t})) \tag{1}$$

The inner function $g_k(\bar{t})$ computes the compatibility of the most recently updated representations of two nodes, $\mathbf{z}^u(\bar{t})$ and $\mathbf{z}^v(\bar{t})$ as follows:

$$g_k(\bar{t}) = \boldsymbol{\omega}_k^T \cdot [\mathbf{z}^u(\bar{t}); \mathbf{z}^v(\bar{t})] \tag{2}$$

Here, k represent the dynamic process (communication vs. association), [;] signifies concatenation and $\boldsymbol{\omega}_k \in \mathbb{R}^{2d}$ serves as the model parameter that learns time-scale specific compatibility. As $g_k(\bar{t})$ is a function of node representations learned through a representation network described in section 3.2, we note that the representation learning network parameterizes the conditional intensity function of the point process model which establishes the first unifying factor of our framework. Note that the dynamics are not two simple point processes dependent to each other, but, they are related through the mediation process and in the embedding space. Moreover, as we'll see a well curated attention mechanism is employed to learn how the past drives future.

The choice of outer function f_k needs to account for two critical criteria: 1.) Intensity needs to be positive. 2.) As mentioned before, the dynamics corresponding to communication and association

processes evolve at different time scales. To account for this, we use a modified version of softplus function parameterized by a dynamics parameter ψ_k to capture this *timescale* dependence:

$$f_k(x) = \psi_k \log(1 + \exp(x/\psi_k)) \tag{3}$$

where, $x = g(\bar{t})$ in our case and ψ_k is scalar time-scale parameter learned as part of training. ψ_k corresponds to the rate of events arising from a corresponding process. In 1D event sequences, the formulation in (3) corresponds to the nonlinear transfer function in (Mei & Eisner, 2017).

3.2 LEARNING LATENT MEDIATION PROCESS VIA TEMPORALLY ATTENTIVE REPRESENTATION NETWORK

We build a deep recurrent architecture that parameterizes the intensity function in Eq. (1) and learns functions to compute node representations. Specifically, after an event has occurred, the representation of both the participating nodes need to be updated to capture the effect of the observed event based on the principles of:

Self-Propagation. Self-propagation can be considered as a minimal component of the dynamics governing an individual node's evolution. A node evolves in the embedded space with respect to its previous position (e.g. set of features) and not in a random fashion.

Exogenous Drive. Some exogenous force may smoothly update the node's current features during the time interval (e.g. between two global events involving that node).

Localized Embedding Propagation. Two nodes involved in an event form a temporary (communication) or a permanent (association) pathway for the information to propagate from the neighborhood of one node to the other node. This corresponds to the influence of the nodes at second-order proximity passing through the other node participating in the event.

To realize the above processes in our setting, we first describe an example setup: Consider nodes u and v participating in any type of event at time t. Let $\mathcal{N}_u$ and $\mathcal{N}_v$ denote the neighborhood of nodes u and v respectively. We discuss two key points here: 1.) Node u serves as a bridge passing information from $\mathcal{N}_u$ to node v and hence v receives the information in an aggregated form through u. 2.) While each neighbor of u wants to pass its information to v, the information that node u captures is governed by an aggregate function parametrized by u's communication and association history with its neighbors.

With this setup, for any event at time t, we update the embeddings for both nodes involved in the event using a recurrent architecture. Specifically, for p-th event of node v, we evolve $\mathbf{z}^v$ as:

$$\mathbf{z}^v(t_p) = \sigma(\ \underbrace{\mathbf{W}^{struct}\mathbf{h}^u_{struct}(\bar{t}_p)}_{\text{Localized Embedding Propagation}}\ + \underbrace{\mathbf{W}^{rec}\mathbf{z}^v(t_{p-1})}_{\text{Self-Propagation}} + \underbrace{\mathbf{W}^t(t_p - t_{p-1})}_{\text{Exogenous Drive}}), \tag{4}$$

where, $\mathbf{h}^u_{struct} \in \mathbb{R}^d$ is the output representation vectors obtained from aggregator function on node u's neighborhood and $\mathbf{z}^v(t_{p-1})$ is the recurrent state obtained from the previous representation of node v. t_p is time point of current event, $\bar{t}_p$ signifies the timepoint just before current event and t_{p-1} represent time point of previous event for node v. $\mathbf{z}^v(t_0)$, the representation of a node v at initial time point t_0 may be initialized either using input node features from dataset or random vector as per the setting. $\mathbf{W}^{struct}, \mathbf{W}^{rec} \in \mathbb{R}^{d \times d}$ and $\mathbf{W}^t \in \mathbb{R}^d$ are model parameters that govern the aggregate effect of all the three processes respectively. The above formulation is flexible in supporting various features. For example, in case of dynamic heterogenous graphs, Eq. 2 and Eq. 4 can be extended to include edge type with corresponding parameters. Similarly, one can support high-dimensional attributes by adding them to Eq. 4 as a feature.

3.2.1 TEMPORALLY ATTENTIVE AGGREGATION

In a dynamic graph, any event would not only affect the nodes involved in that event but also the local neighbors of those nodes. Aligned with our motivation, any event would impact the embeddings of the nodes involved in those events and the updated embeddings would further induce more events across graph. To capture this mutual effect, we design a function that learns structural information local to a node based on historical events on its neighboring nodes. To this effect, we propose a novel *Temporal Point Process based Attentional Mechanism* that empowers the aggregate function (used to compute $\mathbf{h_{struct}}$) to attend to the neighbors based on node's communication and association history.

Let $\mathbf{A}(t) \in \mathbb{R}^{n \times n}$ be the adjacency matrix for graph $\mathcal{G}_t$ at time t. Let $\mathcal{S}(t) \in \mathbb{R}^{n \times n}$ be a stochastic matrix capturing the strength between pair of vertices at time t. One can consider $\mathcal{S}$ as a selection

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Algorithm 1 Update Algorithm for $\mathcal{S}$ and $\mathbf{A}$

Input: Event record $o = (u, v, t, l, k)$, Event Intensity $\lambda_k^{u,v}(t)$ computed in (1), most recently updated $\mathbf{A}(\bar{t})$ and $\mathcal{S}(\bar{t})$. **Output:** $\mathbf{A}(t)$ and $\mathcal{S}(t)$

1. Update $\mathbf{A}$: $\mathbf{A}(t) = \mathbf{A}(\bar{t})$
if $k = 0$ **then** $\mathbf{A}_{uv}(t) = \mathbf{A}_{vu}(t) = 1$ ←{*Association event*}

2. Update $\mathcal{S}$: $\mathcal{S}(t) = \mathcal{S}(\bar{t})$
if $k = 1$ and $l = 0$ **then return** $\mathcal{S}(t), \mathbf{A}(t)$ ←{*Communication event, no Association exists*}
for $j \in \{u, v\}$ **do**
    $b = \frac{1}{|\mathcal{N}_j(t)|}$ where $|\mathcal{N}_j(t)|$ is the size of $\mathcal{N}_j(t) = \{i : \mathbf{A}_{ij}(t) = 1\}$
    $\mathbf{z} \leftarrow \mathcal{S}_j(t)$
    if $k = 1$ and $l = 1$ **then** {   ←{*Communication event, Association exists*}}
        $\mathbf{z}_i = b + \lambda_k^{ji}(t)$ where i is the other node involved in the event.   ←{λ *computed in Eq. 2*}
    else if $k = 0$ and $l = 0$ **then** {   ←{*Association event*}}
        $b' = \frac{1}{|\mathcal{N}_j(t-)|}$ where $|\mathcal{N}_j(t-)|$ is the size of $\mathcal{N}_j(t-) = \{i : \mathbf{A}_{ij}(t-) = 1\}$
        $x = b' - b$
        $\mathbf{z}_i = b + \lambda_k^{ji}(t)$ where i is the other node involved in the event   ←{λ *computed in Eq. 2*}
        $\mathbf{z}_w = \mathbf{z}_w - x; \ \ \forall w \neq u$
    end if
    Normalize $\mathbf{z}$ and set $\mathcal{S}_j(t) \leftarrow \mathbf{z}$
end for
return $\mathcal{S}(t), \mathbf{A}(t)$

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matrix that induces a ***natural selection*** process for a node – it would tend to communicate more with other nodes that it wants to associate with or has recently associated with. And it would want to attend less to non-interesting nodes. We start with following implication required for the construction of $\mathbf{h}_{struct}^u$ in (4): For any two nodes u and v at time t, $\mathcal{S}_{uv}(t) \in [0, 1]$ if $\mathbf{A}_{uv}(t) = 1$ and $\mathcal{S}_{uv}(t) = 0$ if $\mathbf{A}_{uv}(t) = 0$. Denote $\mathcal{N}_u(t) = \{i : \mathbf{A}_{iu}(t) = 1\}$ as the 1-hop neighborhood of node u at time t.

To formally capture the difference in the influence of different neighbors, we propose a novel *conditional intensity based attention layer* that uses the matrix $\mathcal{S}$ to induce a shared attention mechanism to compute attention coefficients over neighborhood. Specifically, we perform *localized attention* for a given node u and compute the coefficients pertaining to the 1-hop neighbors i of node u as: $q_{ui}(t) = \frac{\exp(\mathcal{S}_{ui}(\bar{t}))}{\sum_{i' \in \mathcal{N}_u(t)} \exp(\mathcal{S}_{ui'}(\bar{t}))}$, where q_{vi} signifies the attention weight for the neighbor i at time t and hence it is a temporally evolving quantity. These attention coefficients are then used to compute the aggregate information $\mathbf{h}_{struct}^u(\bar{t})$ for node v by employing an attended aggregation mechanism across neighbors as follows: $\mathbf{h}_{struct}^u(\bar{t}) = \max\left(\left\{\sigma\left(q_{ui}(t) \cdot \mathbf{h}^i(\bar{t})\right), \forall i \in \mathcal{N}_u(\bar{t})\right\}\right)$, where, $\mathbf{h}^i(\bar{t}) = \mathbf{W}^h \mathbf{z}^i(\bar{t}) + \mathbf{b}^h$ and $\mathbf{W}^h \in \mathbb{R}^{d \times d}$ and $\mathbf{b}^h \in \mathbb{R}^d$ are parameters governing the information propagated by each neighbor of u. $\mathbf{z}^i(\bar{t}) \in \mathbb{R}^d$ is the most recent embedding for node i. The use of $\max$ operator is inspired from learning on general point sets (Qi et al., 2017). By applying max-pooling operator element-wise, the model effectively captures different aspects of the neighborhood. We found $\max$ to work slightly better as it considers temporal aspect of neighborhood which would be amortized if $mean$ is used instead.

Connection to Neural Attention over Graphs. Our proposed temporal attention layer shares the motivation of recently proposed Graph Attention Networks (GAT) (Veličković et al., 2018) and Gated Attention Networks (GaAN) (Zhang et al., 2018) in the spirit of applying non-uniform attention over neighborhood. Both GAT and GaAN has demonstrated significant success in static graph setting. GAT advances GraphSage (Hamilton et al., 2017a) by employing multi-head non-uniform attention over neighborhood and GaAN advances GAT by applying different weights to different heads in the multi-head attention formulation. The key innovation in our model is the parameterization of attention mechanism by a point process based temporal quantity $\mathcal{S}$ that is evolving and drives the impact that each neighbor has on the given node. Further, unlike static methods, we use these attention coefficients as input to the aggregator function for computing the temporal-structural effect of neighborhood. Finally, static methods use multi-head attention to stabilize learning by capturing

multiple representation spaces but this is an inherent property in our layer as representations and event intensities update over time and hence new events help capture multiple representation spaces.

Construction and Update of $\mathcal{S}$. We construct a single stochastic matrix $\mathcal{S}$ (used to parameterize attention in the earlier section) to capture complex temporal information. As before, $\mathbf{A}(t) \in \mathbb{R}^{n \times n}$ and $\mathcal{S}(t) \in \mathbb{R}^{n \times n}$ are the adjacency matrix and a right-stochastic matrix at time t. At the initial timepoint $t = t_0$, we construct $\mathcal{S}(t_0)$ directly from $\mathbf{A}(t_0)$. Specifically, for a given node v, we initialize the elements of corresponding row vector $\mathcal{S}_v(t_0)$ as: $\mathcal{S}_{vu}(t_0) = 0$ if $(v = u$ or $\mathbf{A}_{vu}(t_0) = 0)$ and $\mathcal{S}_{vu}(t_0) = \frac{1}{|\mathcal{N}_v(t_0)|}$ if $\mathcal{N}_v(t_0) = \{u : \mathbf{A}_{uv}(t_0) = 1\}$. After observing an event $o = (u, v, t, l, k)$ at time $t > t_0$, we make updates to $\mathbf{A}$ and $\mathcal{S}$ as per the observation of l and k. Specifically, $\mathbf{A}$ only gets updated for association events (k=0, change in structure). $\mathcal{S}$ is used for capturing temporal attention over structure and hence its values are active (non-zero)/updated only for node pairs who have an association edge (k=1, l=1 or k=0, l=0). There will be no effect in $\mathcal{S}$ for communication events happening between disconnected nodes (k=1,l=0). Further, for any two nodes i and j, update to $\mathcal{S}$ is based on respective node's existing structural neighborhood size (b) and the intensity of event between them at time t. From mathematical viewpoint, this update resembles a standard temporal point process formulation where the term coming from b serves as background exogenous intensity while λ can be viewed as endogenous intensity. The updates are asymmetric due to different neighborhood size and hence happen to two rows of $\mathcal{S}$ for any event. Finally, for association event (k=0), the neighborhood size for both nodes increases and hence other nodes (not the ones involved in event) in the neighborhood needs to be updated to accommodate this change (using x). Algorithm 1 outlines complete update scenarios. We provide an extensive ablation study in **Appendix A** that can help discern the contribution of all the above components in achieving our goal.

4 EFFICIENT LEARNING PROCEDURE

The complete parameter space for the current model is $\boldsymbol{\Omega} = \{\mathbf{W}^{struct}, \mathbf{W}^{rec}, \mathbf{W}^t, \mathbf{W}^h, \mathbf{b}^h,$ $\{\boldsymbol{\omega}_k\}_{k=0,1}, \{\psi_k\}_{k=0,1}\}$. For a set $\mathcal{O}$ of P observed events, we learn these parameters by minimizing the negative log likelihood: $\mathcal{L} = -\sum_{p=1}^{P} \log(\lambda_p(t)) + \int_0^T \Lambda(\tau)d\tau$, where $\lambda_p(t) = \lambda_{k_p}^{u_p, v_p}(t)$ represent the intensity of event at time t and $\Lambda(\tau) = \sum_{u=1}^{n} \sum_{v=1}^{n} \sum_{k \in \{0,1\}} \lambda_k^{u,v}(\tau)$ represent total survival probability for events that do not happen. While it is intractable (will require $\mathcal{O}(n^2 k)$ time) and unnecessary to compute the integral in the log-likelihood equation for all possible non-events in a stochastic setting, we can locally optimize $\mathcal{L}$ using mini-batch stochastic gradient descent where we estimate the integral using novel sampling technique. Algorithm 2 in **Appendix F** adopts a simple variant of Monte Carlo trick to compute the survival term of log-likelihood equation. Specifically, in each mini-batch, we sample non-events instead of considering all pairs of non-events (which can be millions). Let m be the mini-batch size and N be the number of samples. The complexity of Algorithm 2 will then be $\mathcal{O}(2mkN)$ for the batch where the factor of 2 accounts for the update happening for two nodes per event which demonstrates linear scalability in number of events which is desired to tackle web-scale dynamic networks (Paranjape et al., 2017). The overall training procedure is adopted from (Trivedi et al., 2017) where the Backpropagation Through Time (BPTT) training is conducted over a global sequence, thereby maintaining the dependencies between events across sequences while avoiding gradient related issues. Implementation details are left to **Appendix E**.

5 EXPERIMENTS

5.1 DATASETS

We evaluate DyRep and baselines on two real world datasets: **Social Evolution Dataset** released by MIT Human Dynamics Lab — #nodes: 100, #Initial Associations: 407, #Final Associations: 809, #Communications: 2020554 and Clustering Coefficient: 0.548. **Github Dataset** available at Github Archive — #nodes: 12328, #Initial Associations: 70640, #Final Associations: 166565, #Communications: 604954 and Clustering Coefficient: 0.087. These datasets cover a range of configurations as Social Dataset is a small network with high clustering coefficient and over 2M events. In contrast, Github dataset forms a large network with low clustering coefficient and sparse events thus allowing us to test the robustness of our model. Further, Github dataset contains several unseen nodes which were never encountered during training.

5.2 TASKS AND METRICS

We study the effectiveness of DyRep by evaluating our model on tasks of dynamic link prediction and event time prediction tasks:

Table 1: Comparison of DyRep with state-of-art approaches

Key Properties	DyRep (Our Method)	Know-Evolve (Dynamic)	DynGem (Dynamic)	GraphSage (Static)	GAT (Static)
Models Association	✓	X	✓	✓	✓
Models Communication	✓	✓	X	X	X
Models Time	✓	✓	X	X	X
Learns Representation	✓	✓	✓	✓	✓
Predicts Time	✓	✓	X	X	X
Graph Information	2nd-order Neighborhood	Single Edge	1st and 2nd-order Neighborhood	2nd-order Neighborhood	1st-order Neighborhood
Attention Mechanism	Temporal Point Process (Non-Uniform)	None	None	Sampling (Uniform)	Multi-head (Non-Uniform)
Learning	Unsupervised	Unsupervised	Semi-Supervised	Unsupervised	Supervised

Dynamic Link Prediction. When any two nodes in a graph has increased rate of interaction events, they are more likely to get involved in further interactions and eventually these interactions may lead to the formation of structural link between them. Similarly, formation of the structural link may lead to increased likelihood of interactions between newly connected nodes. To understand, how well our model captures these phenomenon, we ask questions like: 1.) Which is the most likely node u that would undergo an event of type k with a given node v at time t? or 2.) Given two nodes u and v at time t, which is most likely event type they will participate in? Its conditional density at time t can be computed: $f_k^{u,v}(t) = \lambda_k^{u,v}(t) \cdot \exp\left(\int_{\bar{t}}^{t} \lambda(s)ds\right)$, where $\bar{t}$ is the time of the most recent event on either dimension u or v. We use this conditional density to find most likely node and event type.

For a given test record (u, v, t, l, k), we replace v with other entities in the graph and compute the density as above. We then rank all the entities in descending order of the density and report the rank of the ground truth entity. Please note that the latest embeddings of the nodes update even during the test while the parameters of the model remaining fixed. Hence, when ranking the entities, we remove any entities that creates a pair already seen in the test. We report Mean Average Rank and HITS(@10) metric for dynamic link prediction.

Event Time Prediction. This is a relatively novel application where the aim is to compute the next time point when a particular type of event (structural or interaction) can occur. Given a pair of nodes (u, v) and event type k at time t, we use the above density formulation to compute conditional density at time t. The next time point $\hat{t}$ for the event can then be computed as: $\hat{t} = \int_t^{\infty} t f_k^{u,v}(t)dt$ where the integral does not have an analytic form and hence we estimate it using Monte Carlo trick. For a given test record (u, v, t, l, k), we report the next time this communication event may occur. We compute the next time point using the previous scheme and report MAE against the ground truth.

5.3 BASELINES

Dynamic Link Prediction. We compare the performance of our model against multiple representation learning baselines, four of which has capability to model evolving graphs. Specifically, we compare with **Know-Evolve** (Trivedi et al., 2017)— a state-of-art model for multi-relational dynamic graphs where each edge has time-stamp and type (communication events), **DynGem** (Goyal et al., 2017)—divides timeline into discrete time points and learns embedding for the graph snapshots at these time points. **DynTrd** (Zhou et al., 2018) focuses on specific structure of triad to model how close triads are formed from open triads in dynamic networks. **GraphSage** (Hamilton et al., 2017a)— an inductive representation learning method that learns sample and aggregation functions to learn representations instead of training for individual node. **Node2Vec** (Grover & Leskovec, 2016)—simple transductive baseline to learn graph embeddings over static graphs. Table 1 provides qualitative comparison between state-of-arts and our framework. In our experiments, we compare with GraphSage instead of GAT as we share the unsupervised setting with GraphSage while GAT is designed for supervised learning. In Appendix A (Ablation studies), we show results on one version where we only update attention based on Association events which is temporal analogous to GAT.

Event Time Prediction. We compare our model against (i) Know-Evolve which has the ability to predict time in a multi-relational dynamic graphs (II) Multi-dimensional Hawkes Process (MHP) (Du et al., 2015) model where all events in graph are considered as dyadic.
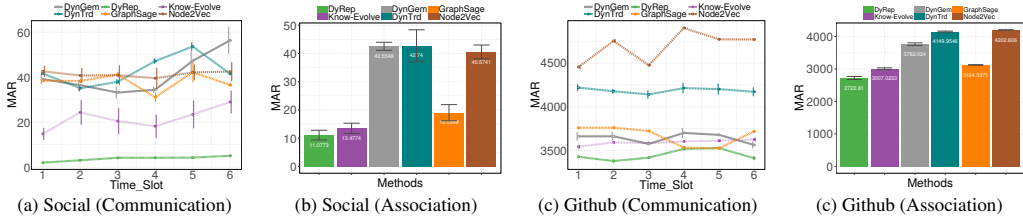
Figure 2: Dynamic Link Prediction Performance for **(a-b)** Social Evolution Dataset **(c-d)** Github Dataset. We report HITS@10 results and zoomed versions in **Appendix C**. Best viewed in pdf.
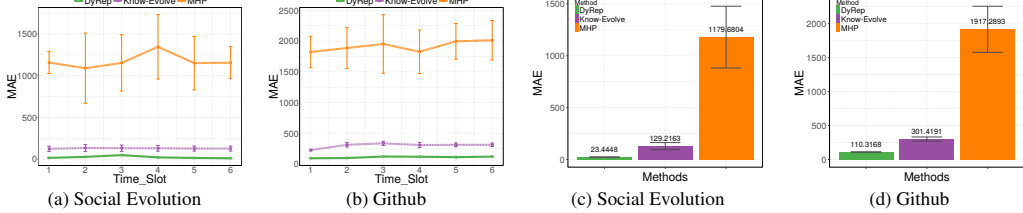

Figure 3: Time Prediction Performance (unit is hrs). Figure best viewed in pdf or colored print.

5.4 EVALUATION SCHEME

We divide our test sets into $n(=6)$ slots based on time and report the performance for each time slot, thus providing comprehensive temporal evaluation of different methods. This method of reporting is expected to provide fine-grained insights on how various methods perform over time as they move farther from the learned training history. For dynamic baselines that do not explicitly model time (DynGem, DynTrd, GraphSage) and static baselines (Node2Vec), we adopt a sliding window training approach with warm-start method where we learn on initial train set and test for the first slot. Then we add the data from first slot in the train set and remove equal amount of data from start of train set and retrain the model using the embeddings from previous train.

5.5 EXPERIMENTAL RESULTS

Communication Event Prediction Performance. We first consider the task of predicting communication events between nodes which may or may not have a permanent edge (association) between them. Figure 2 **(a-b)** shows corresponding results.

Social Evolution. Our method significantly and consistently outperforms all the baselines on both metrics. While the performance of our method drops a little over time, it is expected due to the temporal recency affect on node's evolution. Know-Evolve can capture event dynamics well and shows consistently better rank than others but its performance deteriorates significantly in HITS@10 metric over time. We conjecture that features learned through edge-level modeling limits the predictive capacity of the method over time. The inability of DynGem (snapshot based dynamic), DynTrd and GraphSage (inductive) to significantly outperform Node2vec (transductive static baseline) demonstrate that discrete time snapshot based models fail to capture fine-grained dynamics of communication events.

Github dataset. We demonstrate comparable performance with both Know-Evolve and GraphSage on Rank metric. We would like to note that overall performance for all methods on rank metric is low. As we reported earlier, Github dataset is very sparse with very low clustering coefficient which makes it a challenging dataset to learn. It is expected that for a large number of nodes with no communication history, most of the methods will show comparable performance but our method outperforms all others when there is some history available. This is demonstrated by our significantly better performance for HITS@10 metric where we are able to do highly accurate prediction for nodes where we learn better history. This can also be attributed to our model's ability to capture the effect of evolving topology which is missed by Know-Evolve. Finally, we do not see significant decrease in performance of any method over time in this case which can again be attributed to roughly uniform distribution of nodes with no communication history across time slots.

Association Event Prediction Performance. Association events are not available for all time slots so Figure 2 **(c-d)** report the aggregate number for this task. For both the datasets, our model significantly outperforms the baselines for this task. Specifically, our model's strong performance on HITS@10 metric across both datasets demonstrates its robustness in accurate learning from various

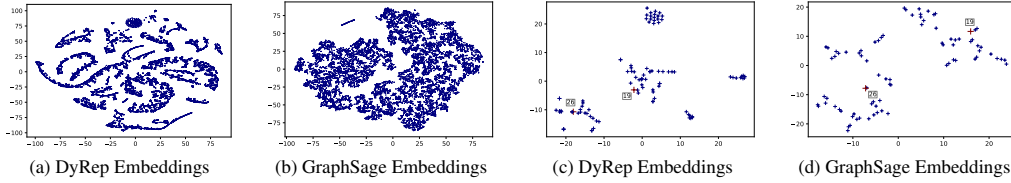(a) DyRep Embeddings   (b) GraphSage Embeddings   (c) DyRep Embeddings   (d) GraphSage Embeddings

Figure 4: tSNE for learned embeddings after training. Figure best viewed in color.

properties of data. On Social evolution dataset, the number of association events are very small (only 485) and hence our strong performance shows that the model is able to capture the influence of communication events on the association events through the learned representations (mediation). On the Github dataset, the network grows through new nodes and our model's strong performance across both metric demonstrates its inductive ability to generalize across new nodes across time. An interesting observation was poor performance of DynTrd which seems to be due to its objective to complete triangles. Github dataset is very sparse and has very few possibilities for triadic closure.

Time Prediction Performance. Figure 3 demonstrates consistently better performance than state-of-art baseline for event time prediction on both datasets. While Know-Evolve models both processes as two different relations between entities, it does not explicitly capture the variance in the time scales of two processes. Further, Know-Evolve does not consider influence of neighborhood which may lead to capturing weaker temporal-structural dynamics across the graph. MHP uses specific parametric intensity function which fails to account for intricate dependencies across graph.

Qualitative Performance. We conducted a series of qualitative analysis to understand the discriminative power of evolving embeddings learned by DyRep. We compare our embeddings against GraphSage embeddings as it is state-of-art embedding method that is also inductive. Figure 4 **(a-b)** shows the tSNE embeddings learned by Dyrep (left) and GraphSage (right) respectively. The visualization demonstrates that DyRep embeddings have more discriminative power as it can effectively capture the distinctive and evolving structural features over time as aligned with empirical evidence. Figure 4 **(c-d)** shows use case of two associated nodes (19 and 26) that has less communication at the two time points for above two methods. DyRep keeps the embeddings nearby although not in same cluster (cos. dist. - 0.649) which demonstrates its ability to learn the association and less communication dynamics between two nodes. For GraphSage the embeddings are on opposite ends of cluster with (cos. dist. - 1.964). We provide more extensive analysis in **Appendix B**.

6 CONCLUSION

We introduced a novel modeling framework for dynamic graphs that effectively and efficiently learns node representations by posing representation learning as latent mediation process bridging dynamic processes of topological evolution and node interactions. We proposed a deep temporal point process model parameterized by temporally attentive representation network that models these complex and nonlinearly evolving dynamic processes and learns to encode structural-temporal information over graph into low dimensional representations. Our superior evaluation performance demonstrates the effectiveness of our approach compared to state-of-arts. We present this work as the first generic and unified representation learning framework that adopts a novel modeling paradigm for dynamic graphs and support wide range of dynamic graph characteristics which can potentially have many exciting adaptations. As a part of our framework, we also propose a novel temporal point process based attention mechanism that can attend over neighborhood based on the history of communications and association events in the graph.

Currently, DyRep does not support network shrinkage due to following reasons: (i) It is difficult to procure data with fine grained deletion time stamps and (ii) The temporal point process model requires more sophistication to support deletion. For example, one can augment the model with a survival process formulation to account for lack of node/edge at future time. Another interesting future directions could be to support encoding higher order dynamic structures.

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APPENDIX

A   ABLATION STUDY

DyRep framework unifies several components that contribute to its effectiveness in learning rich node representation over complex and nonlinear processes in dynamic graphs. In this section, we provide insights on each component and how it is indispensable to the learning mechanism by performing an ablation study on various design choices of our model. Specifically, DyRep can be divided into three main parts: **Multi-time scale point process model**, **Representation Update Formulation** and **Conditional Intensity Based Attention Mechanism**. We focus on design choices available in each component and evaluate them on large github dataset. DyRep in the Figure 5 is the full model.

Multiple Time- Scale Processes. For this component, we perform two major tests:

- **DyRep-Comm.** In this variant, we make Eq 1., time-scale independent (i.e. remove k) and we train on only Communication Events. But we evaluate on both communication and association events. Please note that this is possible as our framework can compute representations for unseen nodes. Hence during training they will only learn representation parameters based on communication events. It is observed that compared to the full model, the performance of model degrades in prediction for both types of events. But the decline is more prominent for the Association events compared to Communication Events.
- **DyRep-Assoc.** In this variant, similar to above, we make Eq 1., time-scale independent and we train on only Association Events. But we evaluate on both communication and association events. It is observed that compared to the full model, the performance of model degrades in prediction for both types of events. But the decline is more prominent for the Communication events compared to Association Events.

The above two experiments show that considering events at a single time scale and not distinguishing between the processes hurt the performance. Although the performance is hurt more when communication events are not considered which may be due to the more availability of communication events due to its rapid frequency. We also performed a small test by training on all events but using a single scale parameter (ψ). The performance for both the dynamics degrades which demonstrates the effectiveness of ψ_k.

Representation Update Formulation. For this component, we focus on Eq. 4 and switch off the components to observe its effect.

- **DyRep-No-SP.** In this variant, we switch off the self-propagation component and we observe that the overall performance is not hurt significantly by not using self-propagation. In general, this term provides a very weak feature and mainly captures the recurrent evolution of one's own latent features independent of others. It is observed that the deviation has increased for Association events which may point to the reason that there are few nodes who have links but highly varying frequency of communication and hence most of their features are either self-propagated or completely associated with others.
- **DyRep-No-Struct.** In this variant, we remove the structural part of the model and as one would expect, the performance drops drastically in both the scenarios. This provides evidence to the necessity of building sophisticated structural encoders for dynamic graphs.

Intensity Attention Mechanism. For this component, we focus on Section 3.2 which builds the novel intensity based attention mechanism. Specifically, we carry following test:

- **DyRep-No-Att.** Here we completely remove the attention from the structural component and we see a significant drop in the performance.
- **DyRep-S-Comm.** In this variant, we focus on Algorithm 1 and we only make update to the $\mathcal{S}$ matrix for Communication events but do not do it for Association events. This leads to slightly worse performance which helps to see how the $\mathcal{S}$ matrix is helping to mediate the two processes and not considering association events leads to loss of information.
- **DyRep-S-Assoc.** In this variant, we focus on Algorithm 1 and we only make update to the $\mathcal{S}$ matrix for Association events but do not do it for Communication events. This leads to

(a) MAR (Communication)

(b) HITS@10 (Communication)
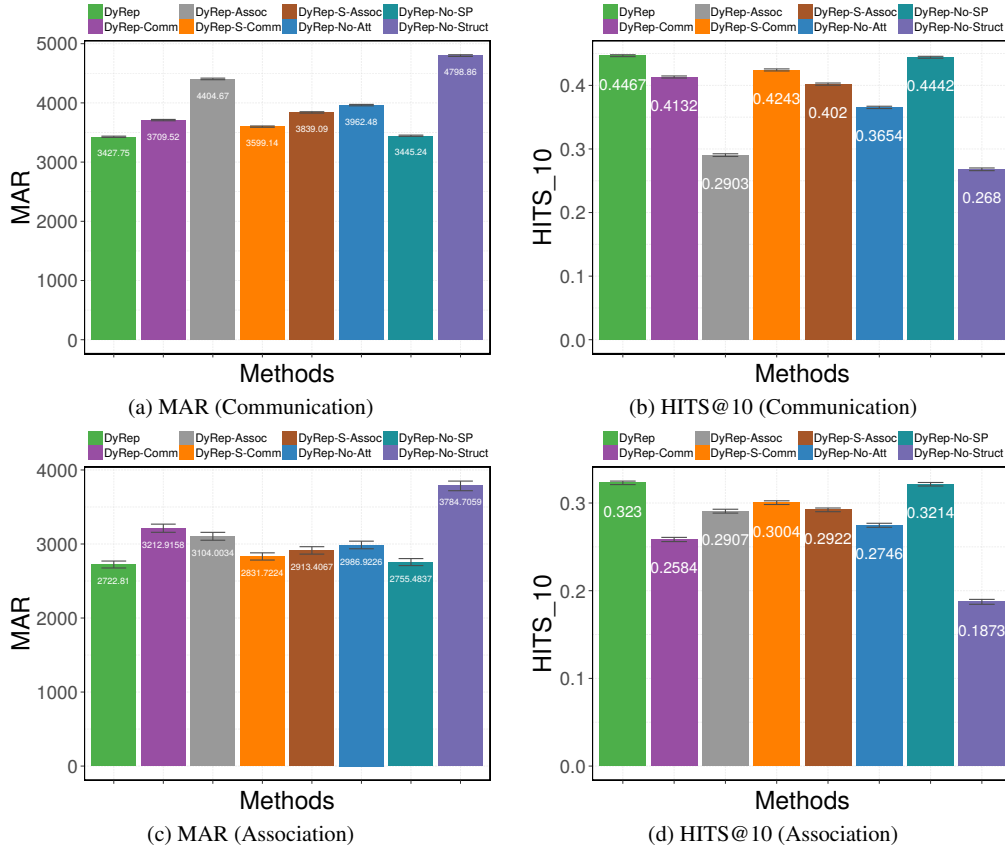
(c) MAR (Association)

(d) HITS@10 (Association)

Figure 5: Ablation Study on Github Dataset

a significant drop in performance again validating the need for using both processes but its prominent effect also shows that communication events (dynamics on the network) is more important while considering the influence of neighbors. Please note that this version is temporal analogous of GAT.

B EXPLORATORY ANALYSIS

We assess the quality of learned embeddings and the ability of model to capture both temporal and structural information. Let t_0 be the time point when train ended. Let t_1 be the timepoint when the first test slot ends.

Effect of Association and Communication on Embeddings. We conducted this experiment on Social dataset. We consider three use cases to demonstrate how the interactions and associations between the nodes changed their representations and visualize them to realize the effect.

- **Nodes that did not have association before test but got linked during first test slot.** Nodes 46 and 76 got associated in test between test points 0 and 1. This reduced the cosine distance in both models but DyRep shows prominent effect of this association which should be the case. DyRep reduces the cosine distance from 1.231 to 0.005. Also, DyRep embeddings for these two points belong to different clusters initially but later converge to same cluster. In GraphSage, the cosine distance reduces from 1.011 to 0.199 and the embeddings still remain in original clusters. Figure 6 shows the visualization of embeddings at the two time points in both the methods. This demonstrates that our embeddings can capture association events effectively.

- **Nodes that did not have association but many communication events (114000).** Nodes 27 and 70 is such a use case. DyRep embeddings consider the nodes to be in top 5 nearest

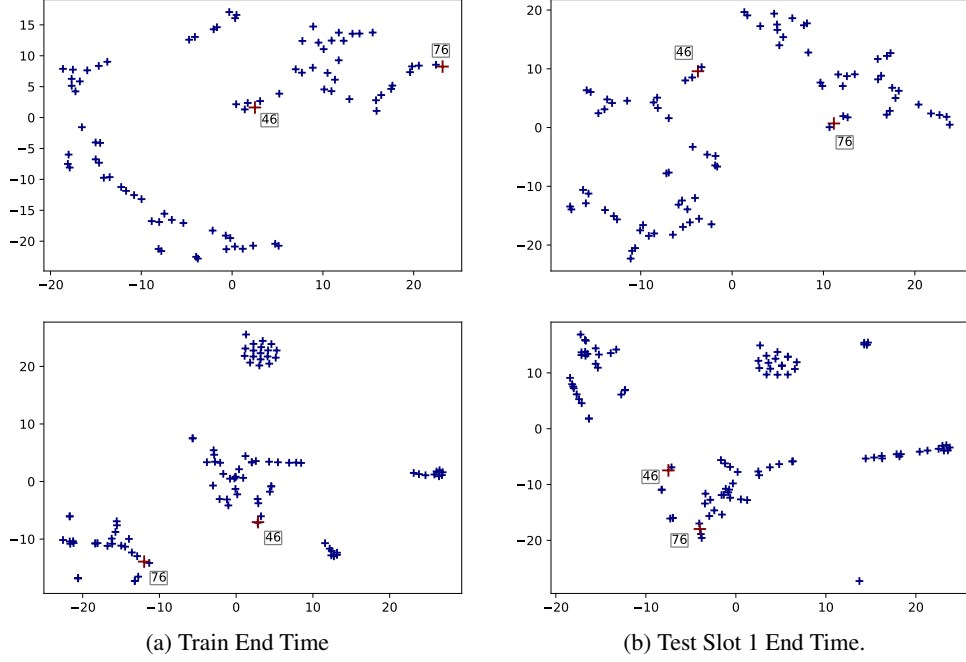(a) Train End Time

(b) Test Slot 1 End Time.

Figure 6: Use Case I. **Top row:** GraphSage Embeddings. **Bottom Row:** DyRep Embeddings.

neighbor of each other, in the same cluster and cosine distance of 0.005 which is aligned with the fact that nodes with large number of events tend to develop similar features over time. Graphsage on the other hand considers them 32nd nearest neighbor, puts them in different clusters with cosine distance - 0.792. Figure 7 shows the visualization of embeddings at the two time points in both the methods. This demonstrates the ability of DyRep's embedding to capture communication events and their temporal effect on embeddings effectively.
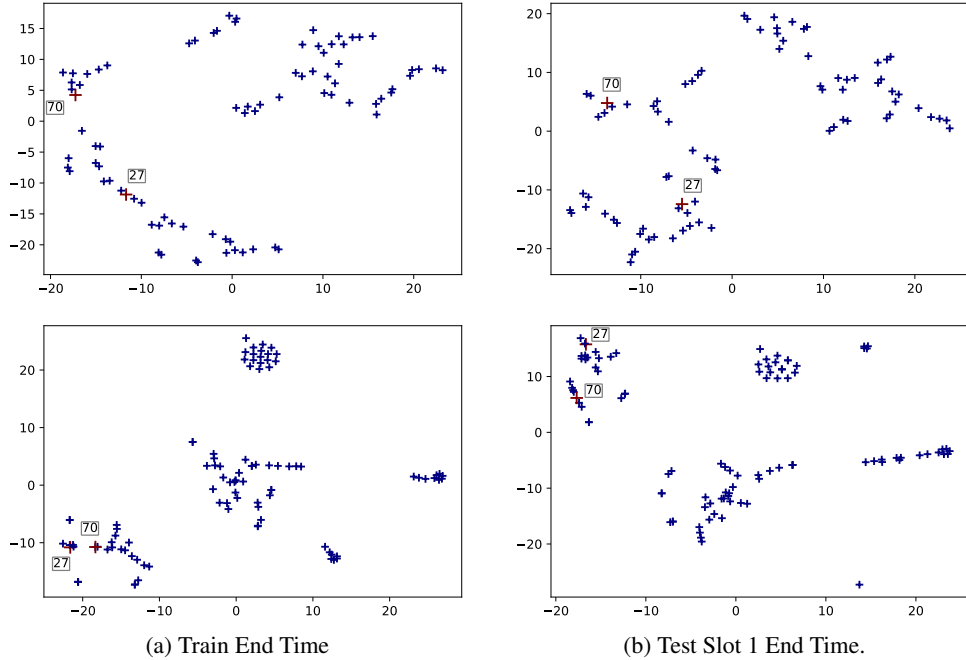

(a) Train End Time

(b) Test Slot 1 End Time.

Figure 7: Use Case II. **Top row:** GraphSage Embeddings. **Bottom Row:** DyRep Embeddings.

- **Temporal evolution of DyRep embeddings.** In figure 8 we visualize the embedding positions of the nodes (tracked in red) as they evolve through time and forms and breaks from clusters.
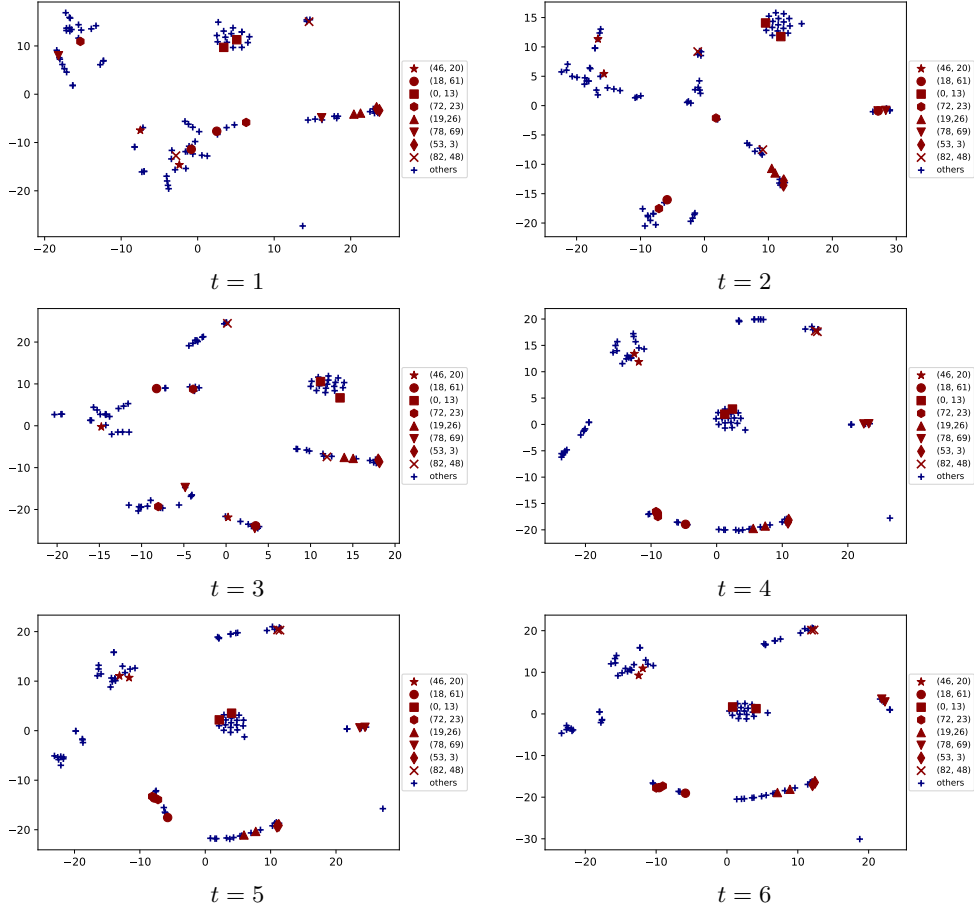

Figure 8: Use Case IV: DyRep Embeddings over time - From left to right and top to bottom. t are the timepoints when test with that id ended. Hence, $t = 1$ means the time when test slot 1 finished.

C Full Experiment Results for both Datasets

Figure 9 provides HITS@10 results in addition to the MAR results reported for Link Prediction in Section 5 (Experiments) of the main paper.
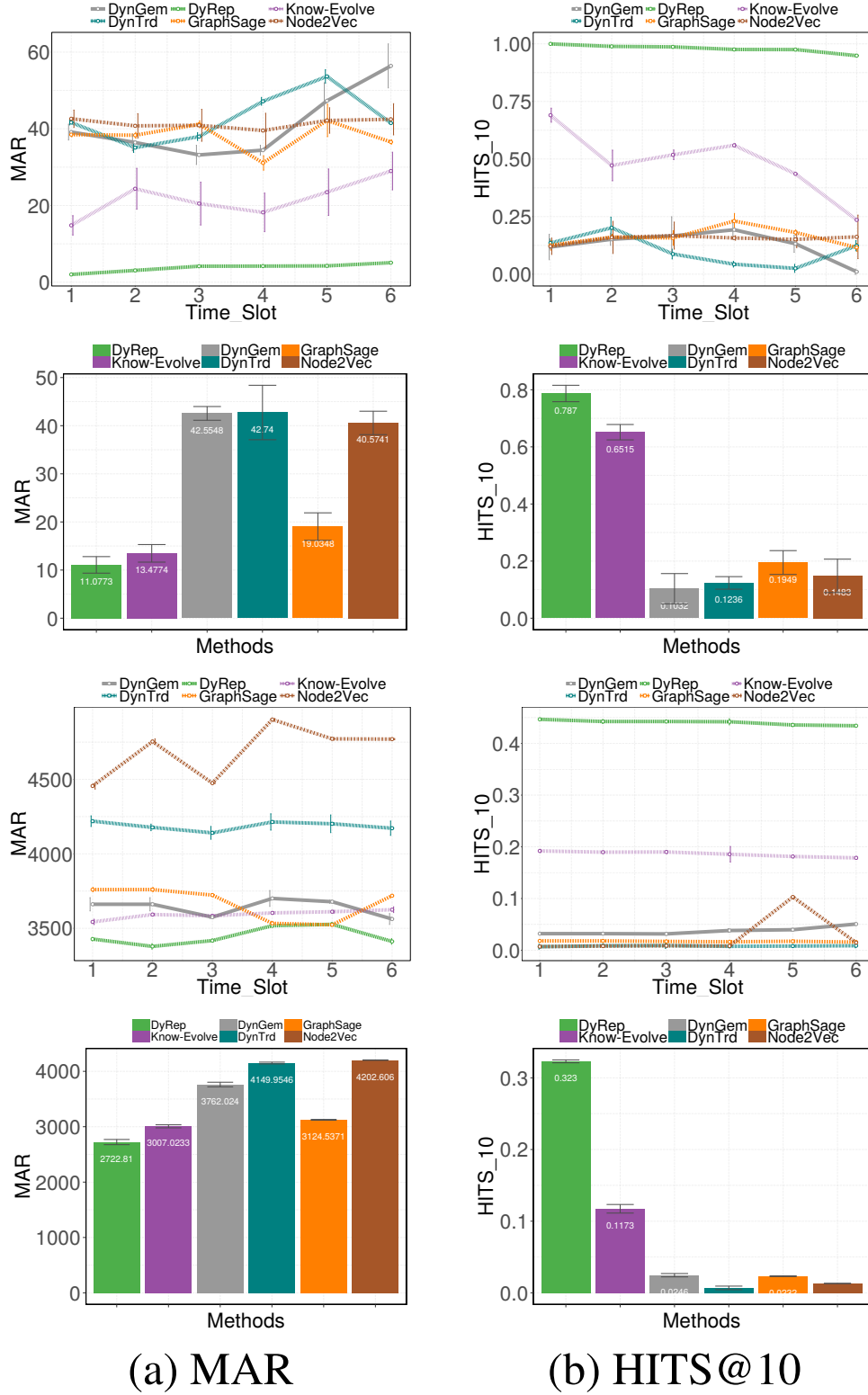
(a) MAR (b) HITS@10

Figure 9: Dynamic Link Prediction Performance: **Top 2 rows** show performance for Social Evolution Dataset. **Bottom 2 rows** show performance for Github Dataset. 1st and 3rd row show performance for Communication Events while 2nd and 4th row show performance for Association Events.

D   DETAILED RELATED WORK

Static Embedding Approaches. Representation Learning approaches for static graphs can be broadly classified into two categories – Node embedding approaches aim to encode structural information pertaining to a node to produce its low-dimensional representation (Cao et al., 2015; Grover & Leskovec, 2016; Perozzi et al., 2014; Tang et al., 2015; Wang et al., 2016a; 2017; Xu et al., 2017). As they learn each individual node's representation, they are inherently transductive. Recently, (Hamilton et al., 2017a) proposed GraphSage, an inductive method for learning functions to compute node representations that can be generalized to unseen nodes. Sub-graph embedding techniques learn to encode higher order graph structures into low dimensional vector representations (Scarselli et al., 2009; Li et al., 2016; Dai et al., 2016). Further, various approaches to use convolutional neural networks (Kipf & Welling, 2017; 2016; Bruna et al., 2014) over graphs have been proposed to capture sophisticated feature information but are generally less scalable. Most of these approaches only work with static graphs or can model evolving graphs without temporal information.

Dynamic Embedding Approaches. Preliminary approaches in dynamic representation learning have considered discrete time approach. (Zhu et al., 2016) propose a temporal latent space model for link prediction using nonnegative matrix factorization. (Goyal et al., 2017) uses a warm start method to train across snapshots and employs a heuristic approach to learn stable embeddings over time but do not model time. (Zhou et al., 2018) focuses on specific structure of triad to model how close triads are formed from open triads in dynamic networks. (Seo et al., 2016) proposes a deep architecture based on combination of CNN to capture spatial characteristics and an RNN to capture temporal characteristics, to model structured sequences which in graph case will lead to discrete time model. (Du et al., 2015) develops extends skip-gram based approaches for network embedding to dynamic setting where the graphs a re observed as discrete time snapshot and the goal is to learn embeddings that can preserve the optimality of skip-gram objective. NetWalk (Yu et al., 2018) is a discrete-time dynamic embedding approach specifically designed for anomaly detection which uses clique based embedding techniques to learn vertex representations. Recently, (Trivedi et al., 2017) proposed Know-Evolve, a deep recurrent architecture to model multi-relational timestamped edges that addresses the communication process. Unlike our approach, Know-Evolve models all edges at a single timescale, works for setting restricted to relational graphs and uses only edge-level structural information with no attention mechanism. DANE (Li et al., 2017) proposes a network embedding method in dynamic environment but their dynamics consists of change in node's attributes over time and their current work can be considered orthogonal to our approach. (Zuo et al., 2018) proposes a dynamic network formation model to learn node representations by employing a Hawkes process to model the temporal evolution of neighborhood for nodes. This work only considers association events. (Ngyuyen et al., 2018) proposes a continuous time embedding framework that employs a temporal version of traditional random walks in a simple manner to capture temporally evolving neighborhood information.

Other models for dynamic networks. There exists a rich body of literature on temporal modeling of dynamic networks (Kim et al., 2017) that focus on link prediction tasks but their goal is orthogonal to us as they build task specific methods and do not focus on representation learning. Further, there are several approaches in graph mining and temporal relational learning community (Loglisci & Malerba, 2017; Loglisci et al., 2015; Esteban et al., 2016; Jiang et al., 2016) that consider dynamic networks but are orthogonal to our current work. Research on learning dynamic embeddings has also progressed in linguistic community where the aim is to learn temporally evolving word embeddings (Bamler & Mandt, 2017; Rudolph & Blei, 2018). (Yang et al., 2017; Sarkar et al., 2007) include some other approaches that propose model of learning dynamic embeddings in graph data but none of these models consider time at finer level and do not capture both topological evolution and interactions.

Deep Temporal Point Process Models. Recently, (Du et al., 2016) has shown that fixed parametric form of point processes lead into the model misspecification issues ultimately affecting performance on real world datasets. (Du et al., 2016) therefore propose a data driven alternative to instead learn the conditional intensity function from observed events and thereby increase its flexibility. Following that work, there have been increased attraction in topic of learning conditional intensity function using deep learning(Mei & Eisner, 2017) and also intensity free approach using GANS (Xiao et al., 2017) for learning with deep generative temporal point process models.

E   IMPLEMENTATION DETAILS

E.1   ADDITIONAL DATASET DETAILS

Table 2: Dataset Statistics for Social Evolution and Github.

Dataset	#Nodes	#Initial Associations	#Final Associations	#Communications	Clustering Coefficient
Social Evolution	100	407	809	2020554	0.548
Github	12328	70640	166565	604954	0.087

For the social evolution dataset, we consider Proximity, Calls and SMS records between users as communication events (k=1) and all Close Friendship records as association events (k=0). For Github dataset, we consider Star/Watch records as communication events (k=1) and Follow records as association events (k=0). The Social Evolution data is collected from Jan 2008 to to July 2009. We consider the association events between user from Jan 2008-Sep 2008 to form the initial network and use the rest of data for our experiments. We collected Github data from Jan 2013 - Dec 2013. For the nodes in 2013, we consider Follow link that existed between them before 2013 to form the initial network. We pre-process both datasets to contain duplicate (not recurrent in time) records and self-loops. We also process Github dataset to only contain users (and not organizations) as nodes and we select nodes that have at least 50 events.

Temporal Train/Test Split: For all the experiments, the data is divided into train and test based on time line. For Social Evolution Dataset, we train on data from Sep 2008 to May 2009 and use the rest of data for test which gives 10 days of time per test slot. This leads to an approximate 70/30 (train/test) split. For Github data, we train from Sep 2013 to Dec 2013 and test set has 15 days of time per slot. This leads to an approximate 65/35 (train/test) split.

E.2   TRAINING CONFIGURATIONS

We performed hyper parameter search for best performance for our method and all the baselines and used the following hyper-parameters to obtain the reported results:
– For social dataset: Num nodes = 100, Num Dynamics = 2, bptt (sequence length) = 200, embed_size = 32, hidden_unit_size = 32, nsamples (for survival) = 5, gradient_clip = 100 and no dropout.
– For github dataset: Num nodes = 12328, Num Dynamics = 2, bptt (sequence length) = 300, embed_size = 256, hidden_unit_size = 256, nsamples (for survival) = 5, gradient_clip = 100 and no dropout.

For baselines, we used the implementations provided by their authors and we report the range of configurations used for baseline here: $max\_iter = \{1000, 5000, 10000\}, bptt = \{100, 200, 300\}, lr = \{0.0005, 0.0050.5, 0.1, 1\}, embed\_E = \{32, 64, 128, 256\}, embed\_R = \{32, 64, 128, 256\}, hidden = \{32, 64, 128, 256\}, warm = 0, t\_scale = 0.0001, w\_scale = 0.1, num\_epochs = \{10, 50, 100, 500, 1000\}$. As mentioned in experiment section, we always train baselines with warmstart in a sliding window training fashion.

Know-Evolve: The code provided by the authors was implemented in C++.
GraphSage: The code was implemented in Tensorflow by the authors. We use only the unsupervised train module to generate embeddings.
Node2Vec: We use the original python code with few changes in the hyper-parameters. We fix q in the node2vec as 0.8 for Social Dataset and 1 for Github dataset.
DynGEM: We experiment on the original code implemented in Keras with Theano backend by the authors.
DynTrd: We use original code provided by the authors.

For tSNE embedding visualization in Figure 4, we used $sklearn.manifold.TSNE$ library to plot this figure with n components = 2, learning rate = 200, perplexity = 30, metric = "euclidean", min_grad_norm = 1e-9, early exaggeration = 4 and ran for 40,000 iterations.

F MONTE CARLO ESTIMATION FOR SURVIVAL TERM IN $\mathcal{L}$ FOR SECTION 4

Algorithm 2 Computation of integral term in $\mathcal{L}$ for a mini-batch

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Input: Minibatch $\mathcal{M} = \{m_q = (u, v, t, l, k)_q\}_{q=1}^{|\mathcal{M}|}$. Minibatch node list $\mathbf{l}$, sample size N.
Output: Minibatch survival loss L_{surv}

$L_{surv} = 0.0$
for $q = 0 \; to \; |\mathcal{M}| - 1$ **do**
    $t_{curr} = m_q \rightarrow t; \; u_{curr} = m_q \rightarrow u$
    $v_{curr} = m_q \rightarrow v \; ; \; u_{surv} = 0 \; ; \; v_{surv} = 0$
    for N samples **do**
        select $u_{other} \in \mathbf{l}$ uniformly randomly s.t. $u_{other} \notin \{u_{curr}, v_{curr}\}$
        select $v_{other} \in \mathbf{l}$ uniformly randomly s.t. $v_{other} \notin \{u_{curr}, v_{curr}\}$
        for $k \in \{0, 1\}$ **do**
            $u_{surv} + = \lambda_k^{u_{curr}, v_{other}}(t_{curr})$
            $v_{surv} + = \lambda_k^{u_{other}, v_{curr}}(t_{curr})$
        end for
    end for
    $L_{surv} + = (u_{surv} + v_{surv})/N$
end for
return L_{surv}

Algorithm 2 is a simple variant of Monte Carlo trick to compute the survival term of log-likelihood equation. Specifically, in each mini-batch, we sample non-events instead of considering all pairs of non-events (which can be millions). Let m be the mini-batch size and N be the number of samples. The complexity of Algorithm 2 will then be $\mathcal{O}(2mkN)$ for the batch where the factor of 2 accounts for the update happening for two nodes per event which demonstrates linear scalability in number of events.