

Embeddings for Signed Weighted, and Temporal Networks

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1 Introduction

Networks provide a rich representation of structured data. From social networks and web-graphs, to protein-protein interaction networks, such data is abundant in our world. A lot of work has been done on how to understand such data on the level of well-studied network features, such as triadic closure, network motifs, and node centrality. Recently, with the great success of machine learning in other domains, embedding network data into the low dimensional vector spaces from which such algorithms often expect their inputs has become an active area of research. In particular, one of the promises of modern embedding algorithms is to be able to move away from hand-crafted features to ones learned automatically, in an unsupervised or semi-supervised manner ([4], [9] [3], [6]).

Such embedding algorithms, however often disregard two potential characteristics of networks occurring in the real world: that networks are *temporal* (also called *dynamic*), meaning that their structures change over time, and that there are naturally occurring *signed weighted* networks, that is, ones whose edges have signed and possibly fractional weights. In this project, I explore ideas random embeddings on such networks: in particular, using temporal random walks, and using what I call “relational weighting” as a modification for embeddings in signed weighted networks. In fact, to the best of my knowledge, embedding methods have as of yet to be developed for signed weighted networks. The quality of the embeddings is to be evaluated on edge-weight prediction tasks. I find that while embedding method my proposed extension of random walk embedding to signed weighted networks isn’t tractable without further improvement, while using temporal random walks improves the performance of regular embedding techniques. The code for this project is available [here](#).

2 Related Work

2.1 Signed edge weight prediction

One of the few current works on signed weighted networks is that of Kumar, et al. [7]. In it, the authors try to predict the ratings users give each other on various platforms. This takes the form of edge-weight prediction in networks, in which the nodes represent

users. Since the ratings can range from negative to positive, the networks are in fact signed and weighted. The insight of this paper is to formulate two properties that seem *a priori* important for networks in which users rate one another — *goodness*, and *fairness* — as well as state axioms that such properties ought to satisfy. (Intuitively, goodness is whether a node is highly rated by other nodes, while fairness is the property measuring how “fairly” a node rates other nodes, that is, whether a node gives other nodes ratings close to their goodnesses.) The authors then give natural recursive definitions of these properties, and show that they fulfill the desired axioms. They then evaluate these definitions of the task of edge-weight prediction, where for two nodes u and v , their predicted edge weight is $f(u)g(v)$, where f is fairness and g is goodness. Comparing in various testing regimes against, previous techniques used for predicting unsigned weights, or just plain sign prediction, the authors find that their theoretically sound approach outperforms the others.

Inspired by this exploration of signed weighted networks, in this project we ask, “can embeddings do better?” In particular, the prior work above: (a) cannot (readily, at least) predict edge existence, (b) is not concerned with dynamic networks, and as such may not be efficiently updateable in an online setting, and lastly (c) assumes that semantically, edges and their weights correspond to agents rating other agents (although, to be fair, this may be the most common type of weighted signed network).

2.2 Embeddings

2.2.1 Random Walk Embedding

One kind of embedding that has recently found success is *random walk embedding*. The now-classical example of this is DeepWalk [9]. This algorithm is inspired by the Skip-Gram model from NLP, in which a model is trained to maximize the probability of seeing the “context” of a word (i.e. a certain window of words occurring together with the word of interest) in a corpus given the word itself. The model is thus trained to learn useful representations of each word w as a vector $f(w)$, thus creating a word-embedding. The paradigm generalizes readily to graphs, as follows. We let nodes be our “words,” and random walks be our “contexts.” Thus, the algorithm is to repeatedly perform random walks v_1, \dots, v_ℓ , and update our model parameters (node representations) such that for each node v in the walk, the probability of encountering the other nodes given that node is maximized. Mathematically, we are maximizing the expression

$$P(\{v_1, \dots, v_{i-1}, v_{i+1}, \dots, v_\ell\} | v_i).$$

The node2vec [3] embedding strategy is algorithmically a strict super-set of DeepWalk, with two tunable parameters governing how the random walks used during training are performed. In particular, these parameters tune the trade-off between how often the walk returns along the edge it just followed, and thus stays close to the source node (parameter p), and how often the walk moves to a node that has a greater distance from the source than the current node (parameter q). The authors here also elaborate two assumptions informing how the probabilities used by the Skip-Gram model are calculated. First is a conditional independence assumption made, meaning that

$$P(\{v_1, \dots, v_{i-1}, v_i, \dots, v_\ell\} | v_i) = P(v_1 | v_i) \cdots P(v_{i-1} | v_i) P(v_{i+1} | v_i) \cdots P(v_\ell | v_i).$$

Second is the “symmetry assumption” that nodes have a “symmetric effect” on each other in feature space, and that thus the conditional probabilities are modeled as a softmax of the dot-products of node embeddings:

$$P(w|v) = \frac{\exp(f(w) \cdot f(v))}{\sum_{w' \in V} \exp(f(w') \cdot f(v))}.$$

With this formulation of the objective function, the embeddings are trained, and then used as inputs to logistic regression for classification of node labels. The authors also suggest extending the framework to allow for link prediction, by using binary operators between node embedding pairs to create “edge embeddings.” The resulting embeddings outperformed state-of-the-art methods on both tasks. Additionally, random walk embeddings have the benefit of being scalable, and embeddings (for nodes *and edges*) are efficiently updateable, by training on new random walks.

Note however that the symmetry assumption above, and subsequent formulation of the conditional probability, are not directly extendable to the signed/weighted network case! A small amount work has in fact been published addressing signed edges in embeddings — see SNE [13] and SiNE [12], of which only the former uses random walk embedding — but these techniques restrict themselves to edge weights of ± 1 . They also turn out to not be naturally extendable to the fractional-weight case (for more, see the Appendix). In addition none of these embedding techniques take into account the temporal nature of networks.

2.2.2 Temporal Embeddings

For embeddings taking into account the temporal nature of edges, I will focus on the “Continuous Time Dynamic Network Embeddings” of Nguyen et al. [5]. The authors present a general framework for extending random walk embeddings to take into account edge arrival times. The intuitive step taken here is to sample temporal random walks rather than regular random walks on some snapshot of the graph. Temporal random walks are walks that respect the time-order of edges. The authors make the contribution of formulating how to sample a start edge/time, as well as how to sample the “temporal neighbors” of each node in the walk. They note that the distributions used to do so may be biased to prefer more recent edges, if such behavior is desirable.

Sampling temporal walks as “contexts,” the authors generalize the Skip-Gram embedding model to train network embeddings. (In this sense, they are applying their temporal framework to node2vec.) Comparing to DeepWalk, node2vec, and other embeddings, they find that temporal embedding outperforms on classification and link prediction (again, formulated as logistic regression, as in node2vec). The authors also note that their approach has the advantage over others of not having to deal with discrete graph “snapshots,” which introduces the problem of having to choose a time granularity for these snapshots. In fact, this, in addition to the high barrier to adaptation for the purposes of this project, is the reason for which we will prefer this approach to those of STWalk [8] and DynGEM [2], although each is interesting in its own right — for example, the latter deals with embedding stability over time.

3 Problem Definition

We will now make precise our problem setting. We are given some graph $G = (V, E)$, where for each edge $e = (u, v) \in E$, we have some weight $\mathcal{W}(e) \in [-1, 1]$. It is important that the edge weight can be any real value, positive or not. It is less important that these weights are bounded, though we should note that this is an assumption being made here. To simplify our problem setting, we scale the edges to ensure that each $[-1, 1]$, and so if a new edge occurs after we first create our embeddings such that its scaled weight is not in $[-1, 1]$, the assumptions being made here may break down. (It is a different question how much this impacts actual performance of our suggested algorithms, which we will not explore here.) We will also assume our graphs are temporal. That is, for each $e \in E$ we have a time-stamp $t_e \in [0, \infty)$, which denotes the time edge e appeared. Furthermore, we will assume that once an edge appears, it does not disappear or change its weight. We will thus notate a temporal, signed weighted network as $G = (V, E, \mathcal{W}, t)$.

Formally, our problem is that of creating a *node embedding* in a temporal signed network $G = (V, E, \mathcal{W}, t)$. For our purposes, a node embedding is a function mapping each node to a representation in d dimensional space $f : V \rightarrow \mathbb{R}^d$, for some chosen constant $d \in \mathbb{N}$. Since we will evaluate our embeddings by prediction of edge-weights, we also need to generate *edge embeddings*. To do so, we use the method suggested in the node2vec paper [3], and let our edge embedding function $g : E \rightarrow \mathbb{R}^d$ be $g((u, v)) = f(u) \odot f(v)$, where \odot denotes the Hadamard (or component-wise) product of two vectors. (In the node2vec paper [3], this is found to be the best performing of several explored ways of combining node embeddings to create edge embeddings. In fact, I believe this is reasonable because the skip-gram objective in essence optimizes similarity as calculated by the sum of the elements of two Hadamard’ed vectors, i.e. their dot-product.)

4 Proposed Embedding Methods

I propose a method in order to take into account signed weighted edges: a modification of the objective function for node2vec, for which I think “relational weighting” is a reasonable name. I will also see if only using time-respecting random walks, as in Nguyen et al. [5], improves performance.

4.1 Time-Respecting Random Walks

A time-respecting (or temporal) random walk is a sequence of edges e_1, \dots, e_k such that $t_{e_1} \leq \dots \leq t_{e_k}$. In order to generate such a walk from a given starting node u_1 , I first sample a node v_2 uniformly from the neighbors of u_1 to get a starting time $t_{(u_1, v_2)}$ for the walk. I then repeatedly sample uniformly neighbors u_{i+1} of the current ending node u_i of the walk, such that $t_{(u_i, u_{i+1})} \geq t_{(u_{i-1}, u_i)}$. Such random walks then form the training corpus for the embedding models.

4.2 Relational Weighting: Accounting for Signed and Weighted Edges in Skip-Gram Models

We now turn to the question of adapting the random walk models above to take into account the fact that edges are signed and weighted (non-integer). Taking a queue from the SNE model of [13], we will try to incorporate information about “where the walk came from/is going to,” in terms of edge weights, when evaluating the similarity between nodes. Recall that `node2vec` defines the conditional probability of seeing node w given node v as

$$P(w|v) = \frac{\exp(f(w) \cdot f(v))}{\sum_{w' \in V} \exp(f(w') \cdot f(v))}.$$

We note that the $f(w) \cdot f(v)$ term in the numerator is in a sense measuring the similarity between nodes w and v . That is, we expect $P(w|v)$ to be large if w and v are similar because we expect our embeddings to be such that $f(w) \cdot f(v)$ is large. (The normalization term in the denominator prevents us from achieving this by just increasing the size of $f(v)$.) We can formulate this by saying that we want the cosine distance between the vectors to be 1

$$\frac{f(w)}{\|f(w)\|} \cdot \frac{f(v)}{\|f(v)\|} = 1.$$

Now suppose that w and v are “dissimilar,” in the sense that they have a negative weight edge between them, $\mathcal{W}(w, v)$. Then, in the sense that the $f(w) \cdot f(v)$ term is measuring similarity, we would expect $f(w) \cdot f(v)$ to be small, or more precisely, be negative. Let us formulate this idea as

$$\frac{f(w)}{\|f(w)\|} \cdot \frac{f(v)}{\|f(v)\|} = \mathcal{W}(w, v).$$

In fact, note that this formulation works for positive edge case as well. In either case, let us re-write this as

$$\frac{1}{\mathcal{W}(w, v)} \frac{f(w)}{\|f(w)\|} \cdot \frac{f(v)}{\|f(v)\|} = 1.$$

By analogy to the initial Skip-Gram probability we are maximizing, we can see that our desired embeddings correspond to a maximization, now of

$$\frac{f(w) \cdot f(v)}{\mathcal{W}(w, v)},$$

rather than of $f(w) \cdot f(v)$.

The first intuitive thrust of this proposal is then to modify our conditional “probabilities” (they are not truly probabilities anymore) we are trying to maximize for each random walk. The next question is how do we determine the equivalent of $\mathcal{W}(w, v)$ if w was not a neighbor of v in the walk. The intuition here is the following. Suppose that there is only one node u_1 between w and v in the random walk. If w “dislikes” u_1 (the edge between them is negative), and u_1 similarly “dislikes” v , then it is likely that w “likes” v . Otherwise stated, this is the principle of transitivity, in the sense of “the enemy of my enemy is my friend.”

Thus, proposal here is to use the product of weights along the path between w and v to determine the *sign* of $\mathcal{W}(w, v)$. However, using the product directly would quickly lead to

minute values of $\mathcal{W}(w, v)$, so the magnitude is instead chosen to be of average \mathcal{W} magnitude along the path. So, if the random walk had node u_1, \dots, u_m occur between w and v in the walk, we now use the term

$$c_{w,v} = \left(\text{sign}(\mathcal{W}(w, u_1)) \prod_{i=1}^m \text{sign}(\mathcal{W}(u_i, v)) \right) \frac{|\mathcal{W}(w, u_1)| + \sum_{i=1}^m |\mathcal{W}(u_i, v)|}{m+1},$$

such that now our numerator for the “probability” is

$$\frac{f(w) \cdot f(v)}{c_{w,v}}.$$

The modeling assumption of transitivity here is indeed strong, and so I explore introducing an attenuation factor $\beta \in (0, 1]$, accounting for the length of the path between w and v , such that

$$c_{w,v} = \left(\text{sign}(\mathcal{W}(w, u_1)) \prod_{i=1}^m \beta \text{sign}(\mathcal{W}(u_i, v)) \right) \frac{|\mathcal{W}(w, u_1)| + \sum_{i=1}^m |\mathcal{W}(u_i, v)|}{m+1}.$$

There is one final question to answer, which is how to modify the partition function, i.e. the denominator of the conditional probability above. The natural approach here is intractable, as it would involve calculating the $c_{w',v}$ for any pair of nodes w', v , possibly by finding the shortest path between them. I will explore using negative sampling — as in the implementation of node2vec [3]. Here, “negative” means we sample edges between nodes that are not likely to have an edge between them, and sample the weight according to the weight distribution of the overall graph. In practice, this means that we simply pick nodes w' unlikely to co-appear on a random walk with v , and let $c_{w',v} = \epsilon$, where $\epsilon \neq 0$ is the average weight of edges in the network, with the expected label being 1, rather than 0 as in negative sampling as normally performed in skip-gram models. (Both the sampling of w' and the choosing of $c_{w',v}$ can be modified, and indeed should be explored further, especially with regards to how the assumptions might need to change given different network properties.) Note that this means that the denominator of the “probability” term needn’t change, since we have

$$P(w|v) = \frac{\exp\left(\frac{f(w) \cdot f(v)}{c_{w,v}}\right)}{\sum_{w' \in V} \exp\left(\frac{f(w') \cdot f(v)}{\epsilon}\right)}.$$

Note that $\epsilon \neq 0$ by assumption (and almost always in practice), and that if $c_{w,v} = 0$, we simply skip the training pair.

5 Data

As in Kumar et al. [7], the main datasets of interest will be [Bitcoin OTC](#) and [Bitcoin Alpha](#), both of which are Bitcoin exchanges in which users can provide ratings for other users (which we normalize to be in $[-1, 1]$). We will take some space here to understand some basic properties about the networks, in order to provide context for the experimental

	OTC	Alpha
Nodes	5881	3783
Edges	35592	24186
Isolates	0	0
Nodes in Largest SCC	4709	3235
Nodes in Largest WCC	5875	3775
Ave. Clustering Coef.	0.1775	0.1776

Table 1: Structural statistics for Bitcoin OTC and Alpha exchange networks (SCCs meaning strongly connected components, and WCCs meaning weakly connected components)

results below. First, basic statistics on the network structures (Table 1) reveal that these are relatively small, and have atypically small average clustering coefficient for social networks (compared, for example, with the social network data used in [11] from Facebook, Twitter, and Google+). This is consistent with the general observation that these networks are not, in fact, “social networks” in the usual sense of “friendship” or “following,” and are instead more transactional in nature. The networks do, however, exhibit the typical structure of having a giant component.

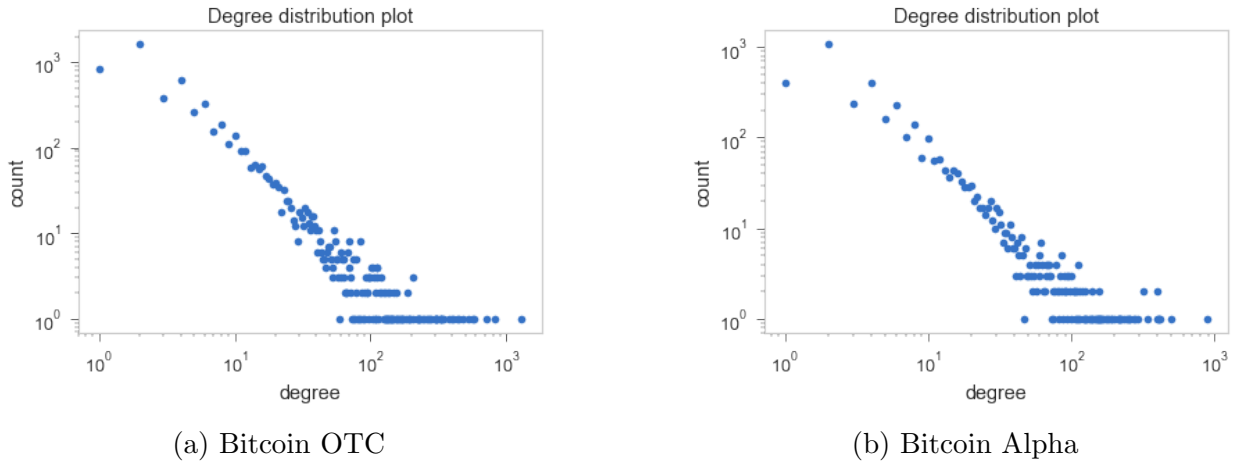


Figure 1: Unweighted degree distributions

We further explore the nature of the networks’ connective structures by examining their degrees, both in the weighted and unweighted sense. The unweighted degree distributions (Figure 1) for both networks follow the “power law” (are generally linear on logarithmic scales), although it is interesting that both exhibit almost a kind of bi-linearity, with what looks like two trend lines in their distributions. Similarly, the weighted degree distributions (Figure 2) look almost identical, with upper bounds that decay in both positive and negative directions as power laws from average weights of slightly above 0. This structure in fact suggests (not necessarily surprisingly) that some sort of aggregation of each node’s incoming edge weights, in the same spirit as in Kumar et al. [7], differentiates nodes well. The fact that ratings are “slightly” skewed in the positive direction is reflected in the fact that edges

are 89% positive for Bitcoin OTC and 93% positive for Bitcoin Alpha.

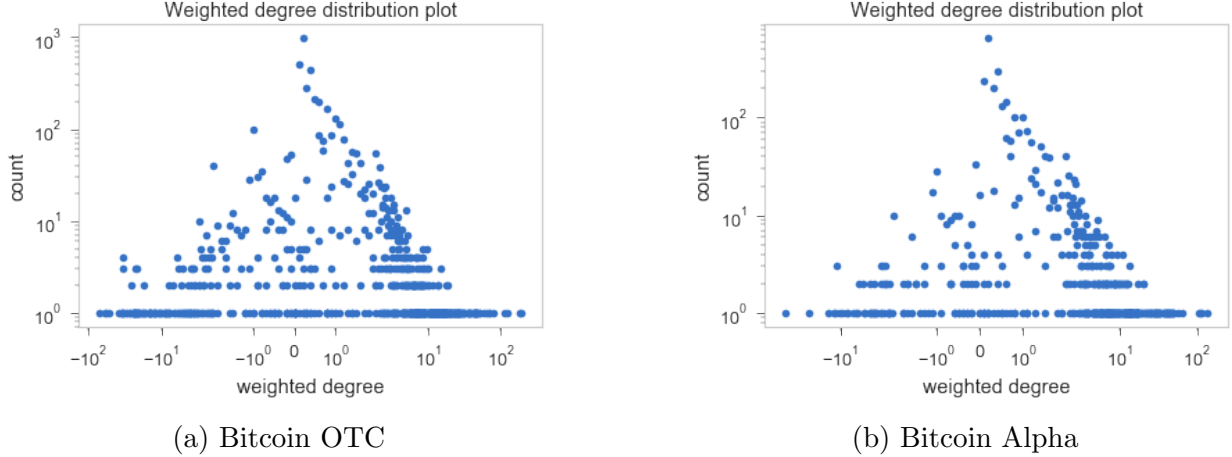


Figure 2: Weighted degree distributions

6 Experiments

6.1 Baselines and Setup

While embeddings can be used for edge-existence prediction, the Fairness Goodness algorithm can be evaluated directly on the edge weight prediction task, and cannot be evaluated on the edge-existence prediction task. We will thus restrict ourselves to the former setting for evaluating our embedding methods. In edge-weight prediction, the embeddings are used as inputs to Support Vector Regression.

We begin by comparing all baselines on the edge-weight prediction task, using the Root Mean Square Error on random test set of the edges as our metric, in order to check against the fairness-goodness algorithm performance in [7], which we do manage to reproduce. For each model, we held out 10% of the data as the validation set. Note though this data was completely unused by Fairness Goodness and DeepWalk, which do not have tunable parameters. However, this hold out was maintained to ensure comparability with the results of node2vec, in which the validation set was used to pick the parameters p, q for each of Bitcoin OTC ($p = 0.25, q = 1$) and Bitcoin Alpha ($p = 0.25, q = 4$). Chosen similarly for SNE, the hyperparameters were the number of walks per node (10), maximum walk length (80), and language model context size (10). We then removed increasing portions of the training set and added them each to the test set, in order to test how the models perform under different data sparsity conditions. In this manner, the test set size begins at 10% and goes up to 70% of the total data.

The results are summarized in Table 2. In all but one setting, the Fairness Goodness algorithm outperforms the embedding baselines. This shouldn’t be surprising, as it’s the only algorithm taking into account edge weights. Note that the simplest embedding approach, DeepWalk performs slightly worse than the Fairness Goodness algorithm, but not by much. This suggests that embedding is indeed a viable approach to signed edge-weight prediction.

Indeed, since DeepWalk does not take edge weights in to account at all, it seems very likely that a lot of information about weights is encoded purely in the topologies of the networks. A second encouraging observation is that the embeddings, and especially node2vec, seem even more resilient to sparse “training” data than the Fairness Goodness algorithm. In fact, resilience to sparse data was one of the aspects of the Fairness Goodness algorithm the authors of [7] advertise.

	OTC			Alpha		
train-test split	80-10	50-40	20-70	80-10	50-40	20-70
FG	0.313	0.321	0.336	0.263	0.275	0.291
DeepWalk	0.332	0.323	0.358	0.288	0.291	0.290
node2vec	0.335	0.330	0.345	0.288	0.288	0.288
SNE	0.336	0.360	0.407	0.313	0.327	0.335
DeepWalk-T	0.321	0.335	0.342	0.273	0.285	0.292
node2vec-T	0.320	0.331	0.336	0.272	0.284	0.289
SNE-T	0.324	0.338	0.350	0.280	0.287	0.290
DeepWalk-RW	0.360	0.365	0.368	0.289	0.296	0.296
DeepWalk-T-RW	0.369	0.366	0.370	0.277	0.285	0.290

Table 2: Performance (RMS error) of models on edge-weight prediction task, grouped as: baselines, embeddings using temporal walks (suffixes with “-T”), embeddings using relational weighting (suffixes with “-RW”) and embeddings using both (suffixes with “-T-RW”)

Among the embeddings, node2vec better than SNE. However, I do not believe any conclusive claim can be made here, since in the SNE paper [13], default hyperparameters are used for node2vec and they find the opposite is true (SNE generally outperforms node2vec). Nonetheless, with some hyperparameter selection for *both* models, I found that 1) SNE performed better when the random walk parameters were set to those used for node2vec (as opposed to those in the paper), and 2) it underperformed.

6.2 Temporal Random Walk Embeddings

Turning to the results of embeddings using temporal walks, we find that in general, there is up to a 10% improvement in root-mean-squared error over the same embeddings using general (non-temporal) random walks. Note however, that the advantage of using temporal walks vanishes when the proportion of training data is decreased. This makes sense, since restricting to only temporal walks (which are a strict subset of all random walks on our graphs) makes the actual trainable sequences of nodes fairly sparse.

Nonetheless, using temporal walks, node2vec manages to perform on-par with Fairness Goodness when both are trained on only 20% of the data, for the OTC dataset. (The hyperparameters for Bitcoin OTC were $p = 4$ and $q = 2$, for and Bitcoin Alpha they were $p = 2$ and $q = 1$.)

6.3 Relational weighting

For the purposed of relational weighting, only DeepWalk was used. Implementing relational weighting involved directly changing the objective optimized in the Gensim [10] implementation of word2vec. This was substantially easier to do in the pure Python routines rather than the highly optimized Gensim Cython code. Thus, relational weighting took considerably longer to train. To account for this fact, only 2 negative samples were drawn for each training pair, and each model was trained for only 3 epochs. It is also worth noting that the negative sampling led to numeric instability, in particular, the exploding of the skip-gram loss term. I will also readily admit to not having 100% certainty that I did not make an error in calculating the change to the backpropagation terms under my modified objective function (although I cannot see how, but please feel free to look at the project code and let me know).

Having noted the above, it is not surprising that no improvement was seen at all from using relational weighting. In fact, I believe it *is* surprising that these models performed as well as they did.

7 Conclusion

In this project, I find that traditional random walk embedding methods come close to the performance of the Fairness Goodness algorithm on edge-weight prediction, while my proposed method of relational weighting is likely numerically unstable. I also confirm the hypothesis that using temporal random walks improves random walk embeddings, but only when there are in fact enough valid temporal walks for the data to not be too sparse. A continuation of this work would, in my opinion, have to include two major points: firstly, and most obviously, more work on making relational weighting tractable, and secondly, on evaluating the embedding methods above on *edge-existence* prediction (as opposed to edge-weight).

In general, I believe an exploration of embedding where we recognize the independent nature of edge weight and edge existence probability is needed. In a sense, these are orthogonal notions of distance. I’ve had the thought, for example, of exploiting embedding vectors’ cosine distance for one, and euclidean distance for the other. I have a sense that in fact, the general approach taken in a recent paper by Chen et al. [1] (in fact, more recent than this project), is the right one. In particular, they take a multi-task learning-like approach, in which a “structural” loss (read: skip-gram) and “relational” (read: neural-network for predicting edge attribute) loss are trained on simultaneously in a random walk embedding procedure. This approach, however is slightly less principled than I would like, since I have a feeling “meaningful” embeddings for signed weighted networks should be possible if we create models with a semantic understanding of edge-weight. This seems to me fruitful ground for embedding weighted signed networks.

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8 Appendix

8.1 Signed network embeddings

One the recent works on signed network embeddings is [13], who call their model SNE. The important note here is that these networks have edge weights of ± 1 . The authors here also perform random-walk embedding, but rather than adopting the Skip-Gram metaphor, elect to predict each node v from the context of previously seen nodes u_1, \dots, u_ℓ in the walk. In order to account for edge sign, the authors adapt the log-bilinear model, predicting the node embedding of v as

$$\hat{f}(v) = \sum_{i=1}^{\ell} c_i \odot g(u_i).$$

We will unpack this equation here. First, c_i is one of two trainable vectors, c_- or c_+ , depending on whether the edge out of u_i in the walk is negative or positive, respectively. This is how the authors take edge sign into account. Second, there are two node embedding functions, the “target embedding” f , and the “source embedding” g . The actual final embedding of any node v , to be used for example in classification or link prediction (by the same method as explained in node2vec), is the concatenation of the two embeddings $[f(v); g(v)]$. The authors claim that having both embeddings is important, as only using one of them leads to lower accuracy on some tasks, although here one can wonder: why not just use a single embedding function from the start? I should note here that it is actually not 100% clear to me from the paper if I’ve interpreted these the use of two embeddings correctly, although I consider the above by far the most reasonable interpretation.

The authors use the trained embeddings to perform classification and link prediction, both with the same paradigm as used in node2vec — with link prediction now having 3 possible classes (-1, 0, 1) — and find that their model outperforms node2vec. This is intriguing (if not surprising) because it suggest to us how we might start thinking about incorporate edge weights and/or signs into a model. Unfortunately, the approach here cannot be readily generalized to the scalar, rather than integral, signed edge case, since we cannot have a c_i vector for each possible edge weight. Another point is that, once again, this model doesn’t take into account the temporal nature of edges. However, vanilla SNE is clearly a baseline to compare against in the project. (It is also interesting to note that when performing parameter sensitivity analysis, the authors find that the best performance on link prediction is achieved when random walks are of length 1. This seems generally surprising, since then the model only captures information about neighbors. This may have something to do with how this model eschews the Skip-Gram models’ sense of “closeness” discussed for node2vec and DeepWalk.)

The other work which I will briefly mention is [12], who propose a model called SiNE for signed network embeddings. Here, the authors develop an objective function based on structural balance theory, the rough intuition being that in the embedding, “friends” should be closer than “enemies” (with friendship being defined in the edge sign sense). A problem arises however when a node has friends but no enemies in their 2-hop network (or vice-versa, although this is less common), since the objective function doesn’t contain such nodes, which makes it impossible to learn embeddings for them. The authors devise a strategy of adding a

“virtual node,” and then making this node the enemy of each node that doesn’t have enemies in its 2-hop network. This approach has the advantage of strong theoretical motivation from social theory, but isn’t suitable for our purposes since it doesn’t take into account either edge weight or even direction. We can also note that the virtual node artificially changes the network topology, which one might imagine could possibly lead to strange effects.