SPACED: A Novel Deep Learning Method for Community Detection in Social Networks

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Abstract: Community detection is a landmark problem in social network analysis. To address this challenge, we propose SPACED: Spaced Positional Autoencoder for Community Embedding Detection, a deep learning-based approach designed to effectively tackle the complexities of community detection in social networks. SPACED generates neighborhood-aware embeddings of network nodes using an autoencoder architecture. These embeddings are then refined through a mixed learning strategy with generated community centers, making them more community-aware. This approach helps unravel network communities through an appropriate clustering strategy. Experimental evaluations across synthetic and real-world networks, as well as comparisons with state-of-the-art methods, demonstrate the high competitiveness and often superiority of SPACED for community detection while maintaining reasonable time complexities.

1 INTRODUCTION

Social networks, connecting vast numbers of individuals and entities, have become pivotal in shaping communication, information dissemination, and social interactions, making the study of their structures and dynamics increasingly crucial. One of the most important social structures that characterizes these networks is their community structure. Although a formal definition of a community in network analysis does not exist, a widely accepted definition describes a community as a subset of social agents, typically represented as graph nodes, between which interactions, typically represented as graph edges, occur more "densely" than with the rest of the net-

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work. This dense intra-community connectivity indicates that members within the community are more closely related or interact more frequently with each other than with nodes outside the community. As a result, such clusters naturally form identifiable groups, known as communities, within the network structure (Figure 1). It immediately appears from this definition that discovering communities inside a network, which is commonly referred to as the community detection problem, has important applications such as targeted advertising, functional group identification, or even terrorist threat prevention (Karataş and Şahin, 2018).

The last decades have witnessed an increase in the use of machine learning (ML) and deep learning (DL) based methods to tackle this problem due to their ability to handle efficiently high dimensional data spaces such as graphs and uncover intricate patterns on them such as community structures. One of the mainstream approaches in the class of DL methods is to embed the graph nodes into a vectorial space in such a way as to reflect their community proximities, i.e. nodes more likely to be in the same community according to the

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Figure 1: An undirected graph representation of a social network used for community detection.

network data are associated with nearby embeddings in the vectorial space, thus transforming the community detection problem on the network into a clustering problem on the vectorial space representation.

Aligning with this class of methods, in this paper, we propose SPACED, a model that aims at generating meaningful node embeddings and community centers that uncover the community structure of social networks based solely on their topologies, i.e. the undirected connections between the network agents. It performs this by generating initial encodings of the nodes based on their neighbourhoods, embedding them using an auto-encoder architecture and generating initial community centre embeddings which will help improve the node embeddings.

The rest of this paper is organized as follows. Section 2 presents the related works outlining the different sources that SPACED inspires from. Then, section 3 details the proposed solution. Section 4 presents the obtained results on the selected benchmarks and their significance. Lastly, section 5 concludes the paper, summarizing our findings and suggesting directions for future research.

2 RELATED WORKS

Various sequence-based node embedding techniques such as Line (Tang et al., 2015), Node2Vec (Grover and Leskovec, 2016) or DeepWalk (Perozzi et al., 2014) base their construction of node embeddings on the "proximity" of the nodes inferred by their cooccurrence on various sampling methods on the network such as truncated random walks, second-order random walks etc., reflecting on the idea that nodes of the same community are more likely to appear simultaneously in small sub-regions of the network. The embeddings generated by this class of techniques have been applied for community detection with competitive results as shown in (Tandon et al., 2021). Also, the SkipGram model (Mikolov et al., 2013) which inspired some of the methods in this class have in the same way motivated parts of SPACED architecure.

Using auto-encoders in community detection has gained significant attention recently. Reference(Xie et al., 2019) introduced CDDTA, which employs a deep auto-encoder for nonlinear feature extraction in networks and incorporates unsupervised transfer learning to refine representations. Reference (Tian et al., 2014) proposed the GraphEncoder method, which uses graph neural networks alongside auto-encoders to encode structural information into low-dimensional embeddings, preserving topological properties and enhancing clustering accuracy in largescale networks. Reference (Bhatia and Rani, 2019) proposed DeCom, which integrates modularity-based community detection with ensemble clustering, combining multiple clustering results into a consensus solution to enhance robustness and stability in community detection.

For the works that consider performing node embedding in conjunction with community embedding with the goal to create better community aware node embeddings, (Rozemberczki et al., 2019) proposed GEMSEC, which aims to mutually enhance both node and community embedding processes. GEMSEC employs sequence-based node embedding techniques, where nodes occurring closely in random walks are embedded close to each other. The optimization objective combines the negative log-likelihood of observed neighbourhood samples with a clustering cost and is solved using a variant of mini-batch gradient descent. Another line of work in this category are comE (Cavallari et al., 2017) and its improved version comeE+ (Cavallari et al., 2019) which automatically detects the number of communities. They consider that community embeddings should reflect the member nodes distribution in the feature space and thus propose embedding the communities as probability distributions by fitting them to a Gaussian mixture model. The results obtained by comE/comE+ proved the benefit of considering community-aware proximities in the embeddings construction process.



Figure 2: SPACED general architecture.

3 PROPOSED SOLVING APPROACH

A social network can be modelled by a graph G = (V, E) where V is the set of nodes representing individuals and E is the set of edges indicating interactions between individuals. The goal is to find the community structure within this graph which is a partition $C = \{C_1, C_2, ..., C_K\}$ of K communities on the node set V.

As stated in the introduction, our method attempts to solve this problem by finding a node embedding that better represents the community structure. A node embedding is a mapping $f: V \to \mathbb{R}^e$ of nodes to an embedding space of dimension *e*. A successful embedding can then turn the problem into a straightforward point cloud clustering. The SPACED approach follows a four-phase pipeline to achieve this as shown in figure 2:

- **Phase 1. Graph Processing** to generate initial high-dimensional encodings for graph nodes.
- **Phase 2. Node embedding** to obtain an initial low-dimensional node embedding using an autoencoder architecture.
- Phase 3. Community centers/leaders initialization aiming to create embeddings for community centers/leaders, in the same space as the nodes, positioning them at the best possible position inside the point cloud formed by the node embed-

ding.

• Phase 4. Mixed learning where the model simultaneously improves node embedding while bringing them and the community centers/leaders closer to each other, thus, producing a more community-aware node embedding with more condensed clusters eventually (Rozemberczki et al., 2019).

The following sections will thoroughly explore these phases, along with their different variants.

3.1 Graph Processing

The first problem encountered in all community detection methods is how to read the graph and explore its topological structure most optimally, here we tried to use a data structure that can better represent the neighbourhood levels of a node while emphasizing the differences between these levels. So, in this paper, we propose two data structures called the Views Matrix and Walks Matrix.

3.1.1 Views Matrix

To produce this structure, the graph is seen as a Markov chain model, where each node is a state and the probability of moving from a node to one of its neighbours is uniform between all neighbours N(i) (since the graph is unweighted). So, *P* is an $N \times N$

matrix (Eq. 1) representing the probability of moving from node i to node j.

$$P_{ij} = \begin{cases} \frac{1}{|N(i)|} & \text{if } j \in N(i) \\ 0 & \text{otherwise} \end{cases}$$
(1)

So the matrix *P* can represent first-order proximity in a probabilistic format. To obtain other levels of neighbourhood we can easily use the property of Markov chains where P^2 represents the probability of going from node *i* to *j* in two steps and so on. So, for a max depth *d* and for all $k \in [1, d]$ we would obtain:

$$P^{(k)} = P^k \tag{2}$$

But you can notice here that the return probability is always considered, for this we propose another variant where we use this formula instead:

$$P_{ij}^{(k)} = \begin{cases} 0 & \text{if } k = -1 \text{ or } (k = 0 \text{ and } i \neq j) \text{ or } P_{ij}^{(k-2)} \neq 0 \\ 1 & \text{if } k = 0 \text{ and } i = j \\ (P^{(k-1)} \times P)_{ij} & \text{otherwise} \end{cases}$$
(3)

Now, we need to differentiate between the different levels of neighborhood, so we add some weights to each level matrix $P^{(k)}$. We use two approaches: (1) Arithmetic approach where we weighted the levels by subtracting one for each level $\{d - 0, d - 1, ..., d - (d - 1)\}$; (2) Harmonic approach where we divide by an incrementing number each level $\{\frac{d}{1}, \frac{d}{2}, ..., \frac{d}{d}\}$.

Finally these resulting matrices $P^{(k)}$ are concatenated along their columns resulting in a $|V| \times (d * |V|)$ matrix as follows:

$$M = \begin{bmatrix} P^{(1)}w_1 & P^{(2)}w_2 & \cdots & P^{(d)}w_d \end{bmatrix}$$
(4)

3.1.2 The Walks Matrix

In an attempt to imitate random walks without performing them, we exploited the already constructed views matrix by combining them with a weighted mean using the same weights used earlier. This creates for each node i a distribution resembling the portion of a node appearance in different random walks of varying lengths started from the node i, which creates some sort of context around each node.

$$W = \frac{\sum_{k=1}^{d} P^{(k)} w_k}{\sum_{k=1}^{d} w_k}$$
(5)

3.2 Node Embedding

We proposed an autoencoder architecture that tries to lower each level of neighbourhood to the embedding space of size e individually and then learn the relation between them. To achieve this, we created a neural network with 3 hidden layers and an input and output. For the input we use the views matrix, where each row represents a sample (which is a node of the graph), thus, a node is represented by a vector of size d * |V|. The first hidden layer, called the Level embedding layer, lowers each level independently to an embedding of size e so the output of this layer is a vector of size d * e. The second layer combines the embedding of each level to a final embedding, here we densely link each dimension from the different levels to their respective dimension in the final embedding, thus, learning for each dimension a weights vector of size |V| for a total of d * |V| weights for all the dimensions. The last layer is a symmetric layer identical to the Level embedding layer. Followed by an output layer symmetric to that of the input layer. This architecture is demonstrated in figure 3

3.3 Communities Centers / Leaders Initialization

This phase aims to position community centers/leaders as optimally as possible inside the point cloud of nodes. Community centers initialization means putting independent points that sit at the center of the communities. In this category, we propose two methods: A neural network architecture, called Community Embedder, which uses a custom loss function (11) for the above-mentioned purpose, and the KMeans method.

Community leaders initialization on the other hand focuses on choosing a node as the leader of its community. Here we crafted two other methods: Leaders SA which is a simulated annealing for choosing the optimal combination of nodes that minimizes the same loss function used earlier (11), and a method named Walk Leaders that exploits the Walks matrix and uses a heuristic we defined to determine the leaders.

3.3.1 Community Embedder

It's a simple neural network containing only densely connected input and output layers. The input layer receives a one-hot encoded vector denoting the community to embed (so the size of this layer is K) and lowers it to the embedding space represented by the output layer (evidently of size e). To define the loss function for this model we needed to put some theories that describe an optimal position of a community center, we came up with four, each having a loss term and then combined them into one loss function as follows:

First, we naively suppose that a good partitioning



Figure 3: The autoencoder architecture.

of a point cloud gives each community an equal share of nodes so each community in this case would get $\frac{|V|}{K}$ nodes. So, if we call I_c the number of nodes that chose community center *c* we can have a loss function that increases as I_c gets higher (Dominance) or lower (Subdominance) and approaches 0 if the number is approximately $\frac{|V|}{K}$, which we can achieve with (6).

$$L_1 = \left(\frac{I_c * K - N}{N(K - 1)}\right)^2 \tag{6}$$

For calculating I_c we need to ensure that the function that calculates I_c is differentiable. To do this we use the Boltzmann Operator for a *SmoothMin* function (7) (α should be negative and large in absolute value), which can help determine the minimum distance for each node to all the other centers producing an almost-one-hot matrix specifying for each node the community it chose. This matrix can then easily be used to get an approximate I_c .

$$SmoothMin(\{x_i\}) = \frac{\sum_{i=1}^{n} x_i e^{\alpha x_i}}{\sum_{i=1}^{n} e^{\alpha x_i}}$$
(7)

Second, inspired by the Modularity metric we want to ensure that the distance between internal nodes is much lower compared to the distance between internal nodes and external nodes:

$$L_{2} = \frac{\sum_{i,j \in c} \|f_{i} - f_{j}\|_{2}}{\sum_{i \in c, j \notin c} \|f_{i} - f_{j}\|_{2}}$$
(8)

Third, to easily approach the nodes to their community center we first need to bring the community center closer to them, so we add another term that penalizes the distance between a community center and the nodes of its community:

$$L_3 = 1 - e^{-\sum_{i \in c} (\|\mu_c - f_i\|_2)} \tag{9}$$

Fourth, we assume that community centers are generally positioned far from each other, so we penalize their closeness:

$$L_4 = e^{-\sum_{c' \in K, c' \neq c} (\|\mu_{c'} - \mu_c\|_2)}$$
(10)

The loss function is then a weighted sum of the previous losses:

$$Loss = \alpha L1 + \beta L2 + \gamma L3 + \delta L4 \tag{11}$$

Where α , β , γ and δ are weights that sum up to 1 and denote the importance of each term.

3.3.2 Leaders SA

This method uses the Simulated Annealing metaheuristic (Kirkpatrick et al., 1983) to minimize as best as possible the loss function (11). It's important to note that here we don't need the derivative of the loss function so we can use the exact way to calculate it by directly using *argmin* instead of using *SmoothMin*. We start by creating an initial random set of leader nodes *S*. We get a neighbor of a solution *S* by choosing a random node from it which we'll be replaced by a randomly chosen node from V - S. This way



Figure 4: Mixed learning architecture.

we can execute Simulated Annealing to progressively improve the initial solution.

3.3.3 Walk Leaders

We propose a heuristic (Algorithm 1) to extract community leaders and automatically detect the number of communities K.

Since the walks matrix closely represents the distribution of nodes in hypothetical varying random walks starting from each node (we call such a distribution the context of the node), we can assume that the leader of a node *i* is more likely to be among the nodes with the most appearances in the context of *i*, so we start by picking for each node the *l* leading nodes in its context (we call them *potential leaders*, see line 1). The goal here is to choose the minimal set of nodes that contains at least one potential leader of each node. So, for selecting the leader of each node we first start with the potential leaders having the maximum number of appearances (see line 8), and then we only take the first one that is already chosen by a previous node (denoting the popularity of the leader among previous nodes, see line 11), in case no max count leader is already chosen by a previous node we choose the one with more appearances in the context of the node (see line 13). After iterating through all the nodes we end up with a current_leaders vector, but this vector can still have some leader nodes that don't choose themselves as leaders which isn't coherent. So, we used the Refinement of the idx vector method proposed by (Taheri and Bouyer, 2020a) to fix this conflict (see line 15).



The result of this method is a vector *current_leaders* assigning for each node *i* a leader node, so to get the number of communities we count the number of unique values in *current_leaders*.

3.4 Mixed Learning

The aim of this phase is to enhance node embeddings to represent community-aware high-order proximities between nodes, i.e. nodes that are in the same community will have embeddings that are close to each other. In this work, such an aim is achieved by improving both node embeddings and community cen-



Figure 5: Node embedder architecture.

ters/leaders in one learning process by bringing them closer to each other, which will form eventually accurate clusters. In both approaches that we'll present in the next sub-sections, the community embedder will either start from the weights it learned if used to initialize the centers, or will receive the centers/leaders embedding coordinates as initial weights.

3.4.1 Autoencoder-Based

One first approach to implement the learning process is to use the same autoencoder that was used to learn initial node embeddings. The autoencoder weights at the end of the initial node embedding phase are used as initial weights in this phase. The autoencoder and the Community Embedder are now trained at the same time. For a given node, the input to the autoencoder is its row in the level matrix while the input to the Community Embedder is the chosen community for that node as a one-hot vector. The training loss is a combined loss between the clustering loss (11) (using the output from the Community Embedder and the embeddings from the autoencoder) and the reconstruction error of the autoencoder as shown in figure 4.

3.4.2 Node Embedder-Based

Another approach proposed in this work is to use a new architecture to enhance node embeddings which we named "Node Embedder". This architecture inspired by Skip-gram (Mikolov et al., 2013) has four layers: an input layer, embedding layer, context layer, and output layer. A node encoded as a one-hot vector of size |V| is passed to the input layer which is connected densely to the embedding layer of size e. The embedding layer is densely connected to the context layer. The context layer contains c vectors of size |V| (so its size is c * |V|). c represents the context size. Each vector is activated with a Softmax function. This layer represents eventually the nodes that are in the context of the current node as one-hot vectors. Then the context layer is connected to the output layer via a one-dimensional convolutional kernel such that the i^{th} node of each vector in the context layer is connected to the i^{th} node in the output layer. Figure 5 summarizes this architecture.

The Node Embedder and Community Embedder are trained at the same time. The input for the Node Embedder is a one-hot vector representing the node, and in the same way as the autoencoder-based approach, the input for the Community Embedder is the chosen community for that node as a one-hot vector. The training loss is a combination of the clustering loss (11) (using the output of the Community Embedder and the embeddings from the Node Embedder) and a mean squared error between the output of the Node Embedder and the corresponsding row from the walk matrix. Figure 4 details the architecture explained in this paragraph.

3.5 Clustering

For community detection, one can use any clustering algorithm on the final node embeddings. Extracted clusters represent the detected communities. In this work, community centers/leaders can be used to detect communities by assigning each node to the closest community center. The results were also tested with Kmeans and Affinity Propagation.

4 COMPUTATIONAL RESULTS AND DISCUSSION

This section presents the results of computational experiments assessing the performance and effectiveness of SPACED for community detection in social networks. All algorithms and tests were developed in Python using TensorFlow and executed on a computer running Windows 11 Pro, equipped with 16 GB of RAM, an Intel(R) Core(TM) i5-10310U CPU at 1.70GHz, and an integrated Intel(R) UHD Graphics. Our experiments cover both synthetic and real-

world networks. The real-world datasets included:

Table 1:	Characteristics	of	the	tested	real-world	bench-
marks.						

Dataset	Comm- unities	Nodes	Edges
Karate (Zachary, 1977)	2	34	78
Dolphins (Lusseau et al., 2003)	2	62	159
Polbooks [V. Krebs, unpublished]	3	105	441
Football (Girvan and Newman, 2002)	12	115	613
Email (Leskovec et al., 2007)	42	1005	25571

For the synthetic networks we used the Lancichinetti-Fortunato-Radicchi (LFR) benchmark. The LFR datasets are characterized by the mixing parameter μ which controls the level of



Figure 6: NMI performance on real-world datasets.

community overlap and noise (the higher the μ value the more the noise). The LFR datasets comprised 128 nodes and 1024 edges, for the mixing parameter we tested on the values : 0.00, 0.25, 0.40, 0.45, and 0.50.

To evaluate the effectiveness of our solution, we employed the Normalized Mutual Information (NMI) metric which is a widely used measure in literature. NMI measures the similarity between the true community structure and the detected community structure, providing a score between 0 and 1, where 1 indicates perfect matching.

4.1 SPACED Performance Analysis on Real-World Datasets

The performance of our method on real-world datasets was assessed under both approaches, AD k (Automatic Detection of the number of communities "k") and NAD k (No-Automatic Detection, "k" is predefined). As shown in Table 2, our results demonstrated good accuracy in detecting the number of communities which is a significant advantage over many traditional methods that require this parameter to be known a priori. Also as summarized in figure 7, the community assignments produced by SPACED showed good NMI results both for the AD k and the NAD k modes. For instance, on the Karate network, our method achieved an NMI of 1.0, perfectly matching the ground truth.



Figures 6a to 6e illustrate the performance in terms of NMI for the Karate, Dolphins, Polbooks, Football, and Email networks respectively, compared against several well-known community detection algorithms, including Louvain (Blondel et al., 2008b),

Table 2: Real-world datasets: detected vs. ground truth number of communities.

Dataset	Ground Truth k	Detected k
Karate	2	2
Dolphins	2	4
Polbooks	3	5
Football	12	15
Email	42	39

Infomap (Rosvall and Bergstrom, 2012), GraphTrees (Dalleau et al., 2020a), and others (see Appendix). Our method, particularly the NAD k variant, consistently performed well across different datasets.

Also, SPACED has demonstrated reasonable execution times, with a direct correlation to network size. Figure 8 presents a comparative analysis of execution times (in seconds) across various real-world datasets, further illustrating that the rate of increase slows as network size grows. This demonstrates SPACED's strong potential for efficient scalability to larger networks.



Figure 8: Execution times in seconds for real-world datasets.

However, the method faces certain limitations in memory usage, primarily due to its reliance on the adjacency matrix, which becomes impractical for largescale networks and may lead to excessive memory consumption. This issue can be mitigated through technical optimizations, such as storing only non-zero values, as adjacency matrices in large-scale networks are typically sparse. Additionally, techniques like graph partitioning and distributed processing can be explored to further optimize the performance on large



Figure 9: SPACED results on synthetic datasets.

networks without compromising accuracy.

4.2 SPACED Performance Analysis on Synthetic Datasets

The synthetic datasets were generated using the LFR benchmark with varying levels of mixing parameters. Figure 9a illustrates the comparative performance of our method between the AD k and NAD k variants on the synthetic datasets for increasing values of the mixing parameter, and Table 3 shows the detected number of communities within the LFR networks. The relatively high NMI values indicate the good performance and robustness of our model, with an expected NMI value decrease as the mixing parameter increases due to the induced community noise. Also our method demonstrated its strength in correctly identifying the number of communities.

Table 3: Synthetic datasets: detected vs. ground truth number of communities.

Dataset	Ground Truth k	Detected k
LFR-0.00	4	5
LFR-0.25	4	4
LFR-0.40	4	4
LFR-0.45	4	3
LFR-0.50	4	4

Our method showed stable execution times independently from the graph complexity, making it also usable for graphs with a complex structure. The line chart in figure 9b shows a comparative study of the execution times in seconds for every synthetic dataset in an increasing order of complexity.

5 CONCLUSION

In this paper, we introduced SPACED, a deep learning model designed to uncover the community structure of networks solely from their topology by generating community-aware node embeddings. We evaluated its performance on several widely recognized benchmarks and compared it to other established methods, showcasing its competitiveness in the field.

Future enhancements for SPACED include exploring and refining the various pipeline variants to identify the optimal configuration. Additionally, while SPACED has shown strong stability and effectiveness on small to medium-sized datasets, it has not yet been tested on extremely large datasets. Thus, future work will need to address challenges in memory usage and execution time to ensure scalability.

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APPENDIX

References for the Compared Against Methods

DC-SBM (Karrer and Newman, 2011) NetMRF(He et al., 2018) ModMRF(Jin et al., 2020) GMRF (Jin et al., 2019b) SVM (Sui et al., 2016) APAS (Taheri and Bouyer, 2020b) ComNet-R(Cai et al., 2020) MRFasGCN(Jin et al., 2019a) DSFCD(Xie et al., 2018) RL (Paim et al., 2020) Louvain (Blondel et al., 2008a) Infomap (Rosvall and Bergstrom, 2012) GraphTrees (Dalleau et al., 2020b) CDMIC (Guo and Zhang, 2016) CMDR (Gong et al., 2014) PCSNMF(Shi et al., 2015).