

# Node Classification in Complex Social Graphs via Knowledge-Graph Embeddings and Convolutional Neural Network\*

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**Abstract.** The interactions between humans and their environment, comprising living and non-living entities, can be studied via Social Network Analysis (SNA). Node classification, as well as community detection tasks, are still open research problems in SNA. Hence, SNA has become an interesting and appealing domain in Artificial Intelligence (AI) research. Immanent facts about social network structures can be effectively harnessed for training AI models in a bid to solve node classification and community detection problems in SNA. Hence, crucial aspects such as the individual attributes of spatial social actors, and the underlying patterns of relationship binding these social actors must be taken into consideration in the course of analyzing the social network. These factors determine the nature and dynamics of a given social network. In this paper, we have proposed a unique framework, Representation Learning via Knowledge-Graph Embeddings and ConvNet (RLVECN), for studying and extracting meaningful facts from social network structures to aid in node classification as well as community detection tasks. Our proposition utilizes an edge sampling approach for exploiting features of the social graph, via learning the context of each actor with respect to neighboring actors/nodes, with the goal of generating vector-space embedding per actor. Successively, these relatively low-dimensional vector embeddings are fed as input features to a downstream classifier for classification tasks about the social graph/network. Herein RLVECN has been trained, tested, and evaluated on real-world social networks.

**Keywords:** Node Classification · Feature Learning · Feature Extraction  
· Dimensionality Reduction · Semi-supervised Learning.

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\* This research was supported by International Business Machines (IBM) and Compute Canada (SHARCNET).

## 1 Introduction and Related Literature

Humans inhabit in a planet comprised of several systems and ecosystems; and interaction is a natural phenomenon and characteristic obtainable in any given system or ecosystem. Thus, relationships between constituent entities in a given system/ecosystem is a strategy for survival and essentiality for the sustenance of a given system/ecosystem. Owing to the recent AI advances, these real-world complex systems and ecosystems can be effectively modelled as social network structures for analysis. Social network graphs [22] are intricate structures which present analytical challenges to Machine Learning (ML) and Deep Learning (DL) models because of their dynamic nature, complex links, and occasionally massive size. In this regard, we have proposed a new hybrid and DL-based model (RLVECN) based on an iterative learning approach [1] for solving (node) classification as well as clustering problems in SNA via an edge sampling strategy.

In SNA, the classification of nodes induces the formation of cluster(s). Consequently, clusters give rise to homophily in social networks. Basically, learning in RLVECN is induced via semi-supervised training. The architecture of RLVECN comprises two (2) distinct Representation Learning (RL) layers, viz: a Knowledge-Graph Embeddings (VE) layer and a Convolutional Neural Network (ConvNet) layer [13]. Both of these layers are trained by means of unsupervised learning. These layers are essentially feature-extraction and dimensionality-reduction layers where underlying knowledge and viable facts are automatically extracted from the social network structure [15]. The vector-embedding layer is responsible for projecting the feature representation of the social graph to a  $q$ -dimensional real-number space,  $\mathbb{R}^q$ . This is done by associating a real-number vector to every unique actor/node in the social network such that the cosine distance of any given tie/edge (a pair of actors) would capture a significant degree of correlation between the two associated actors. Furthermore, the ConvNet layer, which feeds on the vector-embedding layer, is responsible for further extraction of apparent features and/or representations from the social graph. Finally, a classification layer succeeds the RL layers; and it is trained by means of supervised learning. The classifier is based on a Neural Network (NN) architecture assembled using multiple and deep layers of stacked perceptrons (NN units) [6]. Every low-dimensional feature ( $X$ ), extracted by the representation-learning layers, is mapped to a corresponding output label ( $Y$ ); and these ( $X, Y$ ) pairs are used to supervise the training of the classifier such that it can effectively and efficiently learn how to identify clusters and classify actors within a given social network structure. Hence, the novelty of our research contribution are as stated below:

- (1) Proposition of a DL-based and hybrid model, RLVECN, which is aimed at solving node classification problems in SNA.
- (2) Detailed benchmarking reports with respect to classic objective functions used for classification tasks.
- (3) Comparative analysis, between RLVECN and state-of-the-art methodologies, against standard real-world social networks.

RLVECN is capable of learning the non-linear distributed features enmeshed in a social network [9]. We have evaluated RLVECN against an array of state-of-the-art models and RL methodologies which serve as our baselines, viz:

- (i) DeepWalk: Online Learning of Social Representations [19].
- (ii) GCN: Semi-Supervised Classification with Graph Convolutional Networks [11].
- (iii) LINE: Large-scale Information Network Embedding [25].
- (iv) Node2Vec: Scalable Feature Learning for Networks [8].
- (v) SDNE: Structural Deep Network Embedding [26].

## 2 Proposed Methodology and Framework

### 2.1 Definition of Problem

**Definition 1.** *Social Network, SN:* As expressed via equation 1 such that SN is a tuple comprising a set of actors/vertices,  $V$ ; a set of ties/edges,  $E$ ; a metadata function,  $f_V$ , which extends the definition of the vertices' set by mapping it to a given set of attributes,  $V'$ ; and a metadata function,  $f_E$ , which extends the definition of the edges' set by mapping it to a given set of attributes,  $E'$ . Thus, a graph function,  $G(V, E) \subset SN$

$$\begin{aligned}
 SN &= (V, E, f_V, f_E) \equiv (G, f_V, f_E) \\
 V &: |\{V\}| = M && \text{set of actors/vertices with size, } M \\
 E &: E \subset \{U \times V\} \subset \{V \times V\} && \text{set of ties/edges between } V \\
 f_V &: V \rightarrow V' && \text{vertices' metadata function} \\
 f_E &: E \rightarrow E' && \text{edges' metadata function}
 \end{aligned} \tag{1}$$

**Definition 2.** *Knowledge Graph, KG:*  $(\mathbb{E}, \mathbb{R})$  is a set comprising entities,  $\mathbb{E}$ , and relations,  $\mathbb{R}$ , between the entities. Thus, a KG [24][28] is defined via a set of triples,  $t : (u, p, v)$ , where  $u, v \in \mathbb{E}$  and  $p \in \mathbb{R}$ . Also, a KG [27] can be modelled as a social network, SN, such that:  $\mathbb{E} \rightarrow V$  and  $\mathbb{R} \rightarrow E$  and  $(\mathbb{E}, \mathbb{R}) \vdash f_V, f_E$ .

**Definition 3.** *Knowledge-Graph (Vector) Embeddings, X:* The vector-space embeddings,  $X$ , generated by the embedding layer are based on a mapping function,  $f$ , expressed via equation 2.  $f$  projects the representation of the graph's actors to a  $q$ -dimensional real space,  $\mathbb{R}^q$ , such that the existent ties between any given pair of actors,  $(u_i, v_j)$ , remain preserved via the homomorphism from  $V$  to  $X$ .

$$\begin{aligned}
 f &: V \rightarrow X \in \mathbb{R}^q \\
 f &: (u, p, v) \rightarrow X \in \mathbb{R}^q && \text{Knowledge-Graph Embeddings}
 \end{aligned} \tag{2}$$

**Definition 4.** *Node Classification:* Considering, SN, comprising partially labelled actors (or vertices),  $V_{lbl} \subset V : V_{lbl} \rightarrow Y_{lbl}$ ; and unlabelled vertices defined such that:  $V_{ulb} = V - V_{lbl}$ . A node-classification task aims at training a predictive function,  $f : V \rightarrow Y$ , that learns to predict the labels,  $Y$ , for all actors or vertices,  $V \subset SN$ , via knowledge harnessed from the mapping:  $V_{lbl} \rightarrow Y_{lbl}$ .

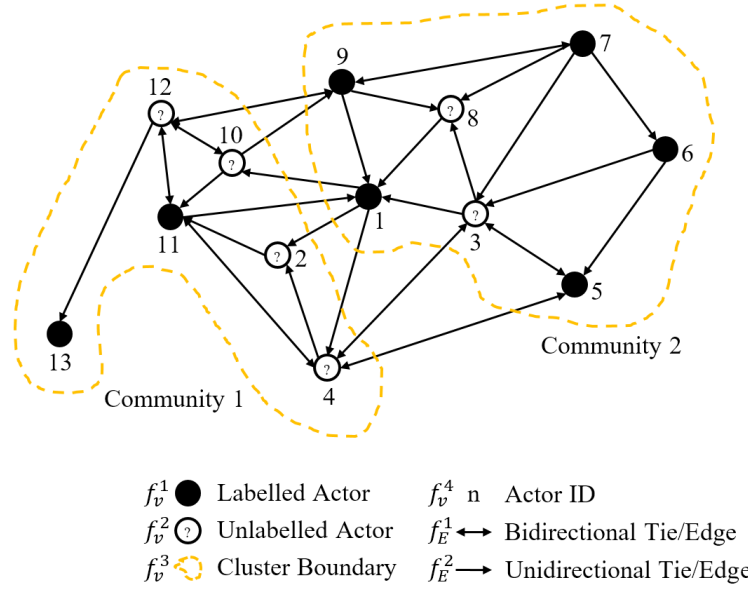


Fig. 1: Node classification task in social graphs

## 2.2 Proposed Methodology

Our proposition, RLVECN, is comprised of two (2) distinct Feature Learning (FL) layers, and one (1) classification layer.

**Representation Learning - Knowledge-Graph Embeddings Layer:** Given a social network,  $SN$ , defined by a set of actors/vertices,  $V : U \subset V \forall \{u_m, v_m\} \in V$ , and  $M : m \in M$  denotes the number of unique actors in  $SN$ . Additionally, let the ties/edges in  $SN$  be defined such that:  $E \subset \{U \times V\}$ ; where  $u_i \in V$  and  $v_j \in V$  represent a source\_vertex and a target\_vertex in  $E$ , respectively.

The objective function of the vector-embedding layer aims at maximizing the average logarithmic probability of the source\_vertex,  $u_i$ , being predicted as neighboring actor to the target\_vertex,  $v_j$ , with respect to all training pairs,  $\forall (u_i, v_j) \in E$ . Formally, the function is expressed as in equation 3:

$$\mu = \frac{1}{M} \sum_{m=1}^M \left( \sum_{(u_i, v_j) \in E} \log Pr(u_i|v_j) \right) \quad (3)$$

Consequently, in order to compute  $Pr(u_i|v_j)$ , we have to quantify the proximity of each target\_vertex,  $v_j$ , with respect to its source\_vertex,  $u_i$ . The vector-embedding model measures this adjacency/proximity as the cosine similarity between  $v_j$  and its corresponding  $u_i$ . Thus, the cosine distance is calculated as the dot product between the target\_vertex and the source\_vertex. Mathematically,  $Pr(u_i|v_j)$  is computed via a softmax function as defined in equation 4:

$$Pr(u_i|v_j) = \frac{\exp(u_i \cdot v_j)}{\sum_{m=1}^M \exp(u_m \cdot v_j)} \quad (4)$$

Hence, the objective function of our vector-embedding model with respect to the  $SN$  is as expressed by equation 5:

$$\sum_{(u_i, v_j) \in E} \log Pr(u_i | v_j) = \sum_{(u_i, v_j) \in E} \log \frac{\exp(u_i \cdot v_j)}{\sum_{m=1}^M \exp(u_m \cdot v_j)} \quad (5)$$

**Representation Learning - ConvNet Layer:** This layer comprises three (3) RL or FL operations, namely: convolution, non-linearity, and pooling operations. RLVECN utilizes a one-dimensional (1D) convolution layer [14] which is sandwiched between the vector-embedding and classification layers. Equation 6 expresses the 1D-convolution operation:

$$\begin{aligned} FeatureMap(F) &= 1D\_InputMatrix(X) * Kernel(K) \\ f_i &= (X * K)_i = (K * X)_i = \sum_{j=0}^{J-1} x_j \cdot k_{i-j} = \sum_{j=0}^{J-1} k_j \cdot x_{i-j} \end{aligned} \quad (6)$$

where  $f_i$  represents a cell/matrix position in the Feature Map;  $k_j$  denotes a cell position in the Kernel; and  $x_{i-j}$  denotes a cell/matrix position in the 1D-Input (data) matrix.

The non-linearity operation is a rectified linear unit (ReLU) function which introduces non-linearity after the convolution operation since real-world problems usually exist in non-linear form(s). As a result, the rectified feature/activation map is computed via:  $r_i \in R = g(f_i \in F) = \max(0, F)$ .

The pooling operation is responsible for reducing the input width of each rectified activation map while retaining its vital properties. In this regard, the *Max Pooling* function is defined such that the resultant pooled (or downsampled) feature map is generated via:  $p_i \in P = h(r_i \in R) = \maxPool(R)$ .

**Classification - Multi-Layer Perceptron (MLP) Classifier Layer:** This is the last layer of our proposed RLVECN’s architecture, and it succeeds the RL layers. The pooled feature maps, generated by the Representation Learning layers, contain high-level features extracted from the constituent actors of the social network structure. Hence, the classification layer utilizes these extracted “high-level features” for identifying clusters, based on the respective classes, contained in the social graph. In this regard, a MLP [5] function is defined as a mathematical function,  $f_c$ , mapping some set of input values,  $P$ , to their respective output labels,  $Y$ . In other words,  $Y = f_c(P, \Theta)$ , and  $\Theta$  denotes a set of parameters. The MLP [4] function learns the values of  $\Theta$  that will result in the best decision,  $Y$ , approximation for the input set,  $P$ . The MLP classifier output is a probability distribution which indicates the likelihood of a feature representation belonging to a particular output class. Our MLP [10] classifier is modelled such that sequential layers of NN units are stacked against each other to form a Deep Neural Network (DNN) structure [3], [16].

**Node Classification Algorithm:** Defined via algorithm 1.

**Algorithm 1** Proposed Algorithm for Node Classification**Input:**  $\{V, E, Y_{lbl}\} \equiv \{\text{Actors, Ties, Ground-Truth Labels}\}$ **Output:**  $\{Y_{ulb}\} \equiv \{\text{Predicted Labels}\}$ **Preprocessing:** $V_{lbl}, V_{ulb} \subset V = V_{lbl} \cup V_{ulb}$  //  $V_{lbl}$  : Labelled actors //  $V_{ulb}$  : Unlabelled actors $E : (u_i, v_j) \in \{U \times V\}$  //  $(u_i, v_j) \equiv (\text{source, target})$  $E_{train} = E_t : u_i, v_j \in V_{lbl}$  //  $|E_{train}| = \sum indegree(V_{lbl}) + \sum outdegree(V_{lbl})$  $E_{pred} = E_p : u_i, v_j \in V_{ulb}$  $f_c \leftarrow \text{Initialize}$  // Construct classifier model**Training:****for**  $t \leftarrow 0$  **to**  $|E_{train}|$  **do** $f : E_t \rightarrow [X \in \mathbb{R}^q]$  // Embedding operation $f_t \in F = (K * X)_t$  // Convolution operation $r_t \in R = g(F) = \max(0, f_t)$  $p_t \in P = h(R) = \maxPool(r_t)$  $f_c | \Theta : p_t \rightarrow Y_{lbl}$  // MLP classification operation**end for****return**  $Y_{ulb} = f_c(E_{pred}, \Theta)$ **2.3 Proposed Architecture/Framework**

Fig. 2 illustrates the architecture of our proposition, RLVECNet.

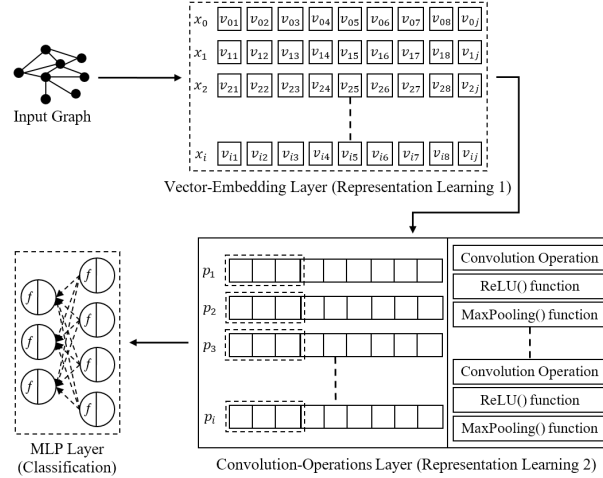


Fig. 2: Proposed system architecture

**3 Data sets and Materials****3.1 Data sets**

With regard to Table 1 herein, six (6) real-world benchmark social-graph data sets were employed for experimentation and evaluation, viz: Cora [23] [20], CiteSeer

[23] [20], Facebook Page-Page webgraph [21], PubMed-Diabetes [17], Internet-Industry partnerships [12] [2], and Terrorists-Relationship [29].

Table 1: Benchmark data sets

Data set	Classes $\rightarrow$ {label: ‘description’}
Cora	$G(V, E) = G(2708, 5429)$ {C1: ‘Case_Based’, C2: ‘Genetic_Algorithms’, C3: ‘Neural_Networks’, C4: ‘Probabilistic_Methods’, C5: ‘Reinforcement_Learning’, C6: ‘Rule_Learning’, C7: ‘Theory’}
CiteSeer	$G(V, E) = G(3312, 4732)$ {C1: ‘Agents’, C2: ‘Artificial Intelligence’, C3: ‘Databases’, C4: ‘Information Retrieval’, C5: ‘Machine Learning’, C6: ‘Human-Computer Interaction’}
Facebook Page2Page	$G(V, E) = G(22470, 171002)$ {C1: ‘Companies’, C2: ‘Governmental Organizations’, C3: ‘Politicians’, C4: ‘Television Shows’}
PubMed Diabetes	$G(V, E) = G(19717, 44338)$ {C1: ‘Diabetes Mellitus - Experimental’, C2: ‘Diabetes Mellitus - Type 1’, C3: ‘Diabetes Mellitus - Type 2’}
Internet Industry	$G(V, E) = G(219, 631)$ {C1: ‘Content Sector’, C2: ‘Infrastructure Sector’, C3: ‘Commerce Sector’}
Terrorists Relation	$G(V, E) = G(851, 8592)$ {C1: ‘Content Sector’, C2: ‘Infrastructure Sector’, C3: ‘Commerce Sector’}

### 3.2 Data Preprocessing

All benchmark data sets ought to be comprised of actors and ties already encoded as discrete data (natural-number format). However, Cora, CiteSeer, Facebook-Page2Page, PubMed-Diabetes, and Terrorists-Relation data sets are made up of nodes and/or edges encoded in mixed formats (categorical and numerical formats). Thus, it is necessary to transcode these non-numeric (categorical) entities to their respective discrete (numeric) data representation, without semantic loss, via an injective function that maps each distinct entry in the categorical-entity domain to distinct numeric values in the discrete-data codomain,  $f_m : \text{categorical} \rightarrow \text{discrete}$ . Thereafter, the numeric representation of all benchmark data sets are normalized,  $f_n : \text{discrete} \rightarrow \text{continuous}$ , prior to training against RLVECN and the benchmark models. Also, only edgelist ties,  $E \subset G$ , whose constituent actors are present in the nodelist,  $V \subset G$ , were used for training/testing our model.

Table 2: Configuration of experimentation hyperparameters

Training Set: 80%	Test Set: 20%	Network Width: 640
Batch Size: 256	Optimizer: <i>AdaMax</i>	Network Depth: 6
Epochs: $1.8 * 10^2$	Activation: <i>ReLU</i>	Dropout: $4.0 * 10^{-1}$
Learning Rate: $1.0 * 10^{-3}$	Learning Decay: 0.0	Embed Dimension: 100

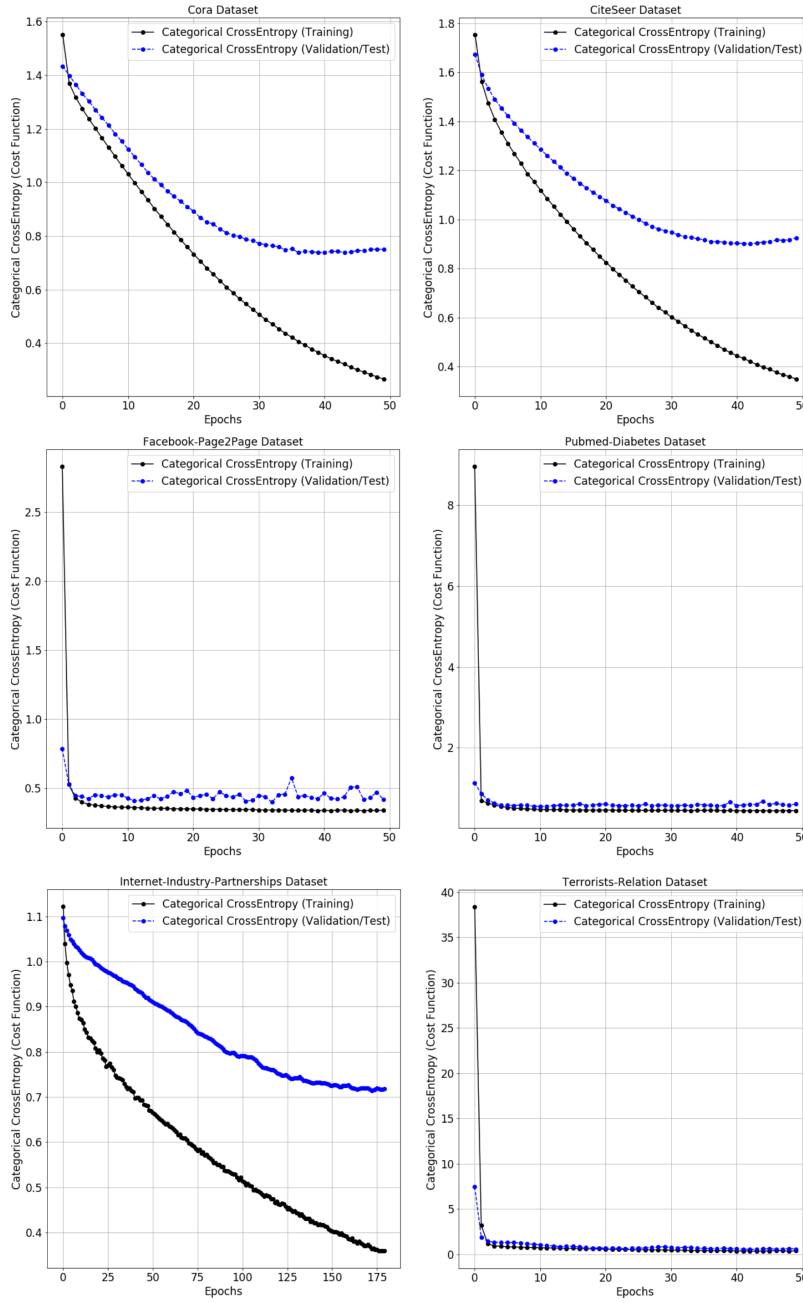


Fig. 3: RLVEC�’s learning curves during node-classification training against Cora, CiteSeer, Facebook-Page2Page, PubMed-Diabetes, Internet-Industry-Partnership, and Terrorists-Relation data sets (loss function *vs* training epochs)

#### 4 Experiment, Results, and Discussions

RLVEC�’s experimentation setup was tuned in accordance with the hyperparameters shown in Table 2. Our evaluations herein were recorded with reference



Table 3: Node-classification over Cora data set. Results are based on the set apart validation sample - data set *vs* models.

Model	Metric	Cora Dataset									Points
		C1	C2	C3	C4	C5	C6	C7	$\mu$		
RLVEC	PC	0.85	0.78	0.80	0.88	0.72	0.90	0.81	0.82	14	
	RC	0.86	0.93	0.81	0.87	0.75	0.91	0.78	0.84		
	F1	<b>0.86</b>	<b>0.85</b>	<b>0.81</b>	<b>0.87</b>	<b>0.74</b>	<b>0.91</b>	<b>0.79</b>	0.83		
	AC	0.93	0.98	0.96	0.95	0.93	0.97	0.96	0.95		
	RO	<b>0.90</b>	<b>0.96</b>	<b>0.90</b>	<b>0.92</b>	<b>0.85</b>	<b>0.95</b>	<b>0.88</b>	0.91		
	SP	541	134	214	405	294	345	237	310		
GCN	PC	0.87	0.95	0.89	0.92	0.85	0.89	0.87	0.89	3	
	RC	0.85	0.73	0.65	0.82	0.58	0.85	0.73	0.74		
	F1	<b>0.86</b>	0.83	0.75	<b>0.87</b>	0.69	0.87	<b>0.79</b>	0.81		
	AC	0.89	0.93	0.91	0.92	0.88	0.93	0.91	0.91		
	RO	0.88	0.83	0.80	0.89	0.75	0.90	0.83	0.84		
	SP	164	36	43	85	70	84	60	77		
Node2Vec	PC	0.58	0.78	0.72	0.81	0.80	0.84	0.82	0.76	0	
	RC	0.85	0.50	0.53	0.68	0.64	0.74	0.60	0.65		
	F1	0.69	0.61	0.61	0.74	0.71	0.78	0.69	0.69		
	AC	0.77	0.96	0.95	0.92	0.93	0.94	0.94	0.92		
	RO	0.79	0.75	0.76	0.83	0.81	0.86	0.79	0.80		
	SP	164	36	43	85	70	84	60	77		
DeepWalk	PC	0.57	0.58	0.72	0.58	0.68	0.72	0.63	0.64	0	
	RC	0.80	0.42	0.42	0.59	0.39	0.63	0.65	0.56		
	F1	0.67	0.48	0.53	0.58	0.49	0.67	0.64	0.58		
	AC	0.76	0.94	0.94	0.87	0.90	0.90	0.92	0.89		
	RO	0.77	0.70	0.70	0.75	0.68	0.79	0.80	0.74		
	SP	164	36	43	85	70	84	60	77		
LINE	PC	0.35	0.86	0.80	0.65	0.50	0.43	0.61	0.60	0	
	RC	0.85	0.17	0.19	0.35	0.20	0.15	0.23	0.31		
	F1	0.50	0.28	0.30	0.46	0.29	0.23	0.34	0.34		
	AC	0.48	0.94	0.93	0.87	0.87	0.84	0.90	0.83		
	RO	0.59	0.58	0.59	0.66	0.59	0.56	0.61	0.60		
	SP	164	36	43	85	70	84	60	77		
SDNE	PC	0.37	0.83	0.70	0.60	0.54	0.64	0.64	0.62	0	
	RC	0.91	0.14	0.16	0.35	0.20	0.27	0.12	0.31		
	F1	0.53	0.24	0.26	0.44	0.29	0.38	0.20	0.33		
	AC	0.50	0.94	0.93	0.86	0.87	0.86	0.89	0.84		
	RO	0.62	0.57	0.58	0.65	0.59	0.62	0.55	0.60		
	SP	164	36	43	85	70	84	60	77		

to a range of objective functions. Thus, Categorical Cross Entropy was employed as the cost/loss function; while the fitness/utility was measured based on the following metrics: Precision (PC), Recall (RC), F-measure or F1-score (F1), Accuracy (AC), and Area Under the Receiver Operating Characteristic Curve (RO). Moreover, the objective functions have been computed against each benchmark data set with regard to the constituent classes (or categories) present in each data set. The Support (SP) represents the number of ground-truth samples per class/category for each data set.

In a bid to avoid sample bias across-the-board, we have used exactly the same SP for all models inclusive of RLVEC model. However, since RLVEC is based on an edge-sampling technique; the SP recorded against RLVEC model represent

Table 4: Node-classification over CiteSeer data set. Results are based on the set apart test sample - data set *vs* models.

Model	Metric	CiteSeer Dataset							Points
		C1	C2	C3	C4	C5	C6	$\mu$	
RLVECN	PC	0.76	0.81	0.78	0.43	0.88	0.60	0.71	12
	RC	0.84	0.83	0.79	0.60	0.79	0.65	0.75	
	F1	<b>0.80</b>	<b>0.82</b>	<b>0.79</b>	<b>0.50</b>	<b>0.83</b>	<b>0.63</b>	0.73	
	AC	0.93	0.88	0.92	0.93	0.96	0.89	0.92	
	RO	<b>0.90</b>	<b>0.87</b>	<b>0.87</b>	<b>0.78</b>	<b>0.89</b>	<b>0.79</b>	0.85	
	SP	304	609	377	107	225	275	316	
GCN	PC	0.80	0.78	0.86	0.95	0.91	0.75	0.84	2
	RC	0.76	0.76	0.73	0.08	0.67	0.54	0.59	
	F1	0.78	0.77	<b>0.79</b>	0.15	0.77	<b>0.63</b>	0.65	
	AC	0.88	0.87	0.88	0.91	0.89	0.83	0.88	
	RO	0.84	0.83	0.83	0.53	0.81	0.72	0.76	
	SP	119	134	140	50	102	118	111	
Node2Vec	PC	0.57	0.55	0.49	0.33	0.55	0.38	0.48	0
	RC	0.55	0.60	0.66	0.06	0.45	0.40	0.45	
	F1	0.56	0.58	0.56	0.10	0.50	0.39	0.45	
	AC	0.85	0.82	0.78	0.92	0.86	0.78	0.84	
	RO	0.73	0.74	0.74	0.53	0.69	0.63	0.68	
	SP	119	134	140	50	102	118	111	
DeepWalk	PC	0.46	0.53	0.43	0.43	0.47	0.33	0.44	0
	RC	0.51	0.54	0.57	0.06	0.41	0.32	0.40	
	F1	0.49	0.54	0.49	0.11	0.44	0.32	0.40	
	AC	0.81	0.81	0.75	0.92	0.84	0.76	0.82	
	RO	0.69	0.71	0.69	0.53	0.66	0.59	0.65	
	SP	119	134	140	50	102	118	111	
SDNE	PC	0.37	0.50	0.24	0.20	0.45	0.31	0.35	0
	RC	0.19	0.27	0.77	0.02	0.14	0.09	0.25	
	F1	0.25	0.35	0.36	0.04	0.21	0.14	0.23	
	AC	0.80	0.80	0.42	0.92	0.84	0.80	0.76	
	RO	0.56	0.60	0.55	0.51	0.55	0.52	0.55	
	SP	119	134	140	50	102	118	111	
LINE	PC	0.18	0.30	0.28	0.60	0.22	0.27	0.31	0
	RC	0.15	0.47	0.39	0.06	0.12	0.21	0.23	
	F1	0.16	0.36	0.32	0.11	0.15	0.24	0.22	
	AC	0.72	0.67	0.65	0.93	0.80	0.76	0.76	
	RO	0.50	0.59	0.56	0.53	0.52	0.55	0.54	
	SP	119	134	140	50	102	118	111	

the numbers of edges/ties used for computation as explained in algorithm 1. With regard to the standard node-classification tasks herein, the performance of our RLVECN model during benchmarking against five(5) popular baselines (DeepWalk, GCN, LINE, Node2Vec, SDNE); and when evaluated against the validation/test samples for the benchmark data sets are as documented in Table 3, Table 4, Table 5, and Table 6 respectively. Consequently, Fig. 3 graphically shows the learning-progress curves during the node-classification tasks using our proposed RLVECN model; and when training over the benchmark data sets. Hence, the dotted-black lines represent learning progress over the training set; and the dotted-blue lines represent learning progress over the test set.

Table 5: Node-classification experiment over Facebook Page-Page webgraph and PubMed-Diabetes data sets. Results are based on the reserved test sample - data set *vs* models. N.B.: Mtc = Metric (fitness function); Pts = Points.

Model	Mtc	Facebook-Page2Page					Pts	PubMed-Diabetes				Pts
		C1	C2	C3	C4	$\mu$		C1	C2	C3	$\mu$	
RLVECN	PC	0.87	0.95	0.91	0.87	0.90	8	0.76	0.83	0.84	0.81	6
	RC	0.84	0.85	0.85	0.86	0.85		0.60	0.88	0.91	0.80	
	F1	<b>0.85</b>	<b>0.90</b>	<b>0.88</b>	<b>0.86</b>	0.87		<b>0.67</b>	<b>0.86</b>	<b>0.87</b>	0.80	
	AC	0.96	0.90	0.94	0.97	0.94		0.89	0.88	0.90	0.89	
	RO	<b>0.97</b>	<b>0.97</b>	<b>0.98</b>	<b>0.98</b>	0.98		<b>0.92</b>	<b>0.94</b>	<b>0.95</b>	0.94	
	SP	9989	33962	16214	6609	16694		3300	7715	7170	6062	
Node2Vec	PC	0.81	0.84	0.81	0.84	0.83	0	0.74	0.47	0.49	0.57	0
	RC	0.82	0.87	0.85	0.67	0.80		0.03	0.65	0.55	0.41	
	F1	0.81	0.85	0.83	0.74	0.81		0.05	0.55	0.52	0.37	
	AC	0.89	0.91	0.91	0.93	0.91		0.80	0.57	0.60	0.66	
	RO	0.87	0.90	0.89	0.82	0.87		0.51	0.58	0.59	0.56	
	SP	1299	1376	1154	665	1124		821	1575	1548	1315	
DeepWalk	PC	0.75	0.84	0.76	0.75	0.78	0	0.65	0.57	0.58	0.60	0
	RC	0.81	0.85	0.82	0.52	0.75		0.15	0.67	0.71	0.51	
	F1	0.78	0.84	0.79	0.62	0.76		0.24	0.62	0.63	0.50	
	AC	0.87	0.90	0.89	0.90	0.89		0.81	0.67	0.68	0.72	
	RO	0.85	0.89	0.87	0.75	0.84		0.56	0.67	0.69	0.64	
	SP	1299	1376	1154	665	1124		821	1575	1548	1315	
LINE	PC	0.53	0.66	0.72	0.66	0.64	0	0.48	0.42	0.44	0.45	0
	RC	0.72	0.71	0.59	0.29	0.58		0.05	0.60	0.46	0.37	
	F1	0.61	0.68	0.65	0.40	0.59		0.08	0.50	0.45	0.34	
	AC	0.73	0.80	0.83	0.87	0.81		0.79	0.51	0.56	0.62	
	RO	0.73	0.77	0.75	0.63	0.72		0.52	0.53	0.54	0.53	
	SP	1299	1376	1154	665	1124		821	1575	1548	1315	
SDNE	PC	0.49	0.80	0.70	0.65	0.66	0	0.65	0.43	0.74	0.61	0
	RC	0.90	0.63	0.50	0.19	0.56		0.05	0.96	0.17	0.39	
	F1	0.64	0.70	0.58	0.29	0.55		0.10	0.59	0.27	0.32	
	AC	0.70	0.84	0.82	0.86	0.81		0.80	0.48	0.65	0.64	
	RO	0.76	0.78	0.71	0.58	0.71		0.52	0.56	0.56	0.55	
	SP	1299	1376	1154	665	1124		821	1575	1548	1315	

Tables 3, 4, 5, and 6 have clearly tabulated our results as a multi-classification task over the benchmark data sets. Thus, for each class per data set, we have laid emphasis on the F1 (the weighted average of the PC and RC metrics) and RO. Therefore, we have highlighted the model which performed best (based on F1 and RO metrics) for each classification task using a **bold font**. Additionally, we have employed a point-based ranking standard to ascertain the fittest model for each node classification task. The model with the highest aggregate point signifies the fittest model for the specified task, and so on in a descending order of aggregate points. Accordingly, as can be seen from our tabular results, our proposed methodology (RLVECN) is at the top with the highest fitness points.

$L2$  regularization ( $L2 = 0.04$ ) [7] and early stopping [18] were used herein as ad-don regularization techniques to overcome overfitting incurred during RLVECN's training. Hence, the application of early stopping with respect to RLVECN's training over the benchmark data sets were, viz: 50 epochs (Cora, CiteSeer, Facebook-Page2Page, PubMed-Diabetes, Terrorists-Relation) and 180 epochs

Table 6: Node-classification experiment over Internet-Industry-Partnership and Terrorists-Relationship data sets. Results are based on the reserved validation sample - data set *vs* models. N.B.: Mtc = Metric (fitness function); Pts = Points.

Model	Mtc	Internet-Industry-Partnership				Pts	Terrorists-Relation Dataset					Pts
		C1	C2	C3	$\mu$		C1	C2	C3	C4	$\mu$	
RLVECN	PC	0.33	0.96	0.29	0.53	5	0.93	0.91	0.46	1.00	0.83	6
	RC	0.65	0.77	0.76	0.73		0.97	0.97	0.42	0.97	0.83	
	F1	<b>0.44</b>	<b>0.86</b>	0.42	0.57		<b>0.95</b>	<b>0.94</b>	0.44	<b>0.98</b>	0.83	
	AC	0.84	0.78	0.87	0.83		0.95	0.98	0.89	0.99	0.95	
	RO	<b>0.76</b>	<b>0.81</b>	<b>0.82</b>	0.80		<b>0.98</b>	<b>1.00</b>	0.85	<b>1.00</b>	0.96	
	SP	26	238	17	94		1706	491	319	561	769	
GCN	PC	No experiment due to the absence of vectorized feature set for this data set				0	0.94	0.74	0.67	0.96	0.83	5
	RC						0.90	0.95	0.60	1.00	0.86	
	F1						0.92	0.83	<b>0.63</b>	<b>0.98</b>	0.84	
	AC						0.92	0.95	0.88	0.99	0.94	
	RO						<b>0.98</b>	0.99	<b>0.91</b>	<b>1.00</b>	0.97	
	SP						92	21	30	27	43	
DeepWalk	PC	0.50	0.81	0.36	0.56	1	0.88	0.82	0.64	0.86	0.80	0
	RC	0.12	0.93	0.44	0.50		0.90	0.86	0.53	0.93	0.81	
	F1	0.20	<b>0.86</b>	0.40	0.49		0.89	0.84	0.58	0.89	0.80	
	AC	0.82	0.82	0.73	0.79		0.88	0.96	0.86	0.96	0.92	
	RO	0.55	0.79	0.62	0.65		0.88	0.92	0.73	0.95	0.87	
	SP	8	27	9	15		92	21	30	27	43	
Node2Vec	PC	0.00	0.68	0.75	0.48	1	0.86	0.82	0.60	0.86	0.79	0
	RC	0.00	1.00	0.33	0.44		0.88	0.86	0.50	0.93	0.79	
	F1	0.00	0.81	<b>0.46</b>	0.42		0.87	0.84	0.55	0.89	0.79	
	AC	0.82	0.70	0.84	0.79		0.86	0.96	0.85	0.96	0.91	
	RO	0.50	0.62	0.65	0.59		0.86	0.92	0.71	0.95	0.86	
	SP	8	27	9	14.67		92	21	30	27	43	
LINE	PC	0.00	0.61	0.00	0.20	0	0.82	0.82	0.58	0.92	0.79	0
	RC	0.00	1.00	0.00	0.33		0.92	0.86	0.37	0.85	0.75	
	F1	0.00	0.76	0.00	0.25		0.87	0.84	0.45	0.88	0.76	
	AC	0.82	0.61	0.80	0.74		0.85	0.96	0.84	0.96	0.90	
	RO	0.50	0.50	0.50	0.50		0.84	0.92	0.65	0.92	0.83	
	SP	8	27	9	15		92	21	30	27	43	
SDNE	PC	0.00	0.61	0.00	0.20	0	0.77	0.90	0.56	1.00	0.81	0
	RC	0.00	1.00	0.00	0.33		0.92	0.86	0.30	0.85	0.73	
	F1	0.00	0.76	0.00	0.25		0.84	0.88	0.39	0.92	0.76	
	AC	0.82	0.61	0.80	0.74		0.81	0.97	0.84	0.98	0.90	
	RO	0.50	0.50	0.50	0.50		0.80	0.92	0.62	0.93	0.82	
	SP	8	27	9	15		92	21	30	27	43	

(Internet-Industry-Partnership). We have used a mini-batch size of 256 for training and testing/validating because we want to ensure that sufficient patterns are extracted by RLVECN during training before its network weights are updated.

## 5 Limitations, Conclusion, Future Work, and Acknowledgements

The benchmark models evaluated herein were implemented using their default parameters. We were not able to evaluate GCN [11] against Facebook Page-Page webgraph, PubMed-Diabetes, and Internet-Industry-Partnership data sets;

because each of these aforementioned data sets does not possess a vectorized feature set which is required by GCN model for input-data processing. Overall, RLVECN’s remarkable performance with respect to our benchmarking results can be attributed to the following:

- (1) The RL kernel of RLVECN is constituted of two (2) distinct layers of FL, viz: Knowledge-Graph Embeddings and ConvNet [13].
- (2) The high-quality data preprocessing techniques employed herein with respect to the benchmark data sets. We ensured that all constituent actors of a given social graph were transcoded to their respective discrete data representations, without any loss in semantics, and normalized prior to training and/or testing.

In conclusion, we intend to expand our experimentation scope to include more social network data sets and benchmark models. This research was made possible by International Business Machines (IBM), SHARCNET and Compute Canada ([www.computecanada.ca](http://www.computecanada.ca)).

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