

Survey

Survey 1-[2018IEEE]

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A comprehensive survey of graph embedding: problems, techniques and applications

[J]. IEEE Transactions on Knowledge and Data Engineering, 2018.

Graph Embedding Problem Settings

- Graph Embedding Input
 - Homogeneous Graph
 - Undirected, unweighted, directed, weighted
 - Challenge: How to capture the diversity of connectivity patterns observed in graphs
 - Heterogeneous Graph
 - Community-based question answering (cQA) sites, Multimedia Networks, Knowledge Graphs
 - Challenge: How to explore global consistency between different types of objects, and how to deal with the imbalances of different-typed objects, if any
 - Graph with Auxiliary Information
 - Label, Attribute, Node feature, Information propagation, Knowledge base
 - Challenge: How to incorporate the rich and unstructured information so that the learnt embeddings are both representing the topological structure and discriminative in terms of the auxiliary information.
 - Graph Constructed from Non-relational Data
 - Challenge: How to construct a graph that encodes the pairwise relations between instances and how to preserve the generated node proximity matrix in the embedded space.
- Graph Embedding Output:
 - how to find a suitable type of embedding output which meets the needs of the specific application task
- Node Embedding
 - Challenge: How to define the node pairwise proximity in various types of input graph and how to encode the proximity in the learnt embeddings.
- Edge Embedding
 - Challenge: How to define the edge-level similarity and how to model the asymmetric property of the edges, if any.
- Hybrid Embedding
 - Substructure embedding, community embedding
 - Challenge: How to generate the target substructure and how to embed different types of graph components in one common space.
- Whole-Graph Embedding
 - Challenge: How to capture the properties of a whole graph and how to make a trade-off between expressivity and efficiency.

Graph Embedding Techniques

- Matrix Factorization
 - Graph Laplacian Eigenmaps
 - The graph property to be preserved can be interpreted as pairwise node similarities, thus a larger penalty is imposed if two nodes with larger similarity are embedded far apart
 - Node Proximity Matrix Factorization
 - Node proximity can be approximated in low-dimensional space using matrix factorization. The objective of preserving the node proximity is to minimize the loss during the approximation
- Deep Learning
 - With Random Walk
 - The second-order proximity in a graph can be preserved in the embedded space by maximizing the probability of observing the neighbourhood of a node conditioned on its embedding
 - There are usually two solutions to approximate the full softmax: hierarchical softmax and negative sampling.
 - Without Random Walk
 - The multi-layered learning architecture is a robust and effective solution to encode the graph into a low dimensional space
 - Autoencoder, Deep Neural Network, Others
- Edge Reconstruction
 - Maximize Edge Reconstruct Probability
 - Good node embedding maximizes the probability of generating the observed edges in a graph
 - Minimize Distance-based Loss
 - The node proximity calculated based on node embedding should be as close to the node proximity calculated based on the observed edges as possible
 - Minimize Margin-based Ranking Loss
 - A node's embedding is more similar to the embedding of relevant nodes than that of any other irrelevant node
- Graph Kernel
- The whole graph structure can be represented as a vector containing the counts of elementary substructures that are decomposed from it
 - Based on Graphlet
 - Based on Subtree Patterns
 - Based on Random Walks
- Generative Model
 - Embed Graph into Latent Space
 - Nodes are embedded into the latent semantic space where the distances among node embeddings can explain the observed graph structure
 - Incorporate Semantics for Embedding
 - Nodes who are not only close in the graph but also having similar semantics should be embedded closer. The node semantics can be detected from node descriptions via a generative model
- Hybrid Techniques and Others

Applications

- Node Related Applications

- Node Classification
- Node Clustering
- Node Recommendation/Retrieval/Ranking
- Edge Related Applications
 - Link Prediction and Graph Reconstruction
 - Triple Classification
- Graph Related Applications
 - Graph Classification
 - Visualization
- Other Applications
 - Knowledge graph related
 - Multimedia network related
 - Information propagation related
 - Social networks alignment

Survey 2-[2017]

Hamilton W L, Ying R, Leskovec J.

Representation Learning on Graphs: Methods and Applications

[J]. arXiv preprint arXiv:1709.05584, 2017.

一、Embedding nodes

1. Overview of approaches: An encoder-decoder perspective

The intuition behind the encoder-decoder idea is the following: if we can learn to decode high-dimensional graph information—such as the global positions of nodes in the graph or the structure of local graph neighborhoods—from encoded low-dimensional embeddings, then, in principle, these embeddings should contain all information necessary for downstream machine learning tasks

encoder是将结点转换成embeddings, decoder是接受一系列embeddings, 然后从中解码出用户指定的统计量, 比如结点属于哪个类, 两个结点之间是否存在边等。

目标是优化encoder和decoder的mapping来最小化误差或损失。损失是decoder的结果和真实结果之间的差异。

四个方面的不同: pairwise similarity function, encoder function, decoder function, loss function

1. Shallow embedding approaches

For these shallow embedding approaches, the encoder function—which maps nodes to vector embeddings—is simply an “embedding lookup”

- Factorization-based approaches
 - Laplacian eigenmaps: decoder是两个向量之差的L2范数的平方, 损失函数是decoder的结果的加权和, 权重是两个结点的相似性。

- Inner-product methods: decoder是两个向量的内积, 例如The Graph Factorization (GF) algorithm, GraRep, and HOPE, 他们的损失函数都是MSE: decoder的结果和实际相似度的差的L2范数的平方。他们的不同之处是相似性的度量。GF直接用邻接矩阵, GraRep用邻接矩阵的平方, HOPE用based on Jaccard neighborhood overlaps
- 他们的共同点是: Loss function基本上是: $\|Z^T Z - S\|_2^2$, embedding矩阵Z和相似性矩阵S
- Random walk approaches
 - DeepWalk and node2vec: decoder是从结点i出发在T步内经过结点j的概率; 交叉熵损失函数
 - Large-scale information network embeddings (LINE)
 - HARP: Extending random-walk embeddings via graph pre-processing: a graph coarsening procedure is used to collapse related nodes in G together into “supernodes”, and then DeepWalk, node2vec, or LINE is run on this coarsened graph. After embedding the coarsened version of G, the learned embedding of each supernode is used as an initial value for the random walk embeddings of the supernode’s constituent nodes
 - Additional variants of the random-walk idea

缺点:

- 1.No parameters are shared between nodes in the encoder
- 2.Shallow embedding also fails to leverage node attributes during encoding
- 3.Shallow embedding methods are inherently transductive

1. Generalized encoder-decoder architectures

- Neighborhood autoencoder methods: they use autoencoders, 例如DNCR, SDNE, Extract high-dimensional neighborhood vector, (si contains vi’s proximity to all other nodes), Compress si to low-dimensional embedding
- Neighborhood aggregation and convolutional encoders: they generate embeddings for a node by aggregating information from its local neighborhood, 例如GCN, column networks, GraphSAGE

1. Incorporating task-specific supervision

cross-entropy loss, backpropagation

1. Extensions to multi-modal graphs

- Dealing with different node and edge types
- Tying node embeddings across layers: 如OhmNet

1. Embedding structural roles

struc2vec, GraphWave

1. Applications of node embeddings

- Visualization and pattern discovery
- Clustering and community detection
- Node classification and semi-supervised learning
- Link prediction

二、Embedding subgraphs

1. Sets of node embeddings and convolutional approaches

The basic intuition behind these approaches is that they equate subgraphs with sets of node embeddings

- Sum-based approaches
- Graph-coarsening approaches
- Further variations

1. Graph neural networks

GNN, MPNNs

1. Applications of subgraph embeddings

DeepWalk-随机游走+Skip-Gram-[2014SIGKDD]

Perozzi, B.; Al-Rfou, R.; and Skiena, S. 2014.

Deepwalk: Online learning of social representations.

In Proceedings of the 20th ACM SIGKDD international conference on Knowledge discovery and data mining

只适用于不加权的图

算法1: $DeepWalk(G, w, d, \gamma, t)$

Input:

graph $G(V;E)$ window size w embedding size d walks per vertex γ walk length t **Output:** matrix of vertex representations $\Phi \in R^{|V| \times d}$ 1: Initialization: Sample Φ from $U^{|V| \times d}$ 2: Build a binary Tree T from V 3: for $i = 0$ to γ do 4: $O = \text{Shuffle}(V)$ 5: for each $v_i \in O$ do 6: $W_{v_i} = \text{RandomWalk}(G, v_i, t)$ 7: $\text{SkipGram}(\Phi, W_{v_i}, w)$ 8: end for 9: end for

算法2: $\text{SkipGram}(\Phi, W_{v_i}, w)$

1: for each $v_j \in W_{v_i}$ do 2: for each $u_k \in W_{v_i}[j-w : j+w]$ do 3: $J(\Phi) = -\log \text{Pr}(u_k | \Phi(v_j))$ 4: $\Phi = \Phi - \alpha * \frac{\partial J}{\partial \Phi}$ (SGD) 5: end for 6: end for

Computing the partition function (normalization factor) is expensive, so instead we will factorize the conditional probability using Hierarchical softmax. We assign the vertices to the leaves of a binary tree, turning the prediction problem into maximizing the probability of a specific path in the hierarchy. If the path to vertex u_k is identified by a sequence of tree nodes $(b_0, b_1, \dots, b_{\lceil \log |V| \rceil})$, ($b_0 = \text{root}$, $b_{\lceil \log |V| \rceil} = u_k$) then

$$\text{Pr}(u_k | \Phi(v_j)) = \prod_{l=1}^{\lceil \log |V| \rceil} \text{Pr}(b_l | \Phi(v_j))$$

$\text{Pr}(b_l | \Phi(v_j))$ could be modeled by a binary classifier that is assigned to the parent of the node b_l as

$$\text{Pr}(b_l | \Phi(v_j)) = 1 / (1 + e^{-\Phi(v_j) \cdot \Psi(b_l)})$$

where $\Psi(b_l) \in R^d$ is the representation assigned to tree node b_l 's parent

实验：多标签分类

参数敏感性分析

Line-1阶&2阶相似度-[2015]

Tang J, Qu M, Wang M, et al.

Line: Large-scale information network embedding

[C]//Proceedings of the 24th International Conference on World Wide Web. International World Wide Web Conferences Steering Committee, 2015: 1067-1077.

以前的方法只适用于小型网络。

Although a few very recent studies approach the embedding of large-scale networks, these methods either use an indirect approach that is not designed for networks or lack a clear objective function tailored for network embedding.

LINE, which is able to scale to very large, arbitrary types of networks: undirected, directed and/or weighted

The model optimizes an objective which preserves both the local and global network structures

两种相似性：

first-order：两个点之间有很强的联系(比如边的权重很大)。

second-order：两个点有很多相同的邻居

they should be represented closely to each other in the embedded space

LINE with First-order Proximity

对于无向边 (i, j) , v_i 和 v_j 的联合概率是 $p_1(v_i, v_j) = \frac{1}{1 + \exp(-\vec{u}_i^T \cdot \vec{u}_j)}$

相应的经验概率是： $\hat{p}_1(i, j) = \frac{w_{ij}}{W}$, $W = \sum_{(i, j) \in E} w_{ij}$

最小化两个概率分布之间的距离，用KL散度距离，因此目标函数是：
 $O_1 = d(\hat{p}_1(\cdot, \cdot), p_1(\cdot, \cdot)) = -\sum_{(i, j) \in E} w_{ij} \log p_1(v_i, v_j)$

LINE with Second-order Proximity

无向和有向图都适用

每个点被看作是"context"，在"context"上有相似分布的结点视作相似。

每个点有两个角色：the vertex itself and a specific "context" of other vertices，对应两个向量 \vec{u}_i 和 \vec{u}_i'

对于每个有向边 (i, j) , 定义由点 v_i 生成的"context" v_j 的概率为： $p_2(v_j | v_i) = \frac{\exp(\vec{u}_j^T \cdot \vec{u}_i')}{\sum_{k=1}^{|V|} \exp(\vec{u}_k^T \cdot \vec{u}_i')}$

where $|V|$ is the number of vertices or "contexts."

最小化目标函数:

$$O_2 = \sum_{i \in V} \lambda_i d(\hat{p}_2(\cdot|v_i), p_2(\cdot|v_i)) = - \sum_{(i,j) \in E} w_{ij} \log p_2(v_j|v_i)$$

其中 λ_i 是 the prestige 声望 of vertex i in the network, which can be measured by the degree or estimated through algorithms such as PageRank, 这里取 $d_i = \sum_{k \in N(i)} w_{ik}$ 即点 i 的出度, $\hat{p}_2(v_j|v_i) = \frac{w_{ij}}{d_i}$

分别训练两个LINE模型, 然后将两个模型的每个点的向量拼接起来, 也可以联合训练两个目标函数(future work)

优化目标函数 O_2 is computationally expensive, 采用负采样方法。对每条边 (i,j) , 目标函数为:

$$\log \sigma(\vec{u}_j^{IT} \cdot \vec{u}_i) + \sum_{i=1}^K E_{v_n \sim P_n(v)} [\log \sigma(\vec{u}_n^{IT} \cdot \vec{u}_i)]$$

set $P_n(v) \propto d_v^{3/4}$, d_v 是点 v 的出度

用 asynchronous stochastic gradient algorithm (ASGD) 优化上述方程。

In each step, the ASGD algorithm samples a mini-batch of edges and then updates the model parameters.

梯度会由权重乘积而来, 而权重的方差很大, 会导致难以找到合适的学习率。

用 the alias table method 根据边的权重进行采样, 将采样到的边视为 binary edges.

对于目标函数 O_1 也采用负采样的方法, 将上式中的 \vec{u}_j^{IT} 换成 \vec{u}_j^T

实验: a language network, two social networks, and two citation networks

second-order proximity suffers when the network is extremely sparse, and it outperforms first-order proximity when there are sufficient nodes in the neighborhood of a node

second-order proximity does not work well for nodes with a low degree

node2vec-捕捉结构相似性和趋同性-[2016]

Grover A, Leskovec J.

node2vec: Scalable feature learning for networks

[C]//Proceedings of the 22nd ACM SIGKDD international conference on Knowledge discovery and data mining. ACM, 2016

a semi-supervised algorithm for scalable feature learning in networks

简单来说就是将原有社交网络中的图结构, 表达成特征向量矩阵, 每一个node (可以是人、物品、内容等) 表示成一个特征向量, 用向量与向量之间的矩阵运算来得到相互的关系。(如向量均值, Hadamard积, Weighted-L1, Weighted-L2)

there is no clear winning sampling strategy that works across all networks and all prediction tasks. This is a major shortcoming of prior work which fail to offer any flexibility in sampling of nodes from a network.

Our algorithm node2vec overcomes this limitation by designing a flexible objective that is not tied to a particular sampling strategy and provides parameters to tune the explored search space.

Algorithm 1 The node2vec algorithm. LearnFeatures (Graph $G = (V, E, W)$, Dimensions d , Walks per node r , Walk length l , Context size k , Return p , In-out q) $\pi = \text{PreprocessModifiedWeights}(G, p, q)$ $G' = (V, E, \pi)$ Initialize walks to Empty for $\text{iter} = 1$ to r do for all nodes $u \in V$ do $\text{walk} = \text{node2vecWalk}(G', u, l)$ Append walk to walks $f = \text{StochasticGradientDescent}(k, d, \text{walks})$ return f **node2vecWalk** (Graph $G' = (V, E, \pi)$, Start node u , Length l) Initialize walk to $[u]$ for $\text{walk_iter} = 1$ to l do $\text{curr} = \text{walk}[-1]$ $V_{\text{curr}} = \text{GetNeighbors}(\text{curr}, G')$ $s = \text{AliasSample}(V_{\text{curr}}, \pi)$ Append s to walk return walk

RandomWalks:

$$P(c_i = x | c_{i-1} = v) = \frac{\pi_{vx}}{Z} \text{ if } (v, x) \in E \text{ otherwise } 0$$

$$\pi_{vx} = \alpha_{pq}(t, x) \cdot w_{vx}$$

$$\alpha_{pq}(t, x) = \frac{1}{p} \text{ if } d_{tx} = 0$$

$$\alpha_{pq}(t, x) = 1 \text{ if } d_{tx} = 1$$

$$\alpha_{pq}(t, x) = \frac{1}{q} \text{ if } d_{tx} = 2$$

d_{tx} denotes the shortest path distance between nodes t and x . t 是当前点的上一个点, v 是当前点, x 是 v 的邻居结点

Parameter p controls the likelihood of immediately revisiting a node in the walk.

Setting it to a high value ($> \max(q, 1)$) ensures that we are less likely to sample an already visited node in the following two steps (unless the next node in the walk had no other neighbor).

if $q > 1$, 倾向于BFS广度优先搜索, if $q < 1$, 倾向于DFS深度优先搜索

极大似然优化, 目标函数是:

$$\max_f \sum_{u \in V} \log \Pr(N_S(u) | f(u))$$

$$\Pr(N_S(u) | f(u)) = \prod_{n_i \in N_S(u)} \Pr(n_i | f(u))$$

$$\Pr(n_i | f(u)) = \frac{\exp(f(n_i) \cdot f(u))}{\sum_{v \in V} \exp(f(v) \cdot f(u))}$$

因此目标函数变为:

$$\max_f \sum_{u \in V} [-\log Z_u + \sum_{n_i \in N_S(u)} f(n_i) \cdot f(u)]$$

$$Z_u = \sum_{v \in V} \exp(f(u) \cdot f(v))$$

可以通过负采样来优化分母的计算量,

用随机梯度下降法优化上述目标函数

实验: 聚类, 多标签的分类, 链接预测

参数敏感分析, 扰动分析, 可扩展性

- 用学到的向量去做分类任务的特征, 结果比其他方法好很多, 并且这种方法很鲁棒! 即使缺少边也没问题。
- 可扩展到大规模 node!

DKRL-基于文本+结构的embedding-[2016AAAI]

R Xie, Z Liu, J Jia, ... - 2016

Representation Learning of Knowledge Graphs with Entity Descriptions

AAAI

考虑实体描述的知识表示学习模型(description-embodied knowledge representation learning, DKRL)提出在知识表示学习中考虑Freebase等知识库中提供的实体描述文本信息。在文本表示方面, DKRL考虑了2种模型:一种是CBOW,将文本中的词向量简单相加作为文本表示;一种是卷积神经网络(convolutional neural network,CNN),能够考虑文本中的词序信息。

优势:除了能够提升实体表示的区分能力外,还能实现对新实体的表示。当新出现一个未曾在知识库中的实体时, DKRL可以根据它的简短描述产生它的实体表示,用于知识图谱补全等任务。

Description-Embodied Knowledge Representation Learning

每个头结点和尾结点有两个向量,分别是基于结构s的向量和基于文本描述d的向量

energy function:

$$E = E_S + E_D, E_D = E_{DD} + E_{DS} + E_{SD}$$

$$E_{DD} = ||h_d + r - t_d||, E_{DS} = ||h_d + r - t_s||, E_{SD} = ||h_s + r - t_d||$$

two encoders to build description-based representations

Continuous Bag-of-words Encoder

为每个实体选择文本描述中的top n个关键词(可以用TF-IDF进行排序)作为输入,将关键词的embeddings加起来作为实体的embedding,用来最小化 E_D

Convolutional Neural Network Encoder

5层,实体的预处理后的文本描述作为输入,输出该实体基于文本描述的embedding.

预处理:去停用词,标记文本描述中的短语,将它们作为词,每个词有一个word embedding,作为CNN的输入。

structure2vec-[2016]

Dai H, Dai B, Song L.

Discriminative embeddings of latent variable models for structured data

[C]//International Conference on Machine Learning. 2016: 2702-2711.

structure2vec, an effective and scalable approach for structured data representation based on the idea of embedding latent variable models into feature spaces, and learning such feature spaces using discriminative information 区分性信息.

Structured data, such as sequences, trees and graphs

GraphSAGE-有效生成新结点embedding-[2017]

采样邻居结点，通过累加器得到结点的embedding，对于unseen nodes可以有效地生成embeddings

累加器有三种：Mean aggregator, LSTM aggregator, Pooling aggregator

Hamilton W, Ying Z, Leskovec J.

Inductive representation learning on large graphs

[C]//Advances in Neural Information Processing Systems. 2017

APP-非对称相似度-阿里-[2017AAAI]

Scalable Graph Embedding for Asymmetric Proximity C Zhou, Y Liu, X Liu, ... - 2017

asymmetric proximity preserving(APP)

we propose an asymmetric proximity preserving(APP) graph embedding method via random walk with restart, which captures both asymmetric and high-order similarities between node pairs. We give theoretical analysis that our method implicitly preserves the Rooted PageRank score for any two vertices.

考虑节点间的非对称相似度，每个节点训练出两个向量：头向量和尾向量。两个节点的相似性用两个节点的头向量和尾向量的内积表示

Higher order proximity :

1.SimRank

SimRank 是一种基于图的拓扑结构来衡量图中任意两个点的相似程度的方法。在基于链接的相似性度量领域中 SimRank被认为与PageRank在信息检索领域具有同样重要的地位。

如果两个点在图中的邻域比较相似（有很多相似邻居），则这两个点也应该比较相似。即两个点是否相似，由他们的邻居是否相似来决定。而他们的邻居是否相似又由他们邻居的邻居的相似性决定。

跟pagerank类似，这也是一个迭代的定义。即通过迭代的方式计算两个点之间的相似度，最终取收敛的相似度。

如果两个节点相同，则相似度是1。如果两个节点不同，那他们的相似度就等于他们两个所有一步邻居的两两相似度的均值，再乘以衰减系数cc。

SimRank的特点：完全基于结构信息，且可以计算图中任意两个节点间的相似度。

Jeh, G., and Widom, J. 2002.

Simrank: a measure of structural-context similarity.

In International Conference on Knowledge Discovery and Data Mining, 538–543. ACM.

2. Rooted PageRank

PageRank的计算充分利用了两个假设：数量假设和质量假设。步骤如下：**1) 在初始阶段：**网页通过链接关系构建起Web图，每个页面设置相同的PageRank值，通过若干轮的计算，会得到每个页面所获得的最终PageRank值。随着每一轮的计算进行，网页当前的PageRank值会不断得到更新。

2) 在一轮中更新页面PageRank得分的计算方法：在一轮更新页面PageRank得分的计算中，每个页面将其当前的PageRank值平均分配到本页面包含的出链上，这样每个链接即获得了相应的权值。而每个页面将所有指向本页面的入链所传入的权值求和，即可得到新的PageRank得分。当每个页面都获得了更新后的PageRank值，就完成了一轮PageRank计算。

PageRank的迭代公式： $R = q \times P * R + (1 - q) * e / N$, e 是单位向量

主题敏感PageRank中： $R = q \times P * R + (1 - q) * s / N$, s 是这样一个向量：对于某topic的 s ，如果网页 k 在此topic中，则 s 中第 k 个元素为1，否则为0。对于每个topic都有一个不同的 s ，而 $|s|$ 表示 s 中1的数量。每个网页归到一个topic。

在PageRank中 e/N 是一个均匀分布，而在PPR中则根据用户的preference指定权重，例如用户指定了10个页面，则可以设置这10个页面对应的权重均为1/10，其余均为0。

一般来说用户会对某些领域感兴趣，同时，当浏览某个页面时，这个页面也是与某个主题相关的（比如体育报道或者娱乐新闻），所以，当用户看完当前页面，希望跳转时，更倾向于点击和当前页面主题类似的链接，即主题敏感PageRank是将用户兴趣、页面主题以及链接所指向网页与当前网页主题的相似程度综合考虑而建立的模型。

PageRank是全局性的网页重要性衡量标准，每个网页会根据链接情况，被赋予一个唯一的PageRank分值。主题敏感PageRank在此点有所不同，该算法引入16种主题类型，对于某个网页来说，对应某个主题类型都有相应的PageRank分值，即每个网页会被赋予16个主题相关PageRank分值。

Haveliwala, T. H. 2002.

Topic-sensitive pagerank.

In Proceedings of the 11th international conference on World Wide Web

3. Katz

https://en.wikipedia.org/wiki/Katz_centrality

两个节点之间所有路径的加权和

Katz, L. 1953.

A new status index derived from sociometric analysis.

Psychometrika

Asymmetric graph embedding

High-Order Proximity preserved Embedding(HOPE for short)

we first derive a general formulation of a class of high-order proximity measurements, then apply generalized SVD to the general formulation, whose time complexity is linear with the size of graph.

Ou, M.; Cui, P.; Pei, J.; and Zhu, W. 2016.

Asymmetric transitivity preserving graph embedding.

In International Conference on Knowledge Discovery and Data Mining. ACM.

the Monte-Carlo End-Point sampling method

Let A denote the adjacency matrix of the web graph with normalized rows and $c \in (0, 1)$ the teleportation probability. In addition, let r be the so-called preference vector inducing a probability distribution over V . PageRank vector p is defined as the solution of the following equation

$$p = (1 - c) \cdot pA + c \cdot r$$

要么以概率 $c \cdot r$ 选择该网页，要么从别的网页跳到该网页来。

按上述式子迭代的话复杂度很高，所以一般用蒙特卡洛模拟。

To compute a rooted pagerank vector for v , the Monte Carlo approach randomly samples N independent paths started from v , with stopping probability of c . Then the rooted pagerank value can be approximated as, $ppr_v(u) = |PathEndsAt(u)|/N$

网页 v 对于主体 u 的 pagerank score

论文中用蒙特卡洛方法模拟 v 到达 u 的概率，用于目标函数中。

Fogaras, D.; R'acz, B.; Csalog'any, K.; and Sarl'os, T. 2005.

Towards scaling fully personalized pagerank: Algorithms, lower bounds, and experiments.

Internet Mathematics

DeepWalk

Perozzi, B.; Al-Rfou, R.; and Skiena, S. 2014.

Deepwalk: Online learning of social representations.

In Proceedings of the 20th ACM SIGKDD international conference on Knowledge discovery and data mining

Line

Tang, J.; Qu, M.; Wang, M.; Zhang, M.; Yan, J.; and Mei, Q. 2015

Line: Large-scale information network embedding.

In Proceedings of the 24th International Conference on World Wide Web

Node2Vec

Grover, A., and Leskovec, J. 2016.

node2vec: Scalable feature learning for networks.

In International Conference on Knowledge Discovery and Data Mining. ACM.

Common Neighbors (CNbrs for short)

最简单的基于局部信息的相似性方法。如果两个节点间的共同邻居结点越多，那么两者存在链接的可能性就越大。得分公式：

$$score(u, v) = |N(u) \cap N(v)|$$

Adamic Adar (Adar for short)

起初用于计算两个用户的个人主页的相似性。在计算两个用户个人主页时，首先要提取两主页的公共关键词，然后计算每个公共关键词的权重，最后对所有的公共关键词进行加权求和。关键词的权重与关键词出现的次数的倒数成反比。

$$\sum_{t \in N(u) \cap N(v)} \frac{1}{\log |N(t)|}$$

Jaccard Coefficient

利用两节点共同邻居的交集与并集个数之比，定义为两节点的相似度。

$$score(u, v) = |N(u) \cap N(v)| / |N(u) \cup N(v)|$$

HIN2Vec-异质结点和边-[2017]

Heterogeneous Information Network to Vector (HIN2Vec)

1527164573047

P5

Fu T, Lee W C, Lei Z. 2017

HIN2Vec: Explore Meta-paths in Heterogeneous Information Networks for Representation Learning

[C]//Proceedings of the 2017 ACM on Conference on Information and Knowledge Management. ACM

IGE-交互图的embedding-[2017]

Zhang Y, Xiong Y, Kong X, et al. 2017

Learning Node Embeddings in Interaction Graphs

[C]//Proceedings of the 2017 ACM on Conference on Information and Knowledge Management. ACM

IGE (Interaction Graph Embedding)

以前的节点嵌入方法都集中在static non-attributed graphs上，现实中attributed interaction graphs.它包含两类实体（如投资者/股票），以及edges of temporal interactions with attributes (e.g. transactions and reviews)

Our model consists of an attributes encoding network and two prediction networks. Three networks are trained simultaneously by back-propagation. Negative sampling method is applied in training prediction networks for acceleration. 将属性转化成固定长度的向量的编码网络（处理属性的异质性）；利用属性向量预测两个实体的暂时事件的子网络如 预测投资者事件的子网络；预测股票事件的子网络 s

Attributed interaction graph:

$G = (X, Y, E)$, X and Y are two disjoint sets of nodes, and E is the set of edges. Each edge $e \in E$ can be represented as a tuple $e = (x, y, t, a)$, where $x \in X$, $y \in Y$, t denotes the timestamp and $a = (a_1, \dots, a_m)$ is the tuple of heterogeneous attributes of the edge. $a_i \in A_i$ is an instance of the i -th attribute.

例如(investor, stock, time, buy_or_sell, price, amount)

Induced Edge List

Given an attributed interaction graph $G = (X, Y, E)$ and a source node $x \in X$, we say $s_x = (e_1, \dots, e_n)$ is an edge list induced by the source node x , if every edge in s_x is incident to x . Similarly, we can define a target node y 's induced edge list s_y .

A Simplified Model

Let $G = (X, Y, E)$ be a given interaction graph where edges are without attributes. Inspired by the Skip-gram model and Paragraph Vector, we formulate the embedding learning as a maximum likelihood problem.

$$L(\theta) = \alpha \sum_{x \in X} \log P(s_x; \theta) + (1 - \alpha) \sum_{y \in Y} \log P(s'_y; \theta)$$

$\alpha \in [0, 1]$ 是超参数，is used to make a trade-off between the importance of source nodes induced lists and target nodes induced lists.

θ 代表所有的模型参数

$$\log P(s_x; \theta) = \sum_i \frac{1}{Z_i} \sum_{j \neq i} e^{\frac{-|t_i - t_j|}{\tau}} \log P(y_j | x, y_i; \theta)$$

where $Z_i = \sum_{j \neq i} e^{\frac{-|t_i - t_j|}{\tau}}$ is a normalizing constant, and τ is a hyperparameter. If τ is small, weights will concentrate on the temporally close pairs. Conversely, the larger the τ is, the smoother the weights are. In this case, more long-term effects will be considered.

$$\log P(y_j = k | x, y_i; \theta) = \frac{\exp(U_X[k, :]^T v_x + U_Y[k, :]^T v_{y_i} + b)}{\sum_l \exp(U_X[l, :]^T v_x + U_Y[l, :]^T v_{y_i} + b)}$$

v_x, v_{y_i} 是 x 和 y_i 的 embeddings，计算下列式子代替前面的式子

$$\log P(s_x; \theta) = \sum_i \frac{1}{N(i)} \sum_{j \in N(i)} \log P(y_j | x, y_i; \theta)$$

where $N(i) = \{i_1, \dots, i_c\}$ is the “context” of i . c is a pre-chosen hyper-parameter indicating the length of context, and i_k is selected randomly with the probability proportional to $e^{\frac{-|t_i - t_j|}{\tau}}$. $N(i)$ is a multiset, which means it allows duplicates of elements.

Embedding Tensors

在不同的场景下有不同的embeddings，例如投资者在买卖股票时有不同的策略。所以结点的embeddings是一个tensor, $\mathcal{T} \in \Re^{V \times K \times D}$, where V is the size of nodes set and D corresponds to the number of tensor slices.

Given a tuple of attributes $a = (a_1, \dots, a_m)$, and an attributes encoder f , we can get an attribute vector $d = f(a) \in \Re^D$. Then we can compute attributed-gated embeddings as $E^d = \sum_{i=1}^D d_i \mathcal{T}[:, :, i]$.

However, fully 3-way tensors are not practical because of enormous size. It is a common way to factor the tensor by introducing three matrices $W^{fv} \in \Re^{F \times V}$, $W^{fd} \in \Re^{F \times D}$, $W^{fk} \in \Re^{F \times K}$, and re-represent E^d by the equation $E^d = (W^{fv})^T \cdot \text{diag}(W^{fd} d) \cdot W^{fk}$

where $\text{diag}(\cdot)$ denotes the matrix with its argument on the diagonal. These matrices are parametrized by a pre-chosen number of factors F . It can be seen as the embeddings conditioned on d , and we let $E^d = (W^{fv})^T W^{fk}$ denote unconditional embeddings.

IGE: A Multiplicative Neural Model

用 $P(y_j|x, y_i, a_i, a_j; \theta)$ 代替上面式子中的 $P(y_j|x, y_i; \theta)$

$$\log P(y_j = k|x, y_i, a_i, a_j; \theta) = \frac{\exp(U_X[k, :]^T v_x + U_Y^{d_j}[k, :]^T v_{y_i}^{d_j} + b)}{\sum_l \exp(U_X[l, :]^T v_x + U_Y^{d_j}[l, :]^T v_{y_i}^{d_j} + b)}$$

交替训练:

1. 选择一个 x , 选择 s_x 的两条边, 根据 $\log P(y_j|x, y_i, a_i, a_j; \theta^{(t-1)})$ 计算出 $\Delta\theta$, 更新 $\theta^{(t)} = \theta^{(t-1)} + \alpha \lambda \Delta\theta$
2. 选择一个 y , 选择 s_y 的两条边, 根据 $\log P(x_t|y, x_s, a_s, a_t; \theta^{(t-1)})$ 计算出 $\Delta\theta$, 更新 $\theta^{(t)} = \theta^{(t-1)} + (1 - \alpha) \lambda \Delta\theta$

Know-Evolve-[2017]

Trivedi R, Dai H, Wang Y, et al.

Know-evolve: Deep temporal reasoning for dynamic knowledge graphs

[C]//International Conference on Machine Learning. 2017: 3462-3471.

Know-Evolve, a novel deep evolutionary knowledge network that learns non-linearly evolving entity representations over time.

四元组 (e_s, r, e_o, t) , 其中 $e^s, e^o \in \{1, \dots, n^e\}$ (实体), $e^s \neq e^o, r \in \{1, \dots, n_r\}, r \in R^+$

SSP-三元组+文本学习-[2017AAAI]

Xiao H, Huang M, Meng L, et al.

SSP: Semantic Space Projection for Knowledge Graph Embedding with Text Descriptions

[C]//AAAI. 2017, 17: 3104-3110.

semantic space projection (SSP) model which jointly learns from the symbolic triples and textual descriptions

损失函数:

$$f_r(h, t) = -\lambda \|e - s^T es\|_2^2 + \|e\|_2^2, \quad e = h + r - t$$

the component of the loss in the normal vector direction is $(s^T es)$, then the other orthogonal one, that is inside the hyperplane, is $(e - s^T es)$.

超平面的法向量 $s = S(s_h, s_t)$

将实体描述看作文档，采用主题模型，得到文档的主题分布作为实体的语义向量s，如(0.1, 0.9, 0.0)表示该实体的主题应该是2

$$S(s_h, s_t) = \frac{s_h + s_t}{\|s_h + s_t\|_2^2}$$

Standard setting: 给定预训练好的语义向量，模型中固定它们然后优化其它参数

Joint setting: 同时实施主题模型和embedding模型，此时三元组也会给文本语义带来正向影响。

Total Loss:

$$L = L_{embed} + \mu L_{topic}$$

$$L_{embed} = \sum_{(h,r,t) \in \Delta, (h',r',t') \in \Delta'} [f_{r'}(h', t') - f_r(h, t) + \gamma]_+$$

$$L_{topic} = \sum_{e \in E, w \in D_e} (C_{e,w} - s_e^T w)^2$$

where E is the set of entities, and D_e is the set of words in the description of entity e. $C_{e,w}$ is the times of the word w occurring in the description e. s_e is the semantic vector of entity e and w is the topic distribution of word w. Similarly, SGD is applied in the optimization.

实验: 知识图谱补全, 实体预测

SaC2Vec-基于结构+内容的embedding-[2018]

Sambaran Bandyopadhyay, Harsh Kara, Anirban Biswas, M N Murty, 2018.4.27

SaC2Vec: Information Network Representation with Structure and Content

Sac2Vec(structure and content to vector), a network representation technique using structure and content. It is a multi-layered graph approach which uses a random walk to generate the node embedding.

$G = (V, E, W, F)$, F can be considered as the content matrix. 其中 f_i 是 the word vector (content) associated with the node

F 可以用词袋模型(bag-of-words)表示, 矩阵的每一行是相应结点的文本内容的tf-idf向量, 所以 F 的维度是 $n \times d$, d 是语料中的unique words的数量。(预处理之后)

给定输入网络 $G=(V,E,W,F)$,分为两层:

Structure Layer: 这一层是 $G_s = (V, E_s, W_s)$, $E_s = E, W_s = W$

Content Layer: 这一层是 有向图 $G_c = (V, E_c, W_c)$, $w_{ij}^c = \text{Cosine}(F_i, F_j)$, 但是只保留 $top \theta \times \lceil avg_s \rceil$ 的出去的边, 其中 θ 是一个正整数, avg_s 是 Structure layer 的结点的平均出度, $|E|/n$ (有向), $2 \times |E|/n$ (无向)

Convex Sum of Embeddings: CSoE

用 node2vec 为结构层和内容层独立地生成 embeddings, 然后取两个 embeddings 的凸组合 (维度一致时才能用)

$$e_{convex}^i = \alpha e_s^i + (1 - \alpha) e_c^i, \quad \forall i \in \{1, 2, \dots, n\} \text{ and } 0 \leq \alpha \leq 1$$

Appended Embeddings: AE

将两个向量拼接起来, 维度不一样也可以

$$e_{appended}^i = [e_s^i || e_c^i], \quad e_s^i \in R^{K_s}, e_c^i \in R^{K_c}, e_{appended}^i \in R^{K_s + K_c}, \forall i \in \{1, 2, \dots, n\}$$

SaC2Vec Model

为 structure layer 的结点 v_i^s 定义 Γ_i^s 如下: content layer 的 Γ_i^c 类似

$$\Gamma_i^s = \{(v_i^s, v_j^s) \in E_s | w_{v_i^s, v_j^s}^s \geq \frac{1}{|E_s|} \sum_{e' \in E_s} w_{e'}^s, v_j^s \in V\}$$

是 i 的那些权重大于该层平均边权重的出边的集合。

两层相同结点之间的权重:

$$w_i^{sc} = \ln(e + |\Gamma_i^s|), w_i^{cs} = \ln(e + |\Gamma_i^c|)$$

random walk

当前所在结点是 v_i , 要么是结构层, 要么是内容层, 下一步走到哪一层呢? 我们的目标是 to move to a layer which is more informative in some sense at node v_i .

$$p(v_i^s | v_i) = \frac{w_i^{cs}}{w_i^{sc} + w_i^{cs}}$$

$$p(v_i^c | v_i) = 1 - p(v_i^s | v_i) = \frac{w_i^{sc}}{w_i^{sc} + w_i^{cs}}$$

考虑第一个式子, w_i^{sc} 越大, 结点 v_i^s 的那些权重大于结构层的相对高的权重的出边越多, 此时 random walk 如果在结构层, 在走下一步时会有很多选择。如果 w_i^{sc} 很小, random walk 处在结构层的话, 下一步选择会很少, 此时的选择会更有信息丰富性, 并且 less random. 所以, 当 w_i^{sc} 很大时, 倾向于选择 content layer, 当 w_i^{cs} 很大时, 倾向于选择 structure layer. 一旦选择了某一层, 在走下一步时就不用考虑另一层了。

Algorithm 1 SaC2Vec - Structure and Content to Vector 1: Input: The network $G = (V, E, W, F)$, K : Dimension of the embedding space where $K \ll \min(n, d)$ (d is distinct words 的数量), r : Number of time to start random walk from each vertex, l : Length of each random walk 2: Output: The node embeddings of the network G 3: Generate the structure layer and the content layer 4: Add the edges between the layers with weights to generate the multi-layered network 5: Corpus = [] . Initialize the corpus 6: for $iter \in \{1, 2, \dots, r\}$ do 7: for $v \in V$ do 8: select the node v as the starting node of the random walk 9: Walk = [v] . Initialize the Walk(sequence of nodes) 10: for $walkIter \in \{1, 2, \dots, l\}$ do 11: Select the layer to move next with probabilities 12: Move 1 step using node2vec to find the next node v_i 13: Append v_i to Walk 14: end for 15: Append Walk to Corpus 16: end for 17: end for 18: Find the node embeddings by running language model on Corpus(SkipGram模型, 最大化给定该结点的向量表示出现context nodes的概率)

实验: node classification, node clustering and network visualization

It means SaC2Vec is able to understand the bad quality of the content layer during the learning process and embeddings were learnt mostly from the structure layer.

feature propagation-特征前向传播+edge2vec-[2018]

We study feature propagation on graph, an inference process involved in graph representation learning tasks。 however few works studied the convergence of feature propagation

Xiang B, Liu Z, Zhou J, et al.

Feature Propagation on Graph: A New Perspective to Graph Representation Learning

[J]. arXiv preprint arXiv:1804.06111, 2018.

Word Embedding的稳定性的影响因素-[2018]

Wendlandt L, Kummerfeld J K, Mihalcea R.

Factors Influencing the Surprising Instability of Word Embeddings

[J]. arXiv preprint arXiv:1804.09692, 2018.

We define stability as the percent overlap between nearest neighbors in an embedding space

两个embedding空间, 找到同一个词的最近的十个邻居, 将两个邻居列表的重叠作为词W的stability。

多个embedding空间时, considering the average overlap between two sets of embedding spaces

建立回归模型:

自变量是(1) properties related to the word itself; (2) properties of the data used to train the embeddings; and (3) properties of the algorithm used to construct these embeddings.

岭回归，最小化下列函数：

$$L(w) = \frac{1}{2} \sum_{n=1}^N (y_n - w^T x_n)^2 + \frac{\lambda}{2} \|w\|^2$$

we set 正则化常数 $\lambda = 1$

Word Properties:

Primary part-of-speech: Adjective Secondary part-of-speech: Noun

Parts-of-speech: 2

WordNet senses: 3

Syllables: 5 音节

Data Properties

Raw frequency in corpus A: 106 Raw frequency in corpus B: 669 Diff. in raw frequency: 563 Vocab. size of corpus A: 10,508 Vocab. size of corpus B: 43,888 Diff. in vocab. size: 33,380 Overlap in corpora vocab.: 17% Domains present: Arts, Europarl Do the domains match?: False Training position in A: 1.02% Training position in B: 0.15% Diff. in training position: 0.86% **Algorithm Properties** Algorithms present: word2vec, PPMI Do the algorithms match?: False Embedding dimension of A: 100 Embedding dimension of B: 100 Diff. in dimension: 0 Do the dimensions match?: True

we show that domain and part-of-speech are key factors of instability

In order to use the most stable embedding spaces for future tasks, we recommend either using GloVe or learning a good curriculum for word2vec training data. We also recommend using in-domain embeddings whenever possible.

Trans系列

TransE

将知识库中的关系看作实体间的某种平移向量，对于每个三元组(h,r,t)，TransE用关系r的向量 l_r ，作为头实体向量 l_h 和尾实体向量 l_t 之间的平移，也可以看作翻译。

对于每个三元组(h,r,t)，TransE希望 $l_h + l_r \approx l_t$

损失函数 $f_r(h,t) = |l_h + l_r - l_t|_{L1/L2}$ ，即向量 $l_h + l_r$ 和 l_t 的 L_1 或 L_2 距离。

在实际学习过程中，为了增强知识表示的区分能力，TransE采用最大间隔方法，定义了如下优化目标函数：

$$L = \sum_{(h,r,t) \in S} \sum_{(h',r',t') \in S^-} \max(0, f_r(h,t) + \gamma - f_{r'}(h',t'))$$

其中S是合法三元组的集合， S^- 为错误三元组的集合（将S中每个三元组的头实体、关系和尾实体其中之一随机替换成其他实体或关系得到）， γ 为合法三元组得分与错误三元组得分之间的间隔距离。

Translating embeddings for modeling multi-relational data

TransH模型

为了解决TransE模型在处理1-N，N-1，N-N复杂关系时的局限性，TransH模型提出让一个实体在不同的关系下拥有不同的表示。

对于关系r, TransH模型同时使用平移向量 l_r 和超平面的法向量 w_r 来表示它。对于一个三元组(h,r,t), TransH首先将头实体向量 l_h 和尾实体向量 l_t 沿法线 w_r 投影到关系r对应的超平面上，用 l_{h_r} 和 l_{t_r} 表示如下：

$$l_{h_r} = l_h - w_r^T l_h w_r$$

$$l_{t_r} = l_t - w_r^T l_t w_r$$

因此TransH定义了如下损失函数： $f_r(h, t) = |l_{h_r} + l_r - l_{t_r}|_{L1/L2}$

由于关系r可能存在无限个超平面，TransH简单地令 l_r 与 w_r 近似正交来选取某一个超平面。

Knowledge graph embedding by translating on hyperplanes

TransR / CTransR模型

虽然TransH模型使每个实体在不同关系下拥有了不同的表示，它仍然假设实体和关系处于相同的语义空间 R^d 中，这一定程度上限制了TransH的表示能力。TransR模型则认为，一个实体是多种属性的综合体，不同关系关注实体的不同属性。TransR认为不同的关系拥有不同的语义空间。对每个三元组，首先应将实体投影到对应的关系空间中，然后再建立从头实体到尾实体的翻译关系。

对于每一个关系r，TransR定义投影矩阵 $M_r \in R^{d \times k}$,将实体向量从实体空间投影到关系r的子空间，用 l_{h_r} 和 l_{t_r} 表示如下：

$$L_{h_r} = l_h M_r, L_{t_r} = l_t M_r$$

然后使 $l_{h_r} + l_r \approx l_{t_r}$.因此，TransR定义了如下损失函数：

$$f_r(h, t) = |l_{h_r} + l_r - l_{t_r}|_{L1/L2}$$

CTransR模型通过把关系r对应的实体对的向量差值 $l_h - l_t$ 进行聚类，将关系r细分为多个子关系 r_c ,CTransR模型为每一个子关系 r_c 分别学习向量表示，对于每个三元组(h,r,t),定义了如下损失函数：

$$f_r(h, t) = |l_{h_r} + l_{r_c} - l_{t_r}|_{L1/L2}$$

Learning entity and relation embeddings for knowledge graph completion

TransD模型

TransR缺点:

- 1.在同一个关系r下，头、尾实体共享相同的投影矩阵。然而，一个关系的头、尾实体的类型或属性可能差异巨大。
- 2.从实体空间到关系空间的投影是实体和关系之间的交互过程，因此TransR让投影矩阵仅与关系有关是不合理的。
- 3.与TransE和TransH相比，TransR由于引入了空间投影，使得TransR模型参数急剧增加，计算复杂度大大提高。

给定三元组(h,r,t), TransD模型设置了2个分别将头实体和尾实体投影到关系空间的投影矩阵 M_{rh} 和 M_{rt} ，具体定义如下:

$$M_{rh} = l_{r_p} l_{h_p} + I^{d \times k}, M_{rt} = l_{r_p} l_{t_p} + I^{d \times k}$$

这里 $l_{h_p}, l_{t_p} \in R^d, l_{r_p} \in R^k$, 下标p代表该向量为投影向量。显然， M_{rh} 和 M_{rt} 与实体和关系均相关。而且，利用2个投影向量构建投影矩阵，解决了原来TransR模型参数过多的问题。最后，TransD模型定义了如下损失函数:

$$f_r(h, t) = \|l_h M_{rh} + l_r - l_t M_{rt}\|_{L1/L2}$$

Knowledge graph embedding via dynamic mapping matrix

TranSparse模型

为了解决实体和关系的异质性(某些关系可能会与大量的实体有连接，而某些关系则可能仅仅与少量实体有连接)，TranSparse提出使用稀疏矩阵代替TransR模型中的稠密矩阵，其中矩阵 M_r 的稀疏度由关系r连接的实体对数量决定。这里头、尾实体共享同一个投影矩阵 M_r 。投影矩阵 $M_r(\theta_r)$ 的稀疏度 θ_r 定义如下: $\theta_r = 1 - (1 - \theta_{min}) N_r / N_r^*$

为了解决关系的不平衡性问题(在某些关系中，头实体和尾实体的种类和数量可能差别巨大)，TranSparse对于头实体和尾实体分别使用2个不同的投影矩阵。

Knowledge graph completion with adaptive sparse transfer matrix

TransA模型

Xiao等人认为TransE及其之后的扩展模型均存在2个重要问题: 1) 损失函数只采用 L_1 或 L_2 距离，灵活性不够; 2) 损失函数过于简单，实体和关系向量的每一维等同考虑。

TransA模型将损失函数中的距离度量改用马氏距离，并为每一维学习不同的权重。对于每个三元组(h,r,t), TransA模型定义了如下评分函数:

$$f_r(h, t) = (l_h + l_r - l_t)^T W_r (l_h + l_r - l_t), \text{ 其中 } W_r \text{ 为与关系 } r \text{ 相关的非负权值矩阵。}$$

TransA: An adaptive approach for knowledge graph embedding

TransG模型

TransG模型提出使用高斯混合模型描述头、尾实体之间的关系。该模型认为，一个关系会对应多种语义，每种语义用一个高斯分布来刻画。

TransG: A generative mixture model for knowledge graph embedding

PTransE

考虑关系路径，Path-based TransE

面临的挑战在于：

- 1) 并不是所有的实体间的关系路径都是可靠的。为此，PTransE提出Path-Constraint Resource Allocation图算法度量关系路径的可靠性。
- 2) PTransE需要建立关系路径的向量表示，参与从头实体到尾实体的翻译过程。这是典型的组合语义问题，需要对路径上所有关系的向量进行语义组合产生路径向量。PTransE尝试了3种代表性的语义组合操作，分别是相加、按位相乘和循环神经网络。相关数据实验表明，相加的组合操作效果最好。

TransNet

TransNet 假设头结点表示向量加上关系表示向量等于尾节点表示向量。其中，通过关键词抽取、命名实体识别等方式，对交互文本抽取出标签集合来表示关系。随后，通过深层自动编码器对标签集合进行压缩，来得到关系的表示向量。该模型能够有效地预测未标注的边上的标签集合，在社会关系抽取任务上取得了显著的提升。

Tu C C, Zhang Z Y, Liu Z Y, et al. TransNet: translation-based network representation learning for social relation extraction. In: Proceedings of International Joint Conference on Artificial Intelligence (IJCAI), Melbourne, 2017

KG2E模型

KG2E使用高斯分布来表示实体和关系，其中高斯分布的均值表示的是实体或关系在语义空间中的中心位置，而高斯分布的协方差则表示该实体或关系的不确定度。

Learning to represent knowledge graphs with Gaussian embedding