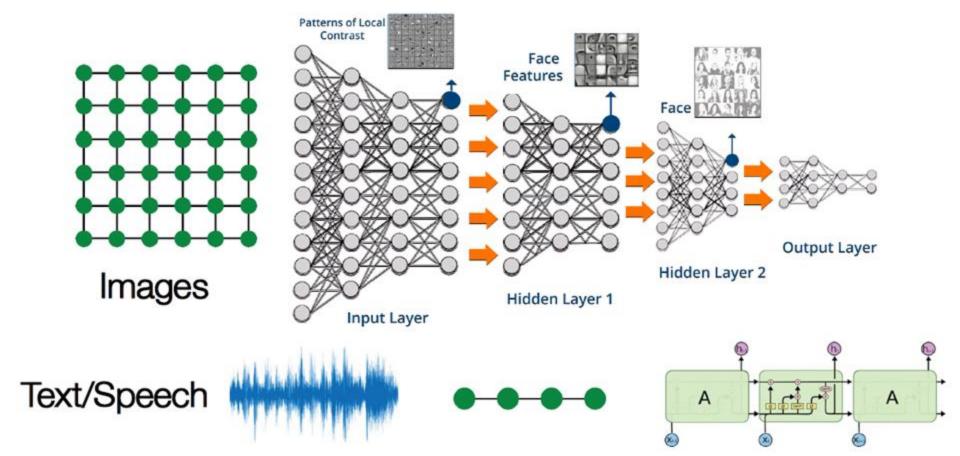
Graph Neural Network

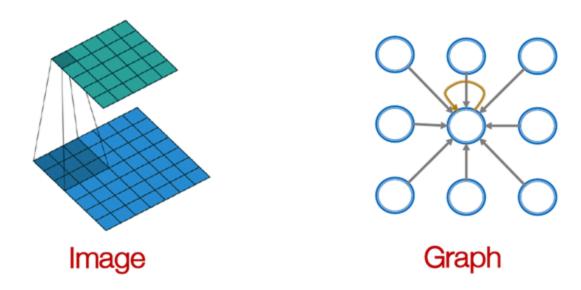
Seagate Technology

Modern DL Toolbox is Designed for Sequences & Grids



From Images to Graphs

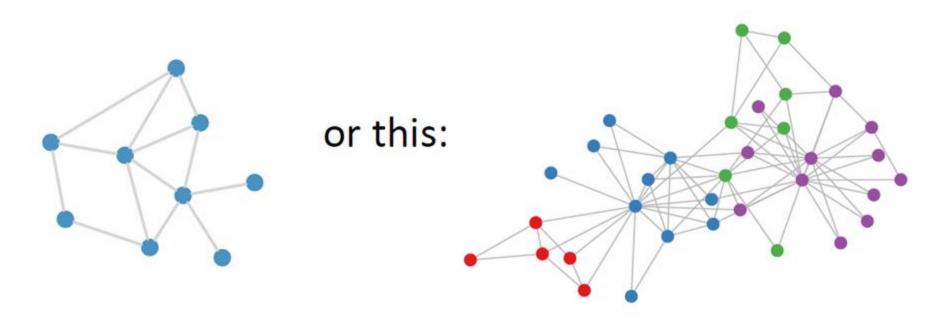
Single CNN layer with 3x3 filter:



Transform information at the neighbors and combine it:

- Transform "messages" h_i from neighbors: W_i h_i
- Add them up: $\sum_i W_i h_i$

How about these graphs?



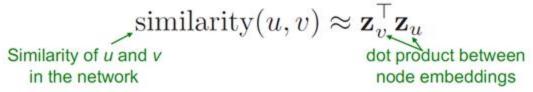
Biological networks, Medical networks, Social networks, Information networks, Knowledge graphs, Communication networks, Web graph

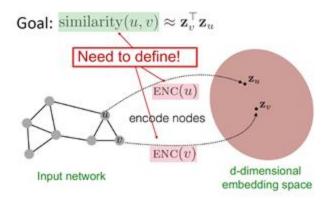
Problem Setup

Encoder: Map a node to a low-dimensional vector:
d-dimensional

$$\mathrm{ENC}(v) = \mathbf{z}_v$$
 embedding node in the input graph

Similarity function defines how relationships in the input network map to relationships in the embedding space:





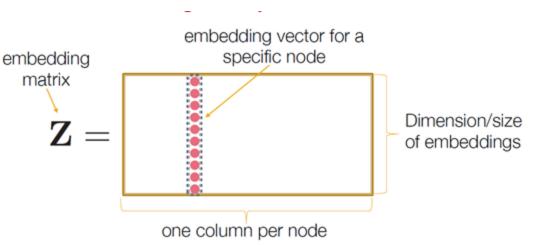
Shallow Encoding

 Simplest encoding approach: encoder is just an embedding-lookup

$$\text{ENC}(v) = \mathbf{Z}\mathbf{v}$$

 $\mathbf{Z} \in \mathbb{R}^{d imes |\mathcal{V}|}$ matrix, each column is a node embedding [what we learn!]

 $\mathbf{v} \in \mathbb{I}^{|\mathcal{V}|}$ indicator vector, all zeroes except a one in column indicating node v

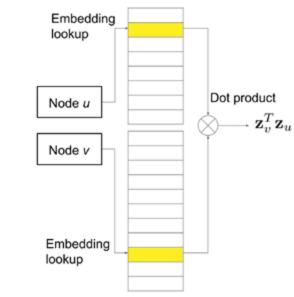


Each node is assigned to a unique embedding vector

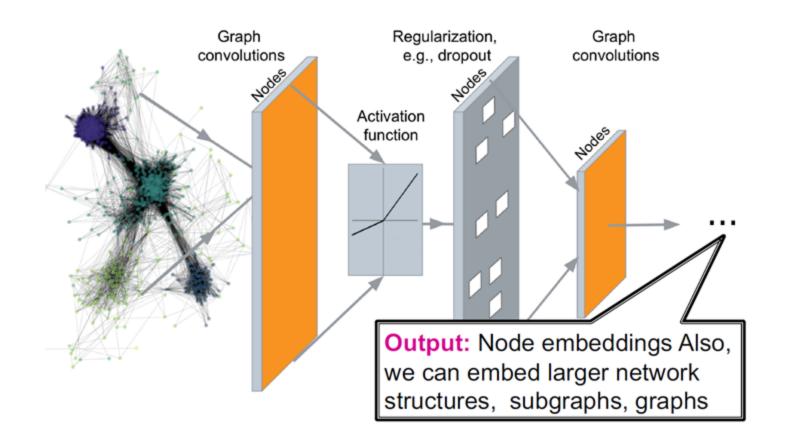
Many methods: DeepWalk, node2vec, TransE

Limitations of Shallow Encoding

- O(|V|) parameters are needed:
 - No sharing of parameters between nodes
 - Every node has its own unique embedding
- Inherently "transductive":
 - Cannot generate embeddings for nodes that are not seen during training
- Do not incorporate node features:
 - Many graphs have features that we can and should leverage

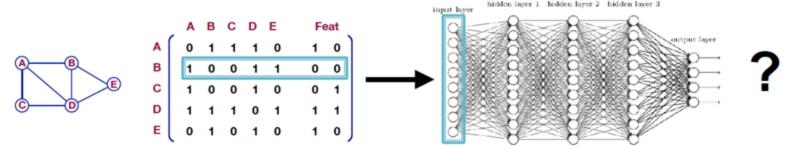


Deep Graph Encoders



A Naive Approach

- Join adjacency matrix and features
- Feed them into a deep neural net:



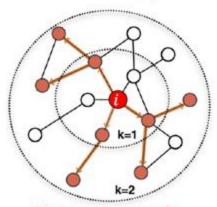
- Issues with this idea:
 - \bullet O(N) parameters
 - Not applicable to graphs of different sizes
 - Not invariant to node ordering

Setup

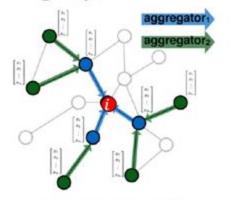
- Assume we have a graph G:
 - V is the vertex set
 - A is the adjacency matrix (assume binary)
 - $X \in \mathbb{R}^{m \times |V|}$ is a matrix of node features
 - Node features:
 - Social networks: User profile, User image
 - Biological networks: Gene expression profiles, gene functional information
 - No features:
 - Indicator vectors (one-hot encoding of a node)
 - Vector of constant 1: [1, 1, ..., 1]

Message Passing of Nodes

Idea: Node's neighborhood defines a computation graph



Determine node computation graph



Propagate and transform information

Advantages:

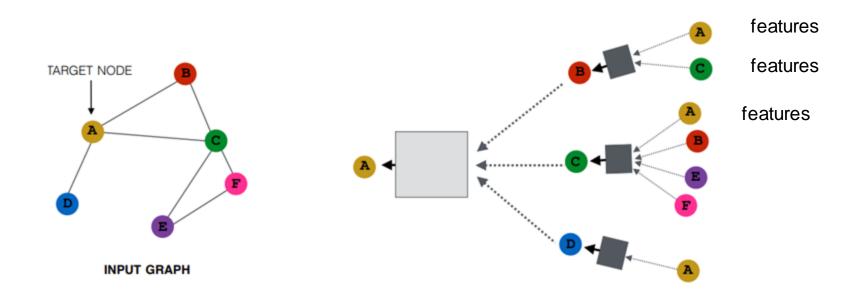
Capturing the Structure

Utilizing feature information

Learn how to propagate information across the graph to compute node features

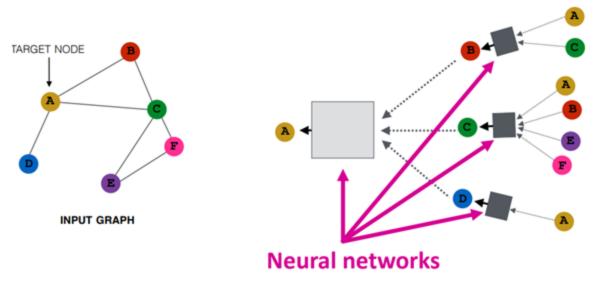
Aggregate Neighbors

 Key idea: Generate node embeddings based on local network neighborhoods



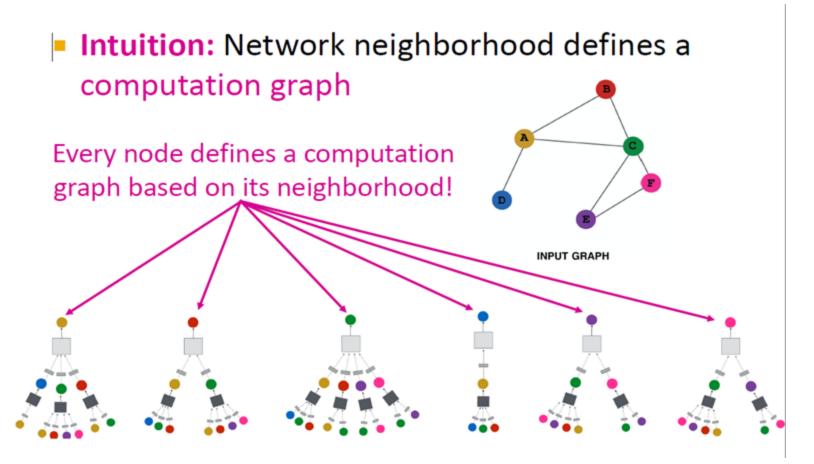
Shared Params of Neural Networks

 Intuition: Nodes aggregate information from their neighbors using neural networks

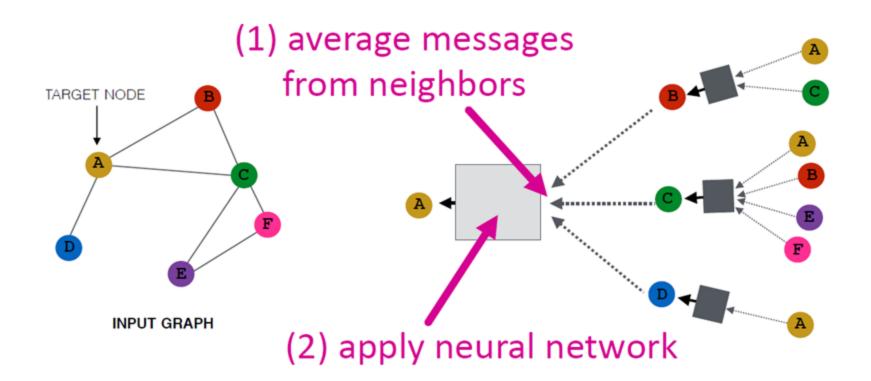


Order invariant

Aggregate Neighbors

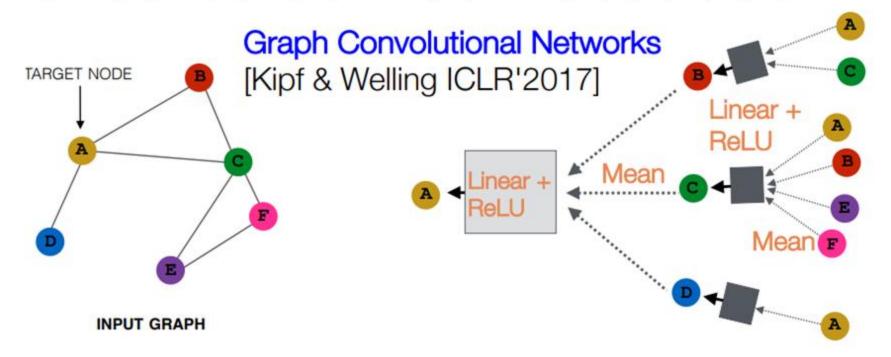


 Basic approach: Average information from neighbors and apply a neural network

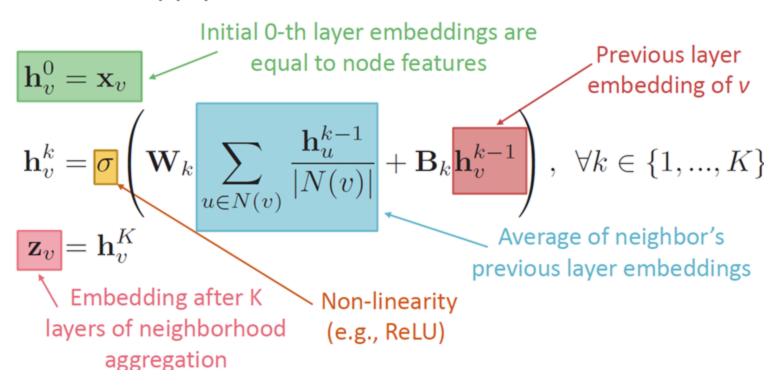


Many model variants have been proposed with difference choice of neural networks.

Scarselli et al., 2009b; Battaglia et al., 2016; Defferrard et al., 2016; Duvenaud et al., 2015; Hamilton et al., 2017a; Kearnes et al., 2016; Kipf & Welling, 2017; Lei et al., 2017; Li et al., 2016; Velickovic et al., 2018; Verma & Zhang, 2018; Ying et al., 2018; Zhang et al., 2018



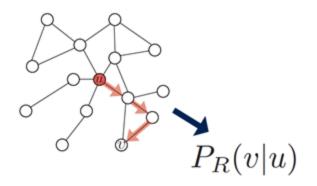
 Basic approach: Average neighbor messages and apply a neural network

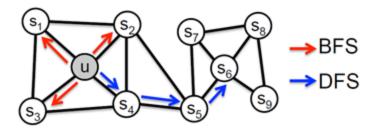


Unsupervised Training

Similar nodes should have similar embeddings

Can define a loss function based on results from node2vec, deepwalk, struc2vec



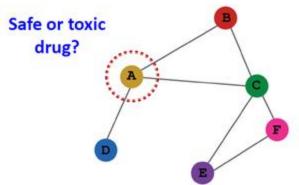


Random Walk

Node2vec: biased random walk

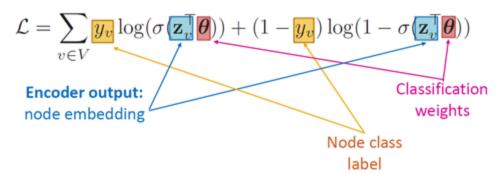
Supervised Training

Node Classification



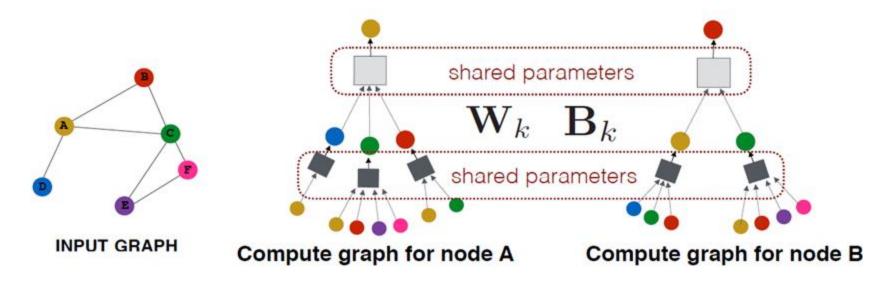
E.g., a drug-drug interaction network

Directly train the model for a supervised task (e.g., node classification)



Inductive Capability

- The same aggregation parameters are shared for all nodes:
 - The number of model parameters is sublinear in |V| and we can generalize to unseen nodes!



Application: Pinterest

Human curated collection of pins



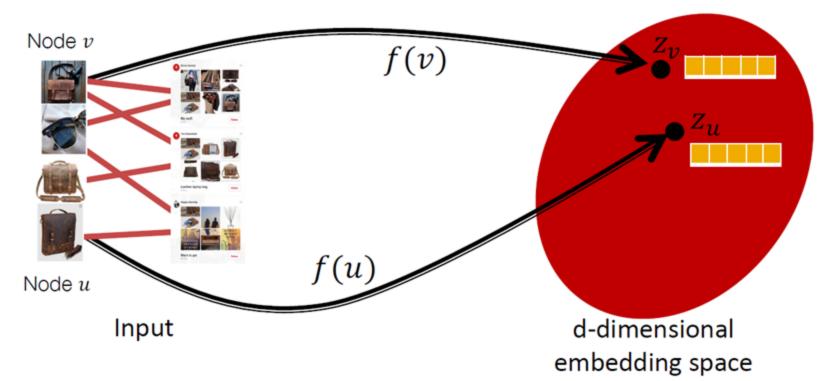
Board: A collection of ideas (pins having something in common)

- PinSage graph convolutional network:
 - Goal: Generate embeddings for nodes (e.g., Pins/images) in a web-scale Pinterest graph containing billions of objects
 - Key Idea: Borrow information from nearby nodes
 - E.g., bed rail Pin might look like a garden fence, but gates and beds are rarely adjacent in the graph





- Pin embeddings are essential to various tasks like recommendation of Pins, classification, clustering, ranking
 - Services like "Related Pins", "Search", "Shopping", "Ads"



Goal: Map nodes to d-dimensional embeddings such that nodes that are related are embedded close together

Task: Recommend related pins to users



Task: Learn node embeddings z_i such that $d(z_{cake1}, z_{cake2})$ $< d(z_{cake1}, z_{sweater})$

Challenges:

- Massive size: 3 billion nodes, 20 billion edges
- Heterogeneous data: Rich image and text features

Goal: Identify target pin among 3B pins

- Issue: Need to learn with resolution of 100 vs. 3B
- Idea: Use harder and harder negative samples
- Include more and more hard negative samples for each epoch



Source pin

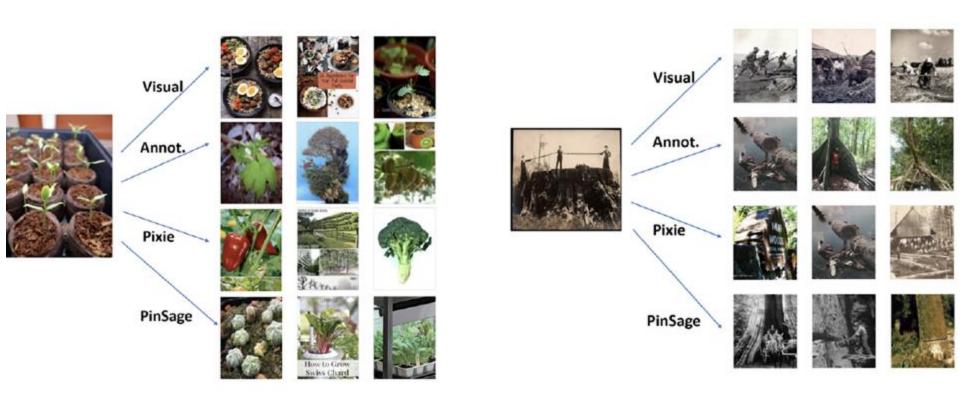


Positive





Easy negative Hard negative



Pixie: Random Walk (Dksombatchai et al., 2018)

Further Reading

Tutorials and overviews:

- Relational inductive biases and graph networks (Battaglia et al., 2018)
- Representation learning on graphs: Methods and applications (Hamilton et al., 2017)

Attention-based neighborhood aggregation:

Graph attention networks (Hoshen, 2017; Velickovic et al., 2018; Liu et al., 2018)

Embedding entire graphs:

- Graph neural nets with edge embeddings (Battaglia et al., 2016; Gilmer et. al., 2017)
- Embedding entire graphs (Duvenaud et al., 2015; Dai et al., 2016; Li et al., 2018) and graph pooling (Ying et al., 2018, Zhang et al., 2018)
- Graph generation and relational inference (You et al., 2018; Kipf et al., 2018)
- How powerful are graph neural networks(Xu et al., 2017)

Embedding nodes:

- Varying neighborhood: Jumping knowledge networks (Xu et al., 2018), GeniePath (Liu et al., 2018)
- Position-aware GNN (You et al. 2019)

Spectral approaches to graph neural networks:

- Spectral graph CNN & ChebNet (Bruna et al., 2015; Defferrard et al., 2016)
- Geometric deep learning (Bronstein et al., 2017; Monti et al., 2017)

Other GNN techniques:

- Pre-training Graph Neural Networks (Hu et al., 2019)
- GNNExplainer: Generating Explanations for Graph Neural Networks (Ying et al., 2019)

PyTorch Geometric