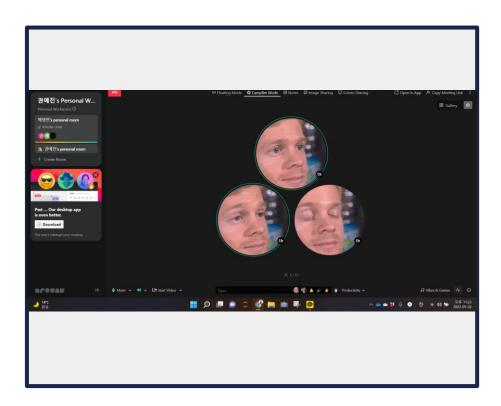
CUAI 스터디 GNN 팀

2022.05.10

발표자 : 권예진

스터디원 소개 및 만남 인증



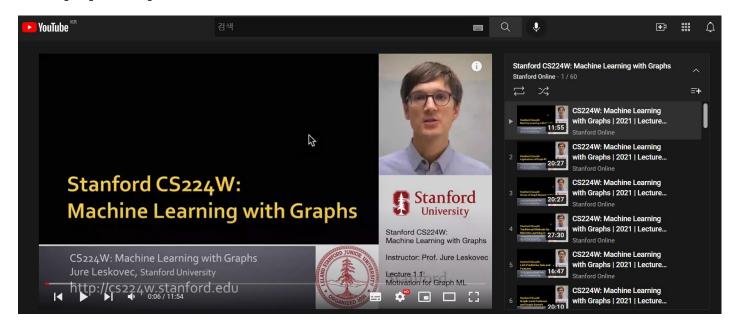
5.2 10:00 PM 세 번째 스터디

<참석자>

스터디원 : 권예진 스터디원 : 이하윤

스터디원: 배병현

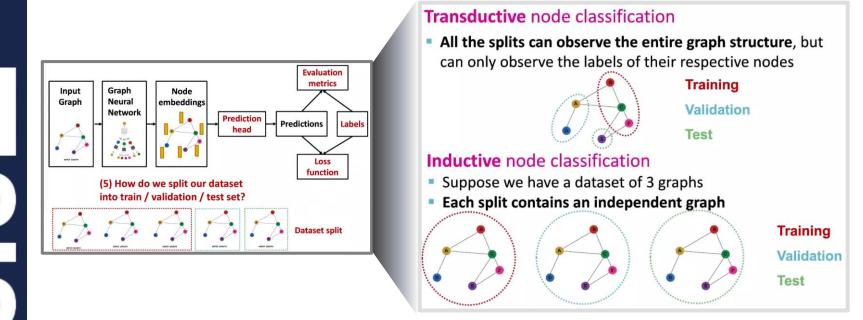
스터디 소개



Stanford CS224W: Machine Learning with Graphs research on the structure and analysis of Graph Neural Network

권예진 스터디원 발표

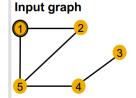
Lecture 8.3 - Setting up GNN Prediction Tasks



이하윤 스터디원 발표

Lecture 9.1 – How Expressive are Graph Neural Networks Lecture 9.2 – Designing the Most Powerful GNNs

To generate a node embedding, GNNs use a computational graph corresponding to a subtree rooted around each node.



Computational graph = Rooted subtree 2 5

Using injective neighbor aggregation

→ distinguish different subtrees

GNN can fully distinguish different subtree structures if every step of its neighbor aggregation is injective.

Next: We analyze aggregation functions of two popular GNN models

- GCN (mean-pool) [Kipf & Welling, ICLR 2017]
 - Uses element-wise mean pooling over neighboring node features

$$Mean(\{x_u\}_{u\in N(v)})$$
 aggregation func.

- GraphSAGE (max-pool) [Hamilton et al. NeurIPS 2017]
 - Uses element-wise max pooling over neighboring node features

$$\operatorname{Max}(\{x_u\}_{u\in N(v)})$$

배병현 스터디원 발표

Lecture 15.1 - Deep Generative Models for Graphs Lecture 15.2 - Graph RNN: Generating Realistic Graphs

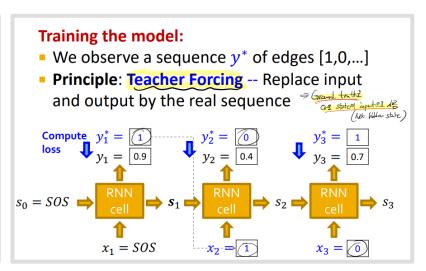
■ Summary: A graph + a node ordering =
A sequence of sequences

■ Node ordering is randomly selected (we will come back to this)

Node-level sequence

Graph G

Adjacency matrix





감사합니다☺