



Representing Long-Range Context for Graph Neural Networks with Global Attention

Zhanghao Wu*, **Paras Jain***, Matthew A. Wright,
Azalia Mirhoseini, Joseph E. Gonzalez, Ion Stoica

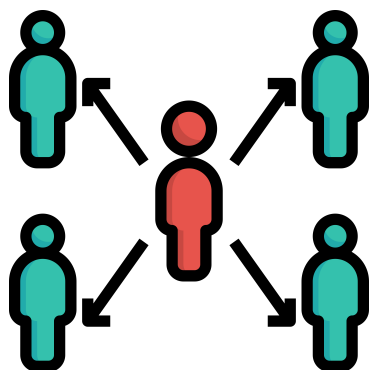
UC Berkeley RISELab, Google Brain



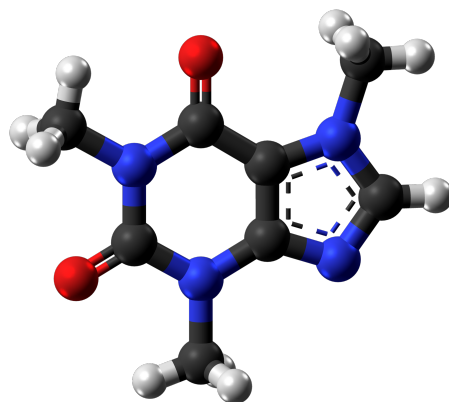
NeurIPS 2021

github.com/ucbrise/graphtrans

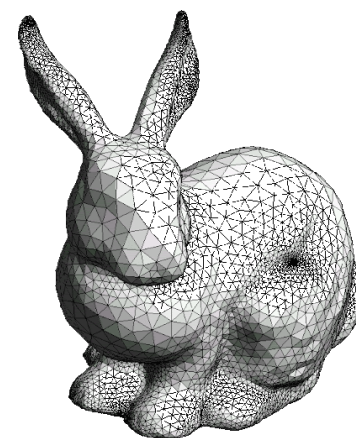
Graphs are an important representation of natural structures



Social networks

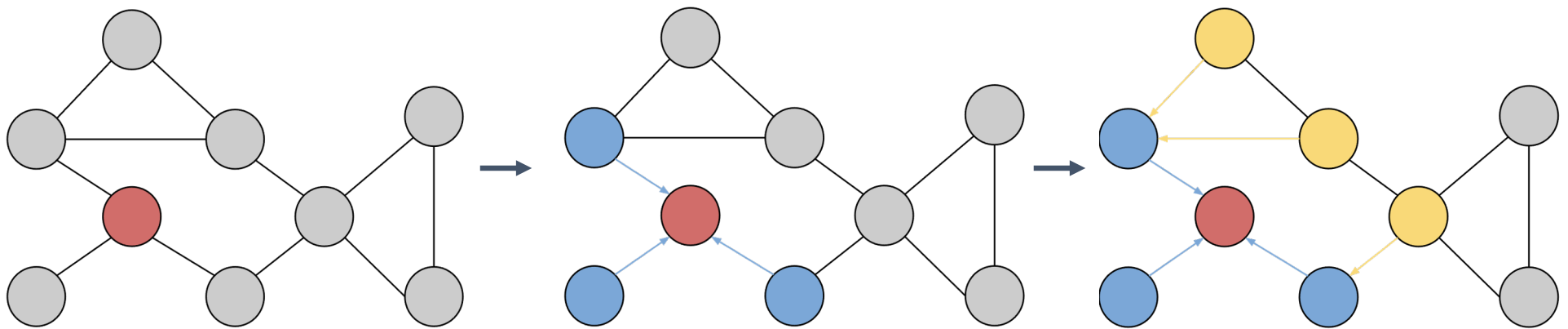


Molecules



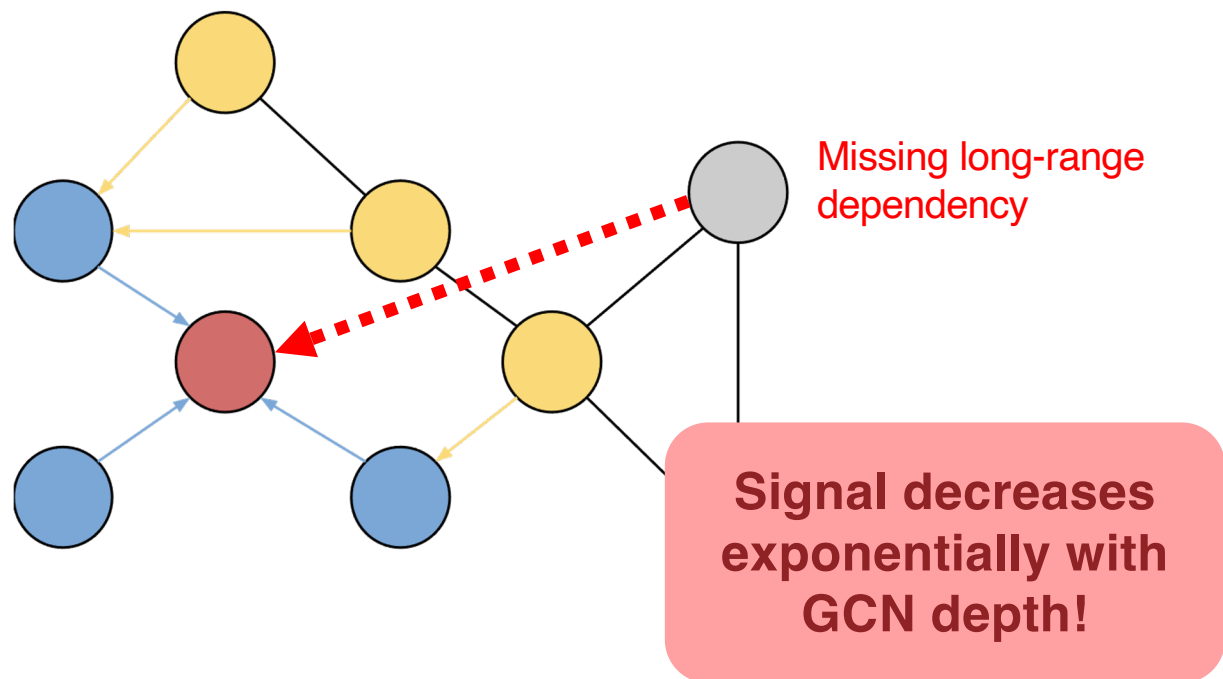
Meshes

Graph neural networks aggregate local neighborhood structure



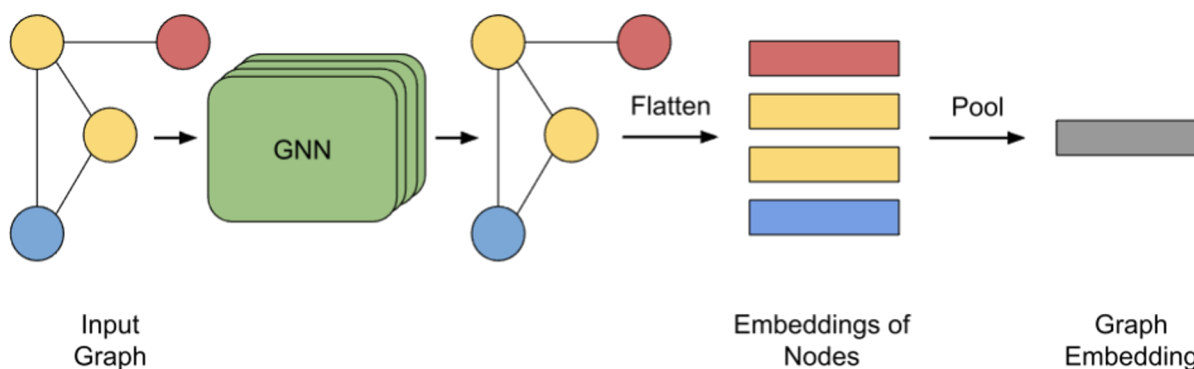
By iteratively pooling immediate neighborhood, GNNs slowly learn to represent local structure

Challenge: long-range dependencies are not represented in GNNs

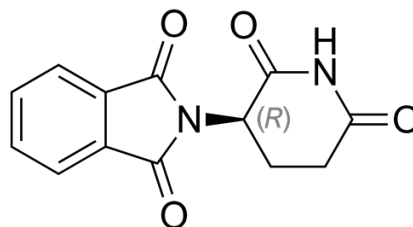


Long-range dependencies important to graph classification

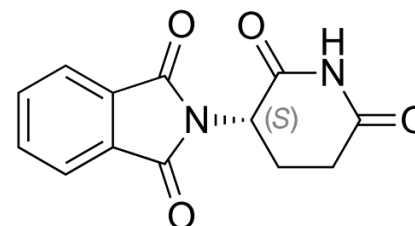
Graph classification considers **pooling embeddings** into single prediction vector.



Small interactions in molecules may result in large changes in function!



(S)-thalidomide = **deadly**

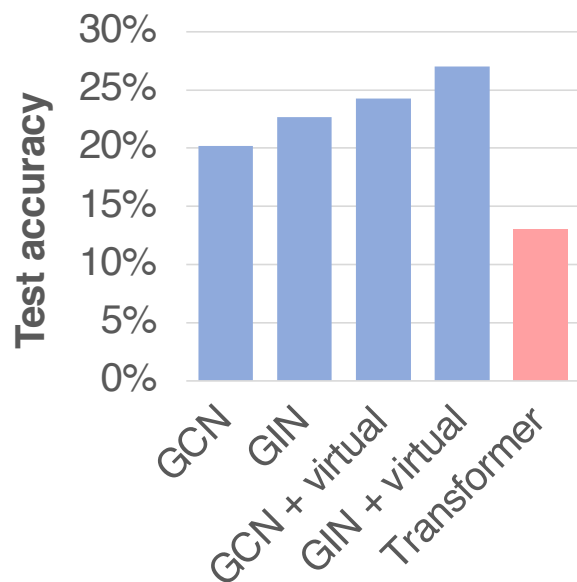


(R)-thalidomide = **safe**

Learning global interactions with GraphTrans



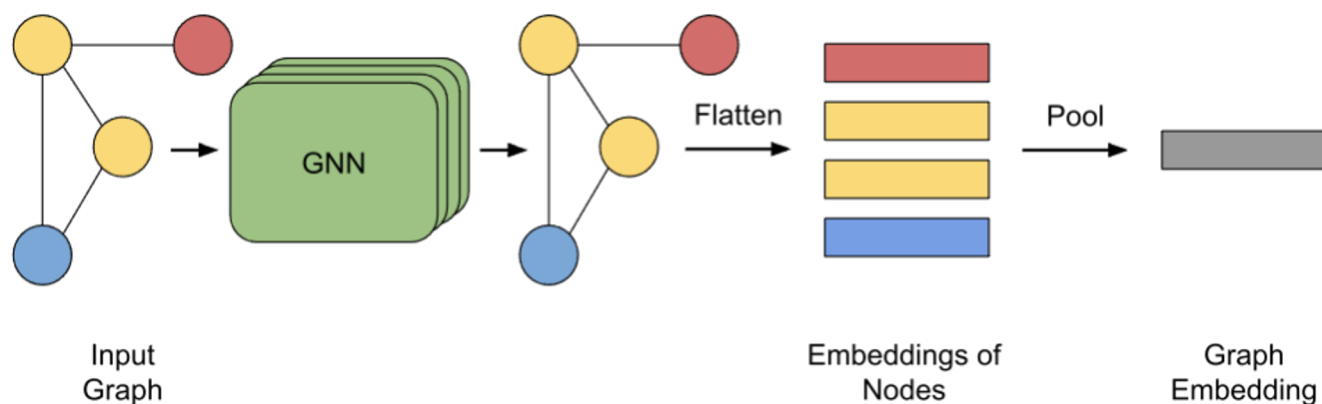
Transformer alone cannot model graph structure



**Transformer results
in -14% test acc. drop!**

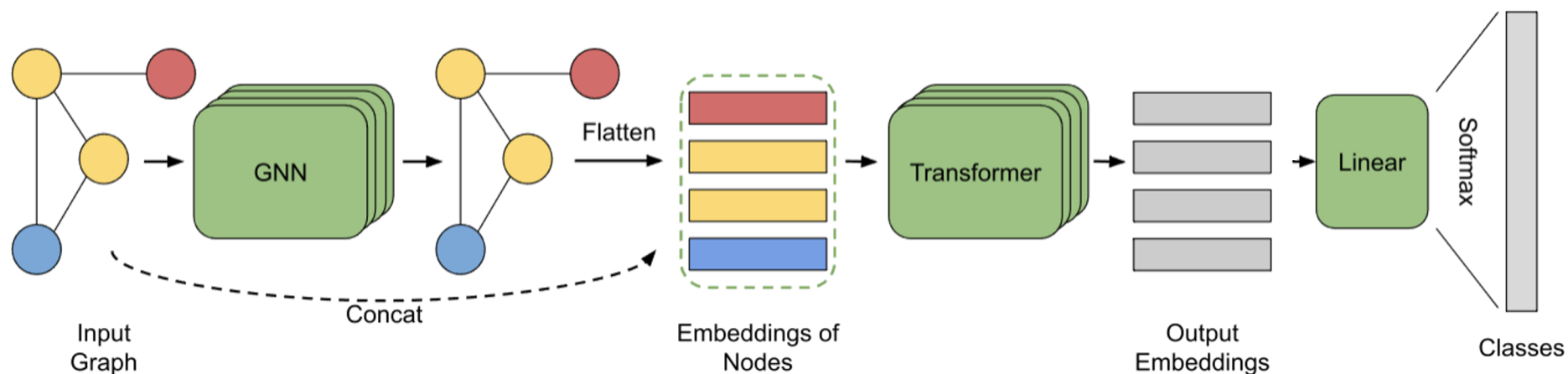
◆ MoleculeNet
OGB-molpcba

Learning global interactions with GraphTrans



We leverage a SOTA
GNN backbone to learn
local structures

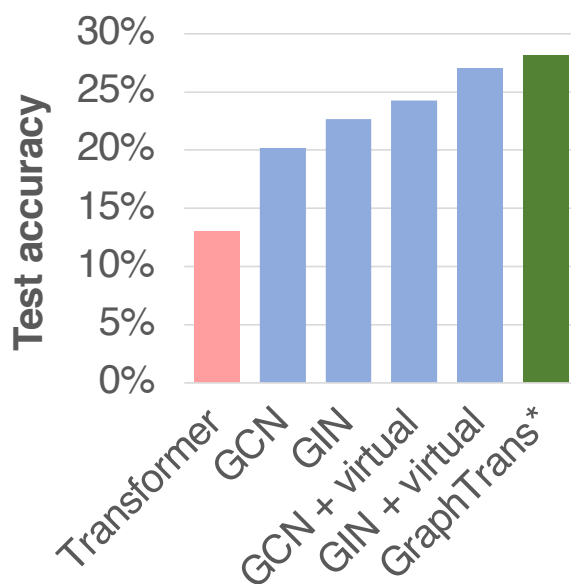
Learning global interactions with GraphTrans



We leverage a SOTA GNN backbone to learn **local structures**

We add a modified Transformer to pool local embeddings to extract **global structures**

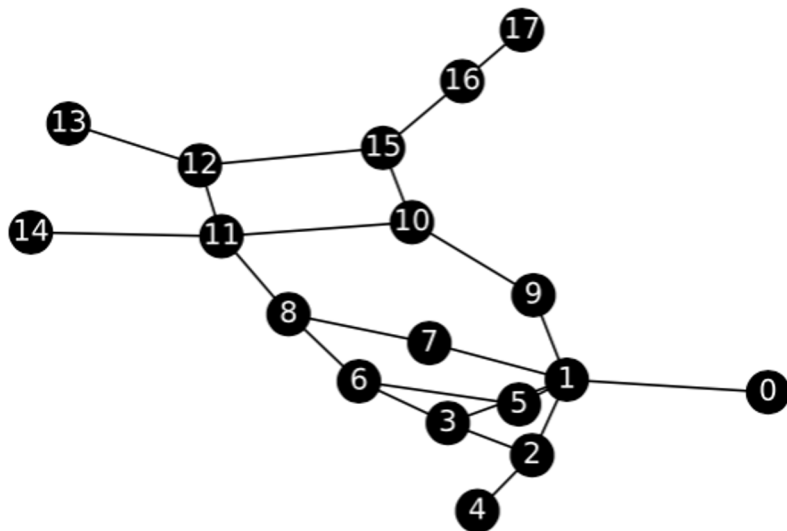
GraphTrans recovers accuracy w/ local + global structure



**Local (GNN) + Global (Transformer) results in
+15% test accuracy over global only,
+1.12% over SOTA**

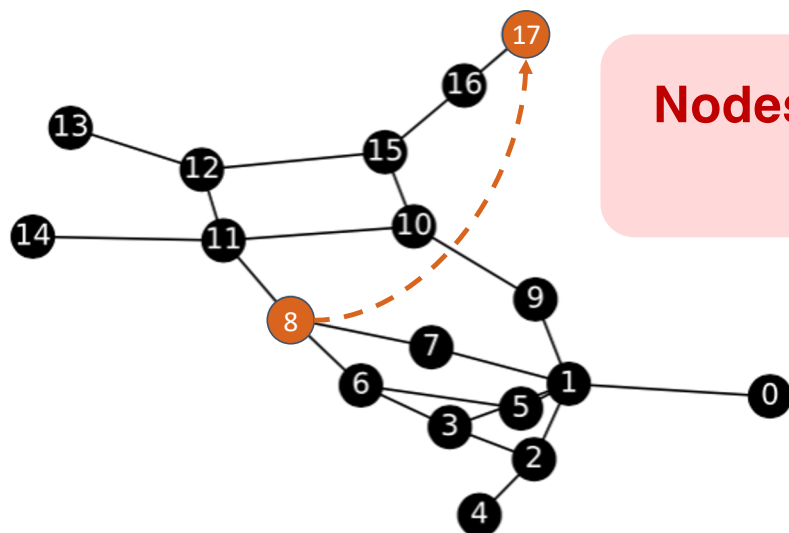
◆ MoleculeNet
OGB-molpcba

Understanding GraphTrans by visualizing attention



Graph from OGB-code2 dataset

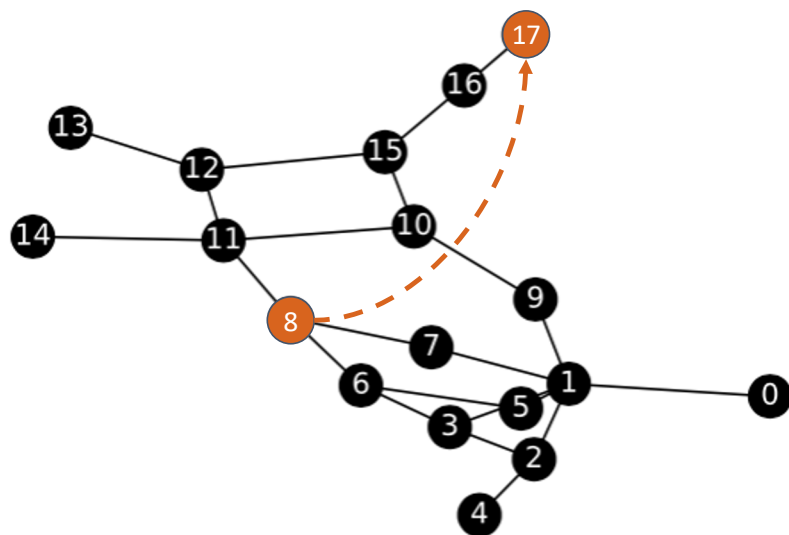
Understanding GraphTrans by visualizing attention



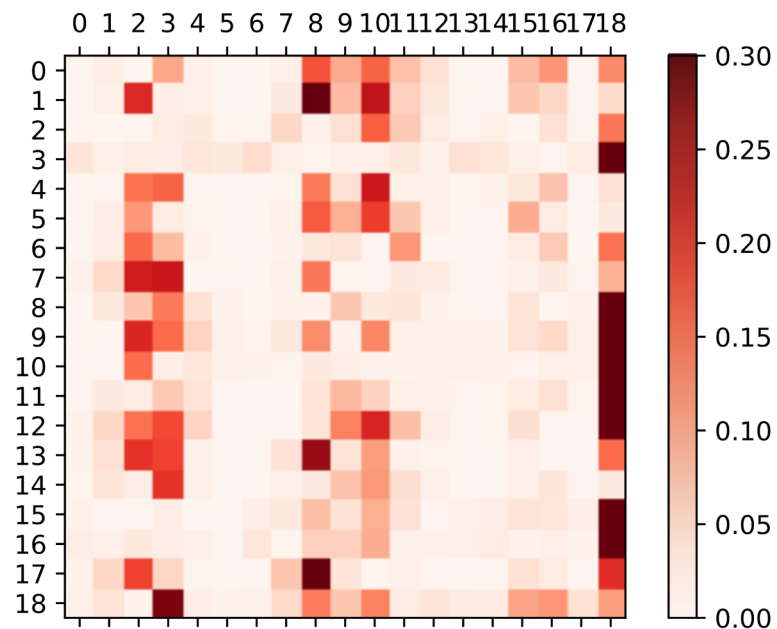
**Nodes 8 and 17 are related,
yet far in graph**

Graph from OGB-code2 dataset

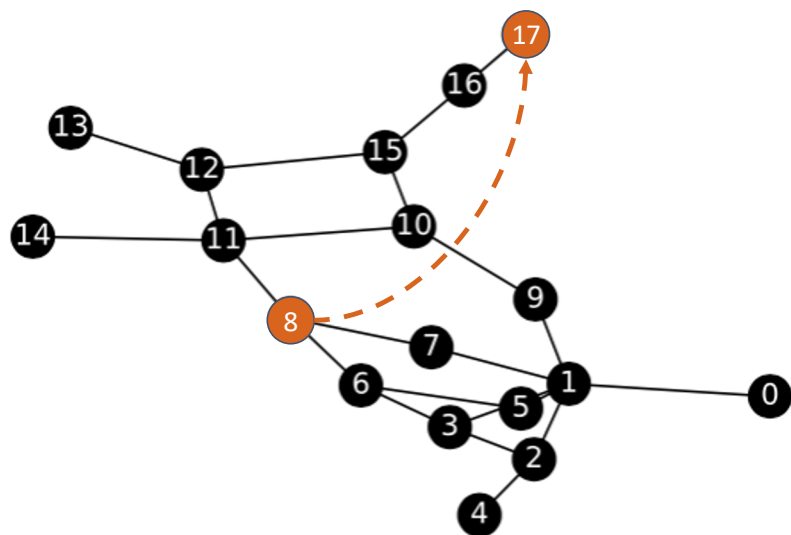
Understanding GraphTrans by visualizing attention



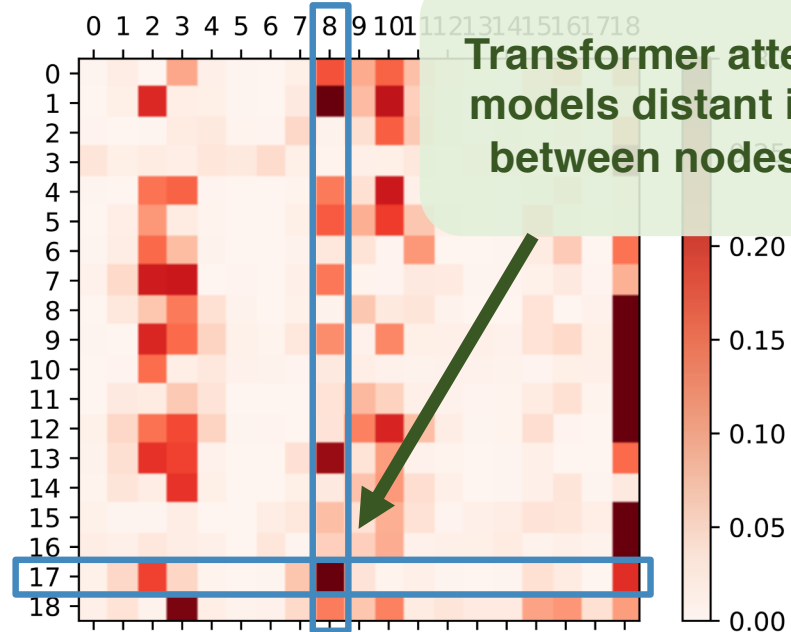
Graph from OGB-code2 dataset



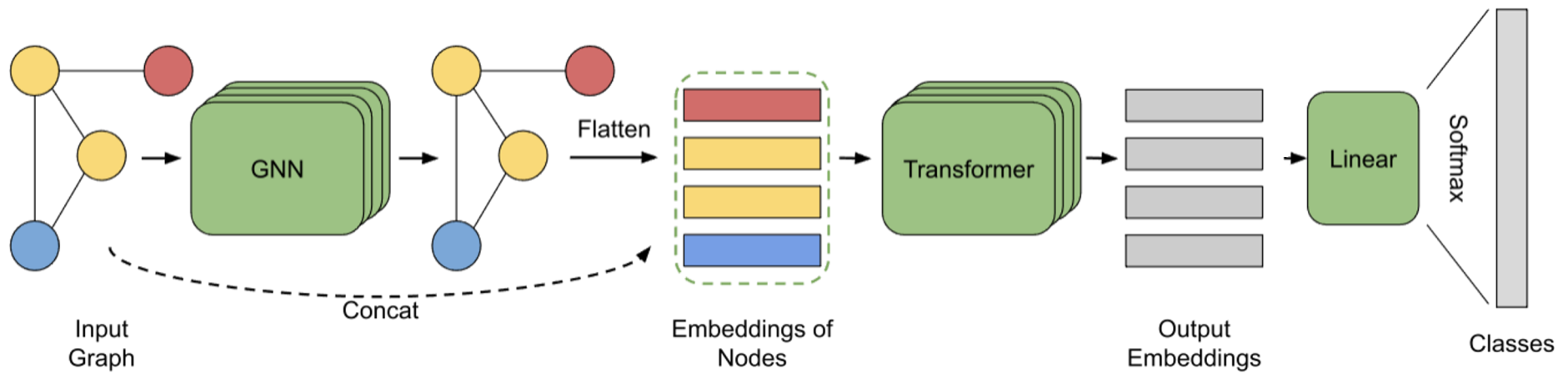
Understanding GraphTrans by visualizing attention



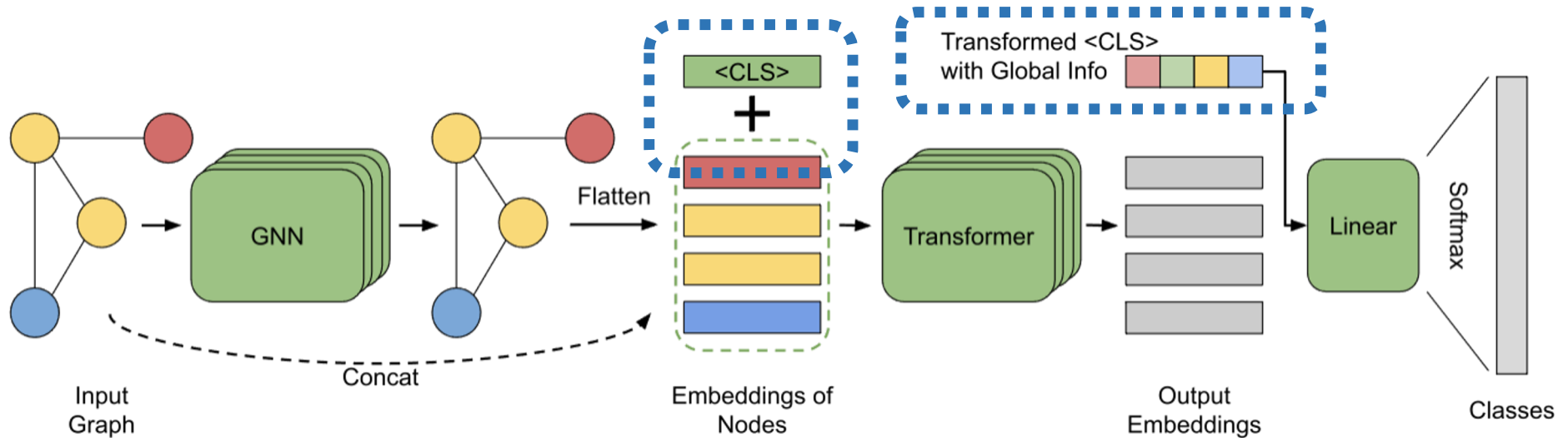
Graph from OGB-code2 dataset



Learning global information with $\langle \text{CLS} \rangle$ token

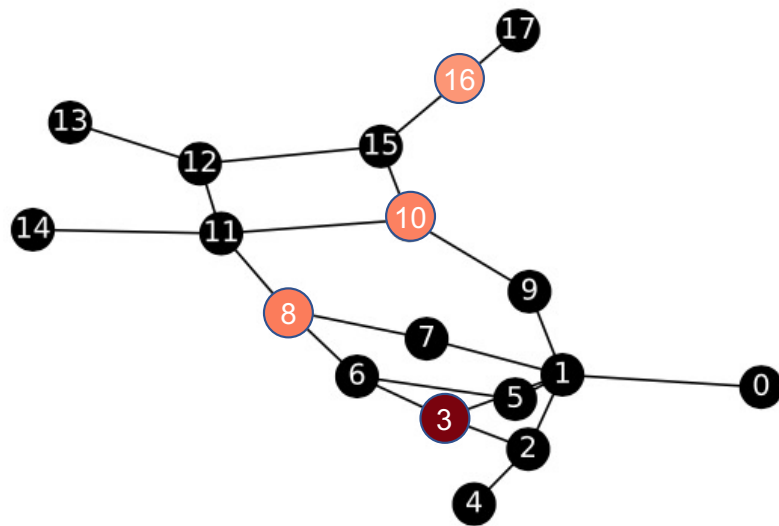


Learning global information with **<CLS>** token

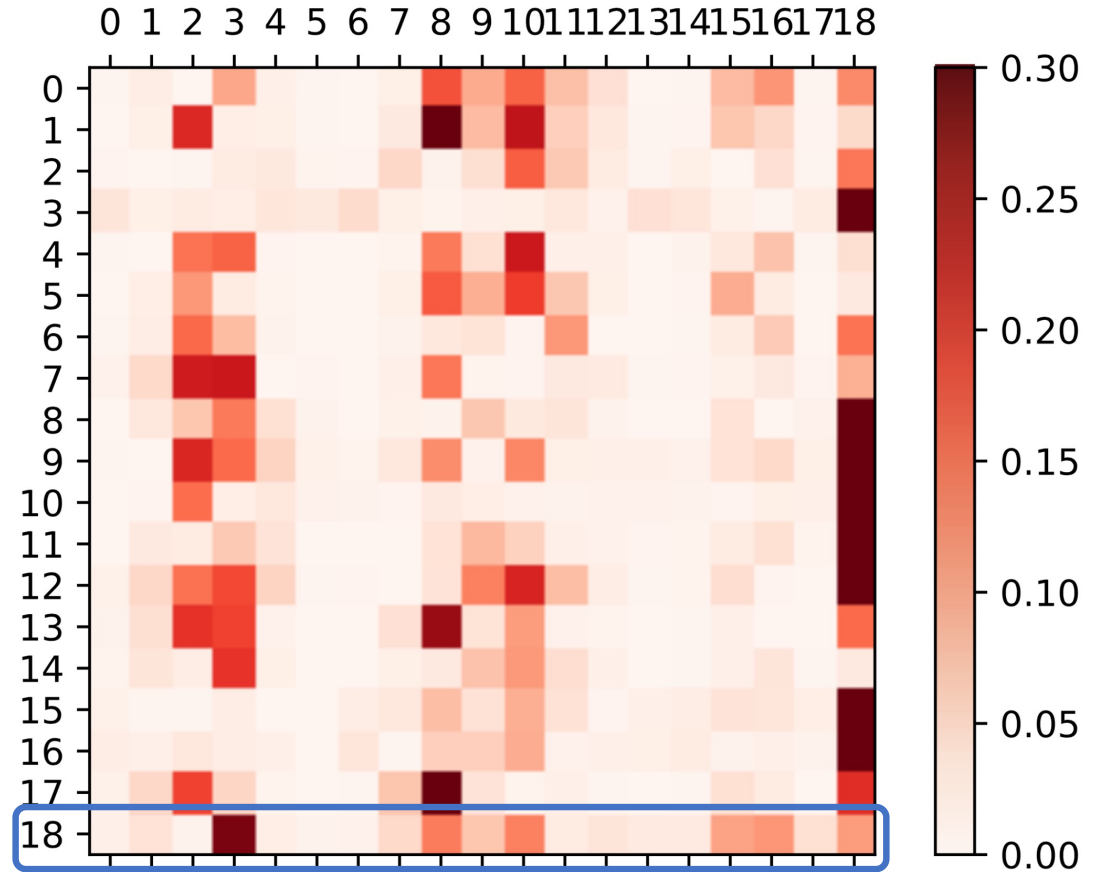


Addition of a single global CLS token aggregates global information into a single vector

Learning global information with **<CLS>** token



<CLS> token learns how
to aggregate global
information together



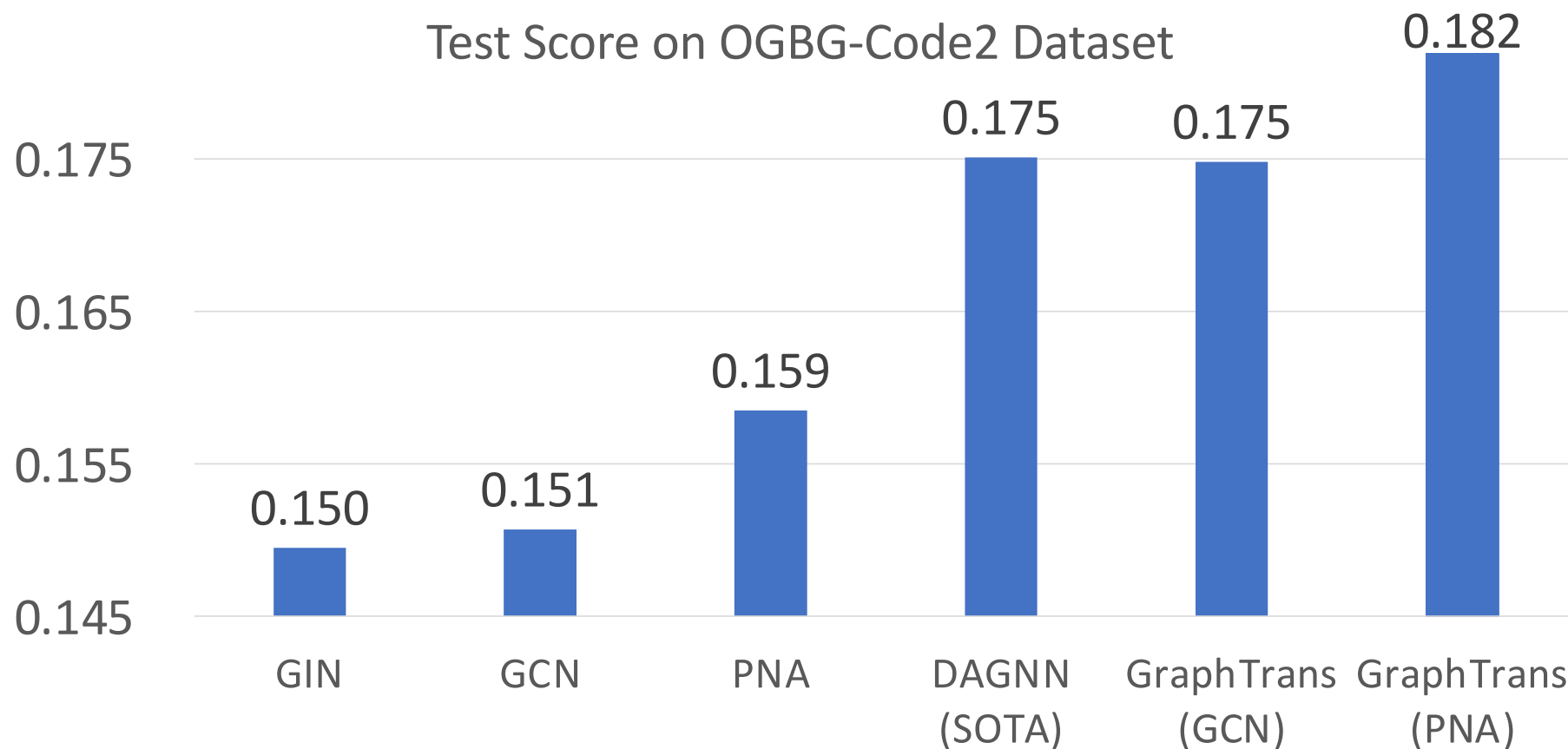
Evaluation: Biological Benchmark

	GNN Type	GNN Layers	NCI1	NCI109
SAGPool _g	GCN	3	74.2	74.1
Strong Baseline	GCN	8	81.5	-
GraphTrans (small)	GCN	3	80.2	79.0
GraphTrans (large)	GIN	4	83.0 (+8.8)	82.5 (+8.4)

Evaluation: Chemical Benchmark

	Valid	Test
GCN-Virtual	0.250	0.242
GIN-Virtual	0.280	0.270
GraphTrans (GIN)	0.288	0.272
GraphTrans (GIN-Virtual)	0.286 (+0.006)	0.282 (+0.012)

Evaluation: Computer Programming Benchmark

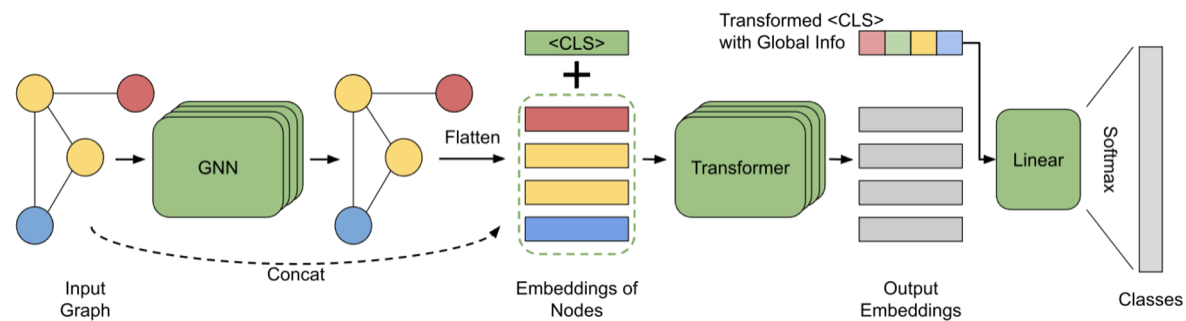


Scalability: Train on Large Graphs

#Nodes	Model	Edge Density			
		20%	40%	60%	80%
500	GCN-Virtual	44.3	58.5	79.3	99.0
	GraphTrans	48.4	57.5	76.4	93.7
1000	GCN-Virtual	90.1	171.8	249.5	OOM
	GraphTrans	96.9	168.4	244.3	OOM
2000	GCN-Virtual	131.8	237.7	OOM	OOM
	GraphTrans	127.9	236.6	OOM	OOM

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- Long-range dependencies crucial to model for graph classification.
- GNNs struggle to learn long-range dependencies.
- GraphTrans combines local representations (from GNN) with global (from Transformer) to learn long-range dependencies.
- We achieve SOTA results on molecular, biological and code prediction datasets.