Old can be Gold: Better Gradient Flow can make Vanilla-GCNs Great Again

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Abstract

Despite the enormous success of Graph Convolutional Networks (GCNs) in modelling graph-structured data, most of the current GCNs are shallow due to the notoriously challenging problems of over-smoothening and information squashing along with conventional difficulty caused by vanishing gradients and over-fitting. Previous works have been primarily focused on the study of over-smoothening and over-squashing phenomenon in training deep GCNs. Surprisingly, in comparison with CNNs/RNNs, very limited attention has been given towards understanding how healthy gradient flow can benefit the trainability of deep GCNs. In this paper, firstly, we provide a new perspective of **gradient flow** to understand the substandard performance of deep GCNs and hypothesize that by facilitating *healthy gradient flow*, we can significantly improve their trainability, as well as achieve state-of-the-art (SOTA) level performance from vanilla-GCNs [1]. Next, we argue that blindly adopting the Glorot initialization for GCNs is not optimal, and derive a **topology-aware isometric initialization** scheme for vanilla-GCNs based on the principles of isometry. Additionally, contrary to ad-hoc addition of skip-connections, we propose to use Dirichlet Energy for **dynamic rewiring of vanilla-GCNs** with skip-connections. Our dynamic rewiring method uses the expressiveness of feature embedding learned by each layer during training to introduce skip-connections *on-demand* basis. We provide extensive empirical evidence across multiple datasets that our methods improves gradient flow in deep vanilla-GCNs and significantly boost their performance to *comfortably compete and outperform* many fancy state-of-the-art methods. Codes are available in the supplement.

1 Introduction

Graphs are omnipresent and Graphs convolutional networks (GCNs)[1] and their variants[2, 3, 4, 5, 6, 7, 8] are a powerful family of neural networks which can learn from graph-structured data. GCNs have been enormously successful in numerous real-world applications such as recommendation systems [9, 10], social and academic networks[11, 9, 5, 4, 12], modeling proteins for drug discovery[13, 14, 10], computer vision[15, 16], etc. Despite their popularity, training deep GCNs are notoriously hard and many classical GCNs[1, 3] achieve their best performance with shallow depth and completely bilk trainability with increasing stacks of layers and non-linearity[17, 18].

Previous works have focused on studying the roadblocks of training deep GCNs from the perspective of *over-smoothening* [17, 18] and *information bolltleneck* [19] and many approaches broadly categorized as *architectural* [20, 21, 17, 22], and *regularization & normalization* [23, 24, 25, 26] tweaks has been proposed for mitigation. In a recent work, [27] theoretically validated that deeper GCN model is at least as expressive as the shallow GCN model, as long as deeper GCNs are trained properly. Furthermore, [28] studied the training difficulty of GCNs from the perspective of graph signal energy

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loss, and proposed modifying the GCN operator, from the energy perspective but it failed to recover the full potential of vanilla-GCNs. Surprisingly, unlike traditional neural networks (CNNs and RNNs), limited effort has been given to understand hurdles in trainability of deep GCNs from the signal propagation perspective, i.e. **gradient flow**. In this paper, we comprehensively show that deep GCNs suffer from *poor gradient flow* (error signals) under back-propagation across multiple datasets and architectures, which significantly hurt their trainability and lead to substandard performance. Further, we are motivated to explore an orthogonal direction: *Can we make vanilla-GCNs [1]* ***go deeper*** *and* ***comparable/better*** *than SOTA, by improvising healthy gradient flow during training?*

In this work, we *firstly* look into the effect of initialization (surprisingly overlooked) in GCNs, and find that blindly adopting the Glorot initialization [29] from CNNs has critical impact on the gradient flow during GCN training, specially for deep GCNs. When the initial weights are not chosen appropriately, the propagation of the input signal into the layers of random weights can result in poor error signals under back-propagation [30]. Inspired from the signal propagation perspective, we leverage the *principle of isometry* [31, 32] to derive a *theoretical initialization scheme* for vanilla-GCNs, by incorporating the graph topological information - which yields a remarkably easy-to-compute form.

Skip-connections have been very resourceful to train deep convolutional networks [33, 34] by improving the gradient flow, and recently some works [20, 21, 35, 17] have identified their benefits for deep GCNs. Although their benefits are known, most of the existing works attempt to add skip-connections in an *ad-hoc and non-optimal* fashion in deep GCNs which creates additional overhead of storing multiple activations during training along with substandard performance. *Secondly*, we study layerwise gradient flow during the training of deep vanilla-GCNs and identify that there exist some layers which receive almost zero error signal (i.e. gradients) during backpropagation and lose trainability and expressiveness. We use *computationally efficient* Dirichlet Energy [36] to dynamically identify such layers during training and propose a principled way of injecting skip-connections on-demand to facilitate healthy gradient flow in deep GCNs. Our main contributions can be summarized as:

- **Deep GCNs suffer from poor gradient flow during training.** We rigorously investigate the *gradient flow* during the training of GCNs and observe that GCNs suffer from poor gradient flow which worsen with increase in depth. We hypothesize that by improving the gradient flow in deep vanilla-GCNs, along with effectively improving their trainability, we can achieve SOTA-level performance from vanilla-GCNs [1].
- **Topology-Aware Isometric Initialization of GCNs.** Unlike blind adoption of Glorot initialization [29] for GCNs, we derive a *theoretical initialization scheme* for vanilla-GCNs [1] based on the principle of isometry, which relates benign initial parameters with graph topology. We experimentally validate that our new initialization strategy significantly improves the gradient flow in deep GCNs and provide huge performance benefits.
- **Dirichlet Energy Guided Dynamic Rewiring of GCNs.** Contrary to *ad-hoc* addition of skip-connections to improve GCNs performance, in this paper, we leverage Dirichlet Energy [36] (which measures the expressiveness of feature embedding) to introduce *dynamic rewiring strategy* of vanilla-GCNs with skip-connections. Our new strategy improves gradient flow, reduces the overhead of storing multiple intermediate activations (dense, residual, jumping skips), as well as outperforms many ad-hoc skip connection mechanisms.

2 Methodology

In this section, we aim to discuss the trainability of deep GCNs from the signal propagation perspective, i.e. gradient flow. We will provide a detailed introduction and theoretical derivation of our topology-aware isometric initialization. Furthermore, we will present a novel way to incorporate skip-connections using Dirichlet Energy to improve the gradient flow and trainability of deep GCNs.

2.1 Preliminaries

We begin by formulating Graph Convolutional Networks (GCN) along the way introducing our notation. Let $G = (\mathcal{V}, \mathcal{E})$ be a simple and undirected graph with vertex set $\mathcal{V}$ and edge set $\mathcal{E}$. Let $\boldsymbol{A} \in \mathbb{R}^{N \times N}$ be the associated adjacency matrix, such that $\boldsymbol{A}_{ij} = 1$ if $(i, j) \in \mathcal{E}$, and $\boldsymbol{A}_{ij} = 0$ otherwise. $N = |\mathcal{V}|$ is the number of nodes. We also define $d_i = \sum_j \boldsymbol{A}_{ij}$ as the node degree of the

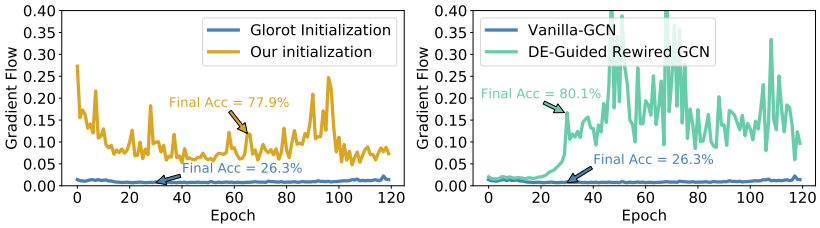

Figure 1: (a) Comparison of Gradient flow in 10-layer vanilla-GCN with default (Glorot initialization) and our Topology-Aware initialization. (b) Comparison of Gradient flow in 10-layer GCN and GCNs with our new Dirichlet Energy guided rewiring.

i-th vertex. Let $\boldsymbol{X} \in \mathbb{R}^{C \times N}$ be the node feature matrix, whose i-th column represents a d-dimension feature vector for the i-th node. GCNs aim to learn an embedding for each node via learnable graph convolutional filters [2, 1]. A graph convolutional layer can be formulated as [1]:

$$f(G, \boldsymbol{X}) = \boldsymbol{W}\boldsymbol{X}(\boldsymbol{A} + \boldsymbol{I}), \tag{1}$$

where adding $\boldsymbol{I}$ to adjacency matrix is known as the self-loop trick [1], $\boldsymbol{W} \in \mathbb{R}^{C' \times C}$ is the weight matrix of this layer, and C' is the dimension of output channel. An L-layer GCN cascades L layer of graph convolutional layers followed by non-linear activations:

$$\boldsymbol{Y} = f^{(L)} \circ \sigma \circ f^{(L-1)} \circ \cdots f^{(2)} \circ \sigma \circ f^{(1)}(G, \boldsymbol{X}), \tag{2}$$

where $\boldsymbol{Y} \in \mathbb{R}^{C'}$ is the final output of the GCN, and $\sigma(\cdot)$ denotes a non-linear activation, which is typically chosen as ReLU. Althrough there exists lots of message-passing based Graph Neural Networks (GNNs), which incorporate with more sophisticate mechanisms, such as attention [3] and gating [37, 38], the focus of our work only lies on the most classical GCN model (Eq. 2 or [1]) that exactly follows the rigorous definition of graph convolutions [39]. As we will show in our experiments, our proposal will bring this theoretical clean GCN back to the SOTA performance.

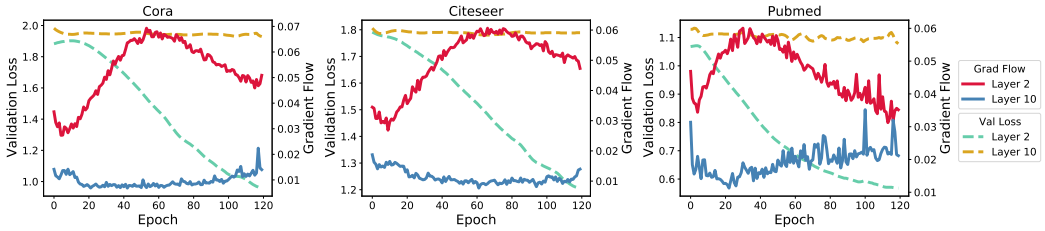

Figure 2: Comparison of validation loss and gradient flow in vanilla-GCNs with 2 and 10 layers.

2.2 Understanding Gradient Flow in Deep GCNs

Gradient flow and its associated problems such as exploding and vanishing gradients have been extensively explored in the dynamics of learning of neural networks (CNNs/RNNs) [29, 30, 32]. GCNs [1] are a special category of neural networks, which uses "graph convolution" operation (linear transformation of all neighbors of a node followed by non-linear activation) to learn the graph representation. Despite being powerful in learning high-quality node representations, GCNs have limited ability to extract information from high-order neighbors and significantly lose their trainability and performance. Recently, many works [17, 18, 19, 21] have been proposed to address this issue from the perspective of over-smoothing and information bottleneck. However, the *signal propagation* perspective to understand the trainability and performance of GCNs has been highly overlooked compared to the attention it has received for CNNs/RNNs.

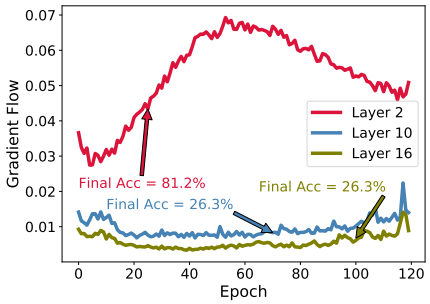

Figure 3: Gradient flow of vanilla-GCNs (layer 4,10,16) trained on Cora.

[1]Here we consider $\boldsymbol{A}$ is symmetrical normalized by default.

Gradient flow is used to study the optimization dynamics of neural networks and it has been widely known that healthy gradient flow facilitates better optimization [29, 30, 32, 40]. Consider an L-layer GCN (see Eq. 2), we denote the weight parameter for the l-th layer $f^{(l)}$ as $\boldsymbol{W}_l$. Suppose we have cost function $\mathcal{L}$, we calculate the gradient across each layer and effective gradient flow (GF) as:

$$\boldsymbol{g}_1 = \frac{\partial \mathcal{L}}{\boldsymbol{W}_1}, \cdots, \boldsymbol{g}_i = \frac{\partial \mathcal{L}}{\boldsymbol{W}_i}, \cdots, \boldsymbol{g}_L = \frac{\partial \mathcal{L}}{\boldsymbol{W}_L} \tag{3}$$

$$\text{Gradient Flow: } \mathrm{GF}_p = \frac{1}{L} \sum_{n=1}^{L} \|\boldsymbol{g}_n\|_p \tag{4}$$

where for every layer l, $\frac{\partial \mathcal{L}}{\partial \boldsymbol{W}_l}$ represents gradients of the learnable parameters $\boldsymbol{W}_l$ and we have used $p = 2$ in our experiments. Figure 3 represents the gradient flow during training of of vanilla-GCNs with layer 4, 6, and 10 on Cora dataset. Figure 2 illustrate the comparison of validation loss and gradient flow in vanilla-GCNs with 2 and 10 layers on Cora, Citeseer, and Pubmed. We consistently observed across each dataset that with increasing depth, the gradient flow across layers decreases significantly along with the performance. In this paper, we hypothesize that by improving the gradient flow in deep vanilla-GCNs, along with effectively improving the their trainability, we can achieve SOTA-level performance from vanilla-GCNs. We propose a new topology-aware initialization strategy based on the principles of isometry and dynamic architecture rewiring for vanilla-GCNs using Dirichlet energy to improvise the gradient flow and subsequently performance (Figure 1(a),(b)).

2.3 Topology-Aware Isometric Initialization

Recent work of [41] has revealed that initializing network paramters near isometric can train very deep CNNs without any normalization. We are inspired to properly initialize each GCN layer to make it an isometric mapping. To begin with, we first give the formal definiton of isometry as below:

Definition 1. *[41] A map $f : \mathcal{X} \to \mathcal{Y}$ between two inner-product spaces is called an isometry if*

$$\langle f(\boldsymbol{x}), f(\boldsymbol{x'}) \rangle = \langle \boldsymbol{x}, \boldsymbol{x'} \rangle \tag{5}$$

for all $\boldsymbol{x}, \boldsymbol{x'} \in \mathcal{X}$.

Existing isometric initialization techniques are not free lunch for GCNs. Qi *et al.* [41] showed that a convolutional operator is isometric if and only if their cross-channel weights are orthogonal to each other. The straightforward implication is that delta kernel can be served as the isometric initialization. However, computing delta kernel in graph signal processing needs to involve complex numbers and eigen-decomposition of the adjacency matrix [42], which is infeasible for a common deep learning framework. To this end, we re-establish theoretical results for GCNs from Definition 1.

Consider one GCN layer (Eq. 1), our goal is to seek a weight matrix such that pair-wise angles between node features are invariant after transforming by the layer. First of all, we denote the output feature of the i-th node as $\boldsymbol{y}_i = \boldsymbol{W}\boldsymbol{X}\boldsymbol{a}_i$ where $\boldsymbol{a}_i$ is the i-th column of $(\boldsymbol{A} + \boldsymbol{I})$. The inner product of two node features $i, j \in [N]$ can be computed as:

$$\langle \boldsymbol{y}_i, \boldsymbol{y}_j \rangle = \left[\left(\boldsymbol{a}_i^\top \otimes \boldsymbol{W} \right) \mathrm{vec}(\boldsymbol{X}) \right]^\top \left[\left(\boldsymbol{a}_j^\top \otimes \boldsymbol{W} \right) \mathrm{vec}(\boldsymbol{X}) \right] \tag{6}$$

$$= \mathrm{Tr}\left(\left[\left(\boldsymbol{a}_i \boldsymbol{a}_j^\top \right) \otimes \left(\boldsymbol{W}^\top \boldsymbol{W} \right) \right] \mathrm{vec}(\boldsymbol{X}) \, \mathrm{vec}(\boldsymbol{X})^\top \right), \tag{7}$$

where $\mathrm{vec}(\cdot)$ denotes vectorization of a matrix, Eq. 6 follows from the equality $\mathrm{vec}(\boldsymbol{ABC}) = (\boldsymbol{C}^\top \otimes \boldsymbol{A}) \, \mathrm{vec} \, \boldsymbol{B}$ [43]. On the other hand, $\boldsymbol{x}_i = \boldsymbol{I}_C \boldsymbol{X} \boldsymbol{e}_i$ where $\boldsymbol{e}_i$ is the i-th canonical basis, we can obtain a similar form of $\langle \boldsymbol{x}_i, \boldsymbol{x}_j \rangle$:

$$\langle \boldsymbol{x}_i, \boldsymbol{x}_j \rangle = \mathrm{Tr}\left(\left[\left(\boldsymbol{e}_i \boldsymbol{e}_j^\top \right) \otimes \boldsymbol{I}_C \right] \mathrm{vec}(\boldsymbol{X}) \, \mathrm{vec}(\boldsymbol{X})^\top \right) \tag{8}$$

To achieve isometry, we hope to solve a $\boldsymbol{W}$ such that Eq. 8 equals to Eq. 7. However, a simple closed-form solution does not exist in general. Instead, we seek a $\boldsymbol{W}$ to minimize the difference between Eq. 8 and Eq. 7.

$$\boldsymbol{W}^* = \underset{\boldsymbol{W} \in \mathbb{R}^{C' \times C}}{\mathrm{argmin}} \sum_{i,j \in [N]} \left| \langle \boldsymbol{y}_i, \boldsymbol{y}_j \rangle - \langle \boldsymbol{x}_i, \boldsymbol{x}_j \rangle \right|^2 \tag{9}$$

$$= \underset{\boldsymbol{W} \in \mathbb{R}^{C' \times C}}{\mathrm{argmin}} \sum_{i,j \in [N]} \mathrm{Tr}^2 \left(\left[\left(\boldsymbol{a}_i \boldsymbol{a}_j^\top \right) \otimes \left(\boldsymbol{W}^\top \boldsymbol{W} \right) - \left(\boldsymbol{e}_i \boldsymbol{e}_j^\top \right) \otimes \boldsymbol{I}_C \right] \mathrm{vec}(\boldsymbol{X}) \, \mathrm{vec}(\boldsymbol{X})^\top \right). \tag{10}$$

However, minimizing objective Eq. 10 will involve data matrix into the solution, which limits computational efficiency and generalization. Hence, we optimize a looser upper bound of Eq. 10 by minimizing the difference between $(\boldsymbol{a}_i\boldsymbol{a}_j^\top) \otimes (\boldsymbol{W}^\top\boldsymbol{W}$ and $(\boldsymbol{e}_i\boldsymbol{e}_j^\top) \otimes \boldsymbol{I}_C$. The intuition is that once this difference reaches zero, Eq. 10 will be minimized to zero as well.

$$\boldsymbol{W}^* = \underset{\boldsymbol{W}\in\mathbb{R}^{C'\times C}}{\operatorname{argmin}} \sum_{i,j\in[N]} \left\| \left(\boldsymbol{a}_i\boldsymbol{a}_j^\top\right) \otimes \left(\boldsymbol{W}^\top\boldsymbol{W}\right) - \left(\boldsymbol{e}_i\boldsymbol{e}_j^\top\right) \otimes \boldsymbol{I}_C \right\|_F^2 . \tag{11}$$

At the first glance, Eq. 17 is non-convex and Kronecker product will produce a computational prohibitive matrix, which makes this problem intractable. However, it has not escaped our notice that the structure of $\boldsymbol{a}_i\boldsymbol{a}_j$ is highly sparse. Hence, the entire norm minimization can be decomposed into block-wise minimization problems. We defer the detailed derivation into the Appendix 1. As the consequence, the optimal solution to Eq. 17 should satisfy:

$$\|\boldsymbol{w}_k\|_2^2 = \frac{N^2}{\sum_{i,j\in[N]}(d_i+1)(d_j+1)} \qquad \forall k \in [C], \tag{12}$$

$$\boldsymbol{w}_k^\top \boldsymbol{w}_l = 0 \qquad \forall k, l \in [C], k \neq l, \tag{13}$$

where $\boldsymbol{w}_k$ denotes the k-th column of optimal weight matrix $\boldsymbol{W}^*$. Eq. 13 suggests that isometry requires weights for different channels to be orthogonal, which conforms with the general results of [41]. But different from [41], our results indicate that the magnitude of channel-wise weights should be constrained by a degree-related constant (Eq. 12), which binds the initial weights with topological information. To randomly initialize $\boldsymbol{W}$ in practice, one can draw each column of $\boldsymbol{W}$ from independent distributions (e.g., white Gaussian) with variance $\Sigma^2 = N^2 / \left[C' \sum_{i,j}(d_i+1)(d_j+1)\right]$, where C' is the number of output channels (a.k.a. fan out). To initialize weights with bounded values, we suggest employ the following uniform distribution:

$$\boxed{\boldsymbol{W} \sim \mathcal{U}\left[-\sqrt{\frac{3N^2}{C'\left(\sum_{i,j\in[N]}(d_i+1)(d_j+1)\right)}}, \sqrt{\frac{3N^2}{C'\left(\sum_{i,j\in[N]}(d_i+1)(d_j+1)\right)}}\right]} \tag{14}$$

2.4 Dirichlet Energy Guided Rewiring

Graph Convolutions can be considered as a type of Laplacian smoothing, and repeatedly applying Laplacian smoothing many times in the case of multi-layer GCNs, leads features of the nodes in the graph to converge to similar values, i.e. lose expressiveness and trainability [17]. Recently, [36] argued to use Dirichlet energy to measure the expressiveness of node embeddings learned by GCNs. Dirichlet Energy is *computationally efficient* since it is calculated across the connected pair of nodes and can be an indicator of expressiveness of learned node embeddings during the training.

We track the Dirichlet energy of feature matrix w.r.t. the (augmented) normalized Laplacian at different layers of the vanilla-GCN during training. Surprisingly, we observed that during training of a deep vanilla-GCN, the Dirichlet energy of some layers decreases exponentially and becomes close to zero within a few epochs of training, i.e lose expressiveness. Detailed gradient flow analysis of these layers (Figure 4 and 5 (a)) further indicate that they receive almost zero error signal during backpropagation (i.e. gradients) and unable to train. Motivated by this finding, we use computationally efficient Dirichlet energy to identify layers that block healthy gradient flow during training and propose to dynamically rewire the vanilla-GCNs using skip-connections (widely known to improve gradient flow), which we call *on-demand dynamic rewiring*. Our dynamic rewiring technique *improves gradient flow, retains Dirichlet Energy, and significantly helps in training* deep vanilla-GCNs with high performance (Figure 1(b) and 4). In our work, Dirichlet energy is defined as:

Definition 2. *[36] Dirichlet energy $E(\boldsymbol{x})$ of a scalar function $\boldsymbol{x} \in \mathbb{R}^N$ on the graph G is defined as:*

$$E(\boldsymbol{x}) = \boldsymbol{x}^T \boldsymbol{L}\boldsymbol{x} = \frac{1}{2}\sum_{(i,j)\in\mathcal{E}} \left(\frac{x_i}{(1+d_i)^{1/2}} - \frac{x_j}{(1+d_j)^{1/2}}\right)^2 \tag{15}$$

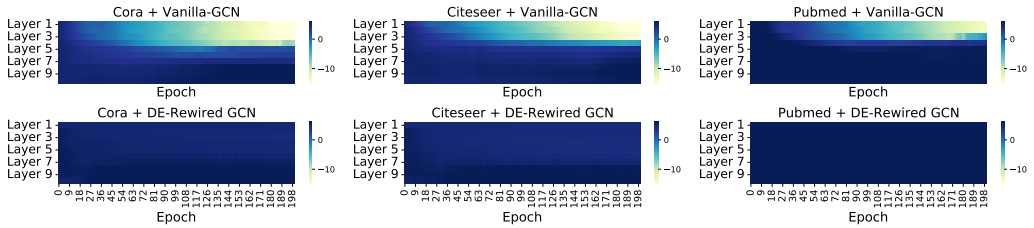

Figure 4: Visualization of Dirichlet Energy of layers during the training of 10-layers vanilla GCN (row 1) and DE-Rewired GCN for Cora, Pubmed, and Citeseer. Plotted using the log scale for better visualization. DE-rewiring prevents Dirichlet Energy to drop to zero during training.

For the vector field $\boldsymbol{X} = [\boldsymbol{x}_1 \quad \cdots \quad \boldsymbol{x}_N] \in \mathbb{R}^{C \times N}$ *, where* $\boldsymbol{x}_i \in \mathbb{R}^C$*, Dirichlet energy is defined as:*

$$E(\boldsymbol{X}) = \mathrm{Tr}(\boldsymbol{X}\boldsymbol{L}\boldsymbol{X}^T) = \frac{1}{2} \sum_{(i,j) \in \mathcal{E}} \left\| \frac{\boldsymbol{x}_i}{(1+d_i)^{1/2}} - \frac{\boldsymbol{x}_j}{(1+d_j)^{1/2}} \right\|_2^2 \tag{16}$$

More specifically, we observe the Dirichlet energy of each vanilla-GCN layer during training, and if the energy drop by $\mathbf{p}\%$ (hyperparameter) of its initial value at the start of training, we assist the layer with a skip-connection. We use a modified version of *initial residual connection* [44, 21], which we define as the output feature matrix of first layer of vanilla-GCN (gives better performance than [44, 21]), to supplement the layers that quickly loses energy. Mathematically, our dynamic rewiring can be written as $X^l = \mathbb{1}[E(X^{(l_t)}) < pE(X^{(l_0)})]\alpha X^{(1)} + \tilde{X}^l$ where $\mathbb{1}[\cdot]$ is an indicator function, t denotes the training epoch, and $\tilde{X}^l = W_{l-1}X^{(l-1)}(A+I)$. Once the layers which are more prone to lose expressiveness are identified and rewired using the skip-connections, the vanilla-GCN training can proceed as normal. In comparison with previous work [20, 21, 35, 17] which blindly inherits skip-connection techniques from CNNs/RNNs, our work provides a principled approach to perform architectural rewiring of vanilla-GCNs. Our extensive experiments on multiple datasets reveal that our approach of guided rewiring not only brings benefits on the efficiency front (memory overhead to store intermediate activations) but also comfortably outperforms all SOTA adhoc skip-connection.

3 Experiment

In this section, we first provide experimental evidence to augment our signal propagation hypothesis and show that our newly proposed methods facilitate healthy gradient flow during the training of deep-vanilla GCNs. Next, we extensively evaluate our methods against state-of-the-art graph neural network models and techniques to improve vanilla-GCNs on on a wide variety of open graph datasets.

Settings	Cora	Citeseer	Pubmed	OGBN-ArXiv
{Learning rate, Weight Decay, Hidden dimesnion, DE energy ratio ($\mathbf{p}$)}	{0.005, $5e-4$, 64, 0.4}	{0.005, $5e-4$, 64, 0.55}	{0.01, $5e-4$, 64, 0.4}	{0.005, 0, 256, 0.5}

Table 1: Hyperparameter configuration for our proposed method on representative datasets.

3.1 Dataset and Experimental Setup

We use three standard citation network datasets Cora, Citeseer, and Pubmed [45] in GNN domain for evaluating our proposed methods against state-of-the-art GNN models and techniques. We have used the basic vanilla-GCN implementation in PyTorch provided by the authors of [1] to incorporate our proposed techniques and show their effectiveness in making traditional GCN comparable/better with SOTA. For our evaluation on Cora, Citeseer, Pubmed, and OGBN-ArXiv, we have closely followed the data split settings and metrics reported by the recent benchmark [46]. For comparison with SOTA models, we have used JKNet [47], InceptionGCN [48], SGC [49], GAT [3], GCNII [21], and DAGNN [50]. We use Adam optimizer for our experiments and performed a grid search to tune hyperparameters for our proposed methods and reported our settings in Table 1. For all our experiments, we have trained our modified GCNs for 1500 epochs and 100 independent repetitions following [46] and reported average performances with the standard deviations of the node classification accuracies. All experiments on large graph datasets, e.g., OGBN-ArXiv, are conducted on single 48G Quadro RTX 8000 GPU, while small graph experiments are completed using a single 16G RTX 5000 GPU.

3.2 Gradient Flow and our proposed methods

Despite significant efforts to improve deep neural networks (CNNs/RNNs) training from the signal propagation perspective, in-depth analysis of deep GCNs training with a focus on the gradient flow during back-propagation is highly overlooked. We use Equation 4 to study gradient flow during the training of GCNs and Figure 3 indicates that with increase in depth, gradient flow in the network drop significantly hurting the trainability of GCNs. Figure 2 indicates the relation between gradient flow and drop in validation loss (performance) for 2-layer and 10-layer GCNs for Cora, Citeseer, and Pubmed. It can be observed that across all datasets, 10-layer GCN has poor gradient flow compared to 2-layer GCN and we observed a negligible drop in validation loss during training.

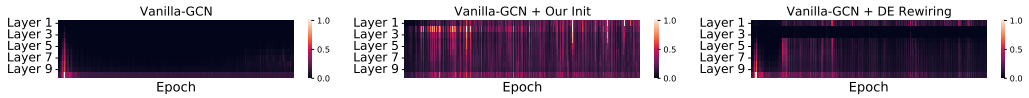

Figure 5: Visualization of gradient flow of layers during the training of 10-layers vanilla GCN, vanilla-GCN with our new topology-aware initialization and dynamic DE-based rewiring for 500 epochs on Cora. Our methods promotes uniform gradient flow across all layers of vanilla-GCNs.

To better understand the behavior of each hidden layer in deep GCN training, we estimated the gradient flow per layer independently of a 10-layer GCN. To our surprise, we found that *many layers receive zero error signals (gradients)* under backpropagation, potentially leading to a catastrophic failure to update during training. Figure 5(a) illustrate layer-wise analysis of gradient flow in 10-layer vanilla-GCN on Cora. We observed that within a few epochs of training, the gradient flow saturates, and then onwards, many layers of GCN receive zero error signal during the backpropagation and they fail to update leading to no drop in validation loss. This conforms very well with our observation in Figure 4 that within a few epochs of training, Dirichlet Energy of many layers drop to zero and they lose expressiveness.

Dataset	**Settings**	**Layer Number**										
		2	3	4	5	6	7	8	9	10	11	12
Cora	GCN	81.1	81.9	80.4	79.9	77.5	77.1	69.5	28.2	27.4	30.1	25.4
	GCN + Our Init	83.0	83.4	80.6	80.0	81.2	80.6	80.4	79.9	80.1	79.2	78.5
CiteSeer	GCN	71.4	67.6	64.6	63.7	63.9	24.1	20.8	22.3	23.2	22.9	21.6
	GCN + Our Init	71.7	78.2	68.3	66.1	66.2	63.3	62.8	63.2	62.9	59.8	61.9
PubMed	GCN	79.0	78.9	76.5	76.7	77.1	76.5	61.2	41.8	40.7	43.2	41.0
	GCN + Our Init	79.3	80.0	78.5	77.6	77.6	76.7	75.8	76.9	76.2	76.3	75.9

Table 2: Performance Comparison (test accuracy %) of of vanilla-GCNs [1] with and without our newly proposed topology-aware isometric initialization. Experiments are conducted on Cora, Citeseer, and PubMed using vanilla-GCNs with layer $l \in \{2, 3, ..., 12\}$.

Initialization of neural networks is tightly coupled with the propagation of error signals [30] and effect their trainability. Our newly proposed topology-aware isometric initialization and Dirichlet Energy guided rewiring significantly improve gradient flow (Figure 1(a) and (b)) as well as mitigate the issue of exponentially dropping expressiveness of layers in deep GCNs (Figure 4). In Figure 5(b) and (c), the detailed layer-wise analysis of vanilla-GCN with our newly proposed methods clearly illustrate how our methods are able to *uniformly restore healthy gradients across all layers* leading to improved trainability and significant performance gains.

3.3 Vanilla-GCNs and our proposed methods

In this section, we conduct a systematic study to understand the performance gain by improving gradient flow using our proposed methods: topology-aware isometric initialization and Dirichlet Energy Guided Dynamic Rewiring, by incorporating them into the training process of deep vanilla-GCNs. Table 2 demonstrate the performance comparison of vanilla-GCNs [1] with and without our new initialization method with increase in depth. Our experiments on Cora, Citeseer, and PubMed using vanilla-GCNs with layers $l \in \{2, 3, ..., 12\}$ illustrate how our new initialization has been successful in retaining the trainability of vanilla-GCNs with increasing depth along with improving performance at shallow depths. The highlighted (red) box represents the depths at which vanilla-GCNs lose trainability, i.e, validation loss shows no improvement during training.

Category	Settings	Cora			Citeseer			PubMed		
		2	16	32	2	16	32	2	16	32
Vanilla-GCN	-	81.10	21.49	21.22	71.46	19.59	20.29	79.76	39.14	38.77
Skip Connection	Residual	74.73	20.05	19.57	66.83	20.77	20.90	75.27	38.84	38.74
	Initial	79.00	78.61	78.74	70.15	68.41	68.36	77.92	77.52	78.18
	Jumping	80.98	76.04	75.57	69.33	58.38	55.03	77.83	75.62	75.36
	Dense	77.86	69.61	67.26	66.18	49.33	41.48	72.53	69.91	62.99
Normalization	BatchNorm	69.91	61.20	29.05	46.27	26.25	21.82	67.15	58.00	55.98
	PairNorm	74.43	55.75	17.67	63.26	27.45	20.67	75.67	71.30	61.54
	NodeNorm	79.87	21.46	21.48	68.96	18.81	19.03	78.14	40.92	40.93
	CombNorm	80.00	55.64	21.44	68.59	18.90	18.53	78.11	40.93	40.90
Random Dropping	DropNode	77.10	27.61	27.65	69.38	21.83	22.18	77.39	40.31	40.38
	DropEdge	79.16	28.00	27.87	70.26	22.92	22.92	78.58	40.61	40.50
	LADIES	77.12	28.07	27.54	68.87	22.52	22.60	78.31	40.07	40.11
Identity Mapping	-	**82.98**	67.23	40.57	68.25	56.39	35.28	79.09	79.55	73.74
DE-Guided Rewiring (Ours)		82.79 ±0.16	**80.40** ±0.23	**79.92** ±0.19	**71.29** ±0.21	**68.59** ±0.09	**68.51** ±0.13	**78.90** ±0.20	**78.10** ±0.11	**78.49** ±0.14

Table 3: Performance Comparison of DE-Guided Rewiring with respect to various proposed fancy techniques to improve the GCN training. Note that our results are generated using vanilla-GCN [1]. Experiments are conducted on Cora, Citeseer, and PubMed with 2/16/32 layers GCN.

Trainability issue of deep GCNs has been extensively studied from the perspective of *over-smoothening* [17, 18] and *information bottleneck* [19] and many approaches broadly categorized as *architectural tweaks* [20, 21, 17, 22], and *regularization & normalization* [23, 24, 25, 26] has been proposed for mitigation. Table 3 demonstrate the performance comparison of our Dirichlet Energy guided dynamic rewiring using skip-connections with respect to various state-of-the-art techniques to improve deep GCNs training. It can be clearly observed that our on-demand dynamic rewiring method of introducing new skip-connections significantly outperforms all fancy normalization and regularization techniques as well as comfortably beats ad-hoc skip-connections techniques.

Method	Cora						PubMed					
	2	4	8	16	32	64	2	4	8	16	32	64
Vanilla-GCN [1]	81.1	80.4	69.5	21.5	21.2	21.9	79.0	76.5	61.2	39.1	38.7	35.3
GAT [3]	81.9	80.3	31.3	30.5	27.1	27.9	78.4	77.4	29.1	26.3	28.7	25.0
JKNet [47]	79.1	79.2	75.0	72.9	73.2	71.5	77.8	68.7	67.7	69.8	68.2	63.4
SGC [49]	79.3	79.0	77.2	75.9	68.5	65.3	78.0	73.1	70.9	69.8	66.6	63.2
InceptionGCN [48]	79.2	77.6	76.5	81.7	81.7	80.0	78.5	77.7	77.9	74.9	74.1	74.3
GCNII [21]	82.2	**82.6**	**84.2**	**84.6**	**85.4**	**85.5**	78.2	78.8	79.3	**80.2**	**79.8**	**79.7**
Ours	**83.1**	82.8	83.0	82.9	82.1	80.3	**80.2**	**79.6**	**79.7**	79.6	79.7	**79.9**
(std: ±)	0.25	0.23	0.10	0.23	0.14	0.27	0.16	0.29	0.03	0.15	0.09	0.14

Table 4: Performance Comparison (test accuracy %) of our proposed method (topology-aware initialization and DE-guided rewiring) with other previous SOTA frameworks using Cora and Pubmed.

3.4 Comparison with state-of-the-art methods

To further validate the effectiveness improving gradient flow in vanilla-GCNs, we perform comparisons with previous state-of-the-art frameworks, including SGC[49], GAT[3], JKNet[47], APPNP[44], InceptionGCN[48], GPRGNN[51], and GCNII[21]. We first observe that Dirichlet Energy guided rewiring and Topology-aware isometric initialization brings orthogonal benefits and combining them helps vanilla-GCNs to achieve better performance. Table 4 reports the mean classification accuracy and the standard deviation on the test set of Cora and Pubmed of our methods using vanilla-GCNs with layer $l \in \{2, 4, 8, 16, 32, 64\}$. One key benefit to note is that with the help of our methods, we can train vanilla-GCNs as deep as 64 layers. It can be observed that vanilla-GCNs with our methods outperform all SOTA methods on Cora (layers 2 and 4) and Pubmed (layers 2, 4, 8, and 64), while holding the second position for all other layers. To evaluate on a large graph, we have chosen OGBN-AxRiv dataset, and Table 5 represents the performance comparison of vanilla-GCNs with our method against SOTA baselines. It can be observed that with the help of our methods, vanilla-GCNs can be efficiently trained deeper without any significant loss in performance. Very similar to Cora and Pubmed, our methods beats all SOTA methods while training 2-layer vanilla-GCNs and holds second position for all other layers. Note that we are NOT trying to beat SOTA, but instead our objective is to reveal the correct initialization and training scheme for vanilla-GCNs, and scale deep efficiently.

Layer	GCN[1]	SGC[49]	DAGNN [22]	GCNII[21]	JKNet[47]	APPNP[44]	GPRGNN[51]	**Ours**
2	69.46±0.22	61.98±0.08	67.65±0.52	**71.24±0.17**	63.73±0.38	65.31±0.23	69.31±0.09	**71.81±0.08**
16	67.96±0.38	41.58±0.27	**71.82±0.28**	**72.61±0.29**	66.41±0.56	66.95±0.24	70.30±0.15	71.76±0.03
32	45.46±4.50	34.22±0.04	71.46±0.27	**72.60±0.25**	66.31±0.63	66.94±0.26	70.18±0.16	**71.73±0.11**

Table 5: Performance Comparison (test accuracy %) of our proposed method (topology-aware initialization and DE-guided rewiring) with other previous SOTA frameworks using OGBN-AxRiv.

The last sanity check is whether our proposed methods can make deep vanilla-GCNs effective across multiple different graph datasets. Specially, we evaluate it on seven other open-source graph datasets: (i) one Co-author datasets [52] (CS), (ii) two Amazon datasets [52] (Computers and Photo), (iii) three WebKB datasets [53] (Texas, Wisconsin, Cornell), and (iv) the Actor dataset [53]. Table 6 reports the performance comparison of deep vanilla-GCN with 32-layers with state-of-the-art methods having the same depth. Our methods universally encourage significant trainability benefits to deep vanilla-GCNs across all datasets achieving state-of-the-art performance on Computers dataset while performing in top-2 for all remaining datasets.

Category	CS [52]	Computers [52]	Photo [52]	Texas [53]	Winconsin [53]	Cornell [53]	Actors [53]
GCN [1]	24.01±3.42	58.72±4.97	58.64±8.40	60.12±4.22	52.94±3.99	54.05±7.11	25.46±1.43
GAT [3]	11.04±0.94	9.42±0.54	17.11±1.02	11.54±0.72	14.01±0.88	19.78±1.42	6.42±0.39
SGC [49]	70.52±3.96	37.53±0.20	26.60±4.64	56.41±4.25	51.29±6.44	58.57±3.44	26.17±1.15
GCNII [21]	71.67±2.68	37.56±0.43	62.95±9.41	**69.19±6.56**	**70.31±4.75**	**74.16±6.48**	**34.28±1.12**
JKNet [47]	81.82±3.32	**67.99±5.07**	**78.42±6.95**	61.08±6.23	52.76±5.69	57.30±4.95	28.80±0.97
APPNP [44]	**91.61±0.49**	43.02±10.16	59.62±23.27	60.68±4.50	54.24±5.94	58.43±3.74	28.65±1.28
Ours	**88.92±0.13**	**79.50±0.25**	**72.77±2.27**	**63.92±3.76**	**59.11±7.65**	**58.60±1.89**	**30.52±0.50**

Table 6: Transfer studies of our proposed method (topology-aware initialization and DE-guided rewiring) with deep vanilla-GCNs (32-layers). Comparisons are conducted on seven open-source widely adopted datasets with other previous state-of-the-art frameworks.

4 Other Related Works

Deep Graph Neural Networks. GNNs have established state-of-the-art performance in numerous real-world applications [9, 10, 11, 9, 5, 4, 12, 13, 14, 10, 15, 16]. While deep architectures imporves the representational power of neural networks, not every useful GNN has to be deep. Many real-world graph are "small-world"[54], where a node can reach any other node in a few hops, hence a few layers would suffice to provide the global coverage. However, when the graph data doesn't have small-world properties, or the related task requires long-range information, then deeper GNNs become very necessary. Many works [17, 18] revealed that when we start stacking spatial aggregations recursively in GNNs, the node representations will collapse to indistinguishable vectors and it will hamper the training of deep GNNs. Recently, there has been a series of techniques developed to handle the over-smoothing issue, which can be broadly categorized under skip connection, graph normalization, random dropping. Skip-connections [20, 21, 35, 17] have been applied to GNNs to exploit node embeddings from the preceding layers, to relieve the over-smoothing issue. Graph normalization techniques [25, 55, 56, 57, 58] re-scale node embeddings over an input graph to constrain pairwise node distance and thus alleviate over-smoothing. Dropout methods [59, 23, 24, 60] can be regarded as data augmentation methods, which help relieve both the over-fitting and over-smoothing issues in training very deep GNNs.

5 Conclusion

This paper makes an important step towards understanding the substandard performance of deep vanilla-GCNs from the signal propagation perspective and hypothesizes that by facilitating healthy gradient flow, we can significantly improve their trainability, as well as achieve state-of-the-art (SOTA) level performance or close using merely vanilla-GCNs. This paper derives a topology-aware isometric initialization scheme for vanilla-GCNs based on the principles of isometry. Additionally, this paper proposes to use Dirichlet Energy for the dynamic rewiring of vanilla-GCNs with skip-connections. Extensive experiments across multiple datasets illustrate that these methods improvise gradient flow in deep vanilla-GCNs and significantly boost their performance. An interesting direction for future work includes building theoretical relation between our proposed methods and gradient flow.

References

[1] Thomas Kipf and Max Welling. Semi-supervised classification with graph convolutional networks. *ArXiv*, abs/1609.02907, 2017.

[2] Michaël Defferrard, Xavier Bresson, and Pierre Vandergheynst. Convolutional neural networks on graphs with fast localized spectral filtering. *Advances in neural information processing systems*, 29, 2016.

[3] Petar Veličković, Guillem Cucurull, Arantxa Casanova, Adriana Romero, Pietro Lio, and Yoshua Bengio. Graph attention networks. *arXiv preprint arXiv:1710.10903*, 2017.

[4] Yuning You, Tianlong Chen, Zhangyang Wang, and Yang Shen. L2-gcn: Layer-wise and learned efficient training of graph convolutional networks. *2020 IEEE/CVF Conference on Computer Vision and Pattern Recognition (CVPR)*, pages 2124–2132, 2020.

[5] Hongyang Gao, Zhengyang Wang, and Shuiwang Ji. Large-scale learnable graph convolutional networks. *Proceedings of the 24th ACM SIGKDD International Conference on Knowledge Discovery & Data Mining*, 2018.

[6] Wei-Lin Chiang, Xuanqing Liu, Si Si, Yang Li, Samy Bengio, and Cho-Jui Hsieh. Cluster-gcn: An efficient algorithm for training deep and large graph convolutional networks. In *Proceedings of the 25th ACM SIGKDD International Conference on Knowledge Discovery & Data Mining*, pages 257–266, 2019.

[7] Jie Chen, Tengfei Ma, and Cao Xiao. Fastgcn: fast learning with graph convolutional networks via importance sampling. *arXiv preprint arXiv:1801.10247*, 2018.

[8] Kiran K Thekumparampil, Chong Wang, Sewoong Oh, and Li-Jia Li. Attention-based graph neural network for semi-supervised learning. *arXiv preprint arXiv:1803.03735*, 2018.

[9] Rex Ying, Ruining He, Kaifeng Chen, Pong Eksombatchai, William L Hamilton, and Jure Leskovec. Graph convolutional neural networks for web-scale recommender systems. In *Proceedings of the 24th ACM SIGKDD international conference on knowledge discovery & data mining*, pages 974–983, 2018.

[10] Junyuan Shang, Cao Xiao, Tengfei Ma, Hongyan Li, and Jimeng Sun. Gamenet: Graph augmented memory networks for recommending medication combination. *ArXiv*, abs/1809.01852, 2019.

[11] Lei Tang and Huan Liu. Relational learning via latent social dimensions. In *KDD*, 2009.

[12] Yuning You, Tianlong Chen, Zhangyang Wang, and Yang Shen. When does self-supervision help graph convolutional networks? *Proceedings of machine learning research*, 119:10871–10880, 2020.

[13] Alex Fout, Jonathon Byrd, Basir Shariat, and Asa Ben-Hur. Protein interface prediction using graph convolutional networks. In *NIPS*, 2017.

[14] Marinka Zitnik and Jure Leskovec. Predicting multicellular function through multi-layer tissue networks. *Bioinformatics*, 33:i190 – i198, 2017.

[15] Long Zhao, Xi Peng, Yu Tian, Mubbasir Kapadia, and Dimitris N. Metaxas. Semantic graph convolutional networks for 3d human pose regression. *2019 IEEE/CVF Conference on Computer Vision and Pattern Recognition (CVPR)*, pages 3420–3430, 2019.

[16] Yixiao Zhang, Xiaosong Wang, Ziyue Xu, Qihang Yu, Alan Loddon Yuille, and Daguang Xu. When radiology report generation meets knowledge graph. In *AAAI*, 2020.

[17] Qimai Li, Zhichao Han, and Xiao-Ming Wu. Deeper insights into graph convolutional networks for semi-supervised learning. In *Thirty-Second AAAI conference on artificial intelligence*, 2018.

[18] Hoang NT and Takanori Maehara. Revisiting graph neural networks: All we have is low-pass filters. *ArXiv*, abs/1905.09550, 2019.

[19] Uri Alon and Eran Yahav. On the bottleneck of graph neural networks and its practical implications. *ArXiv*, abs/2006.05205, 2021.

[20] Guohao Li, Matthias Müller, Ali K. Thabet, and Bernard Ghanem. Can gcns go as deep as cnns? *ArXiv*, abs/1904.03751, 2019.

[21] Ming Chen, Zhewei Wei, Zengfeng Huang, Bolin Ding, and Yaliang Li. Simple and deep graph convolutional networks. *ArXiv*, abs/2007.02133, 2020.

[22] Meng Liu, Hongyang Gao, and Shuiwang Ji. Towards deeper graph neural networks. *Proceedings of the 26th ACM SIGKDD International Conference on Knowledge Discovery & Data Mining*, 2020.

[23] Yu Rong, Wenbing Huang, Tingyang Xu, and Junzhou Huang. Dropedge: Towards deep graph convolutional networks on node classification. In *ICLR*, 2020.

[24] Wenbing Huang, Yu Rong, Tingyang Xu, Fuchun Sun, and Junzhou Huang. Tackling over-smoothing for general graph convolutional networks. *ArXiv*, abs/2008.09864, 2020.

[25] Lingxiao Zhao and Leman Akoglu. Pairnorm: Tackling oversmoothing in gnns. *ArXiv*, abs/1909.12223, 2020.

[26] Kuangqi Zhou, Yanfei Dong, Kaixin Wang, Wee Sun Lee, Bryan Hooi, Huan Xu, and Jiashi Feng. Understanding and resolving performance degradation in deep graph convolutional networks. *Proceedings of the 30th ACM International Conference on Information & Knowledge Management*, 2021.

[27] Weilin Cong, Morteza Ramezani, and Mehrdad Mahdavi. On provable benefits of depth in training graph convolutional networks. *Advances in Neural Information Processing Systems*, 34:9936–9949, 2021.

[28] Sitao Luan, Mingde Zhao, Xiao-Wen Chang, and Doina Precup. Training matters: Unlocking potentials of deeper graph convolutional neural networks. *arXiv preprint arXiv:2008.08838*, 2020.

[29] Xavier Glorot and Yoshua Bengio. Understanding the difficulty of training deep feedforward neural networks. In *AISTATS*, 2010.

[30] Kaiming He, X. Zhang, Shaoqing Ren, and Jian Sun. Delving deep into rectifiers: Surpassing human-level performance on imagenet classification. *2015 IEEE International Conference on Computer Vision (ICCV)*, pages 1026–1034, 2015.

[31] Jeffrey Pennington, Samuel S. Schoenholz, and Surya Ganguli. Resurrecting the sigmoid in deep learning through dynamical isometry: theory and practice. In *NIPS*, 2017.

[32] Samuel S. Schoenholz, Justin Gilmer, Surya Ganguli, and Jascha Sohl-Dickstein. Deep information propagation. *ArXiv*, abs/1611.01232, 2017.

[33] Kaiming He, X. Zhang, Shaoqing Ren, and Jian Sun. Deep residual learning for image recognition. *2016 IEEE Conference on Computer Vision and Pattern Recognition (CVPR)*, pages 770–778, 2016.

[34] Gao Huang, Zhuang Liu, and Kilian Q. Weinberger. Densely connected convolutional networks. *2017 IEEE Conference on Computer Vision and Pattern Recognition (CVPR)*, pages 2261–2269, 2017.

[35] Keyulu Xu, Chengtao Li, Yonglong Tian, Tomohiro Sonobe, Ken ichi Kawarabayashi, and Stefanie Jegelka. Representation learning on graphs with jumping knowledge networks. *ArXiv*, abs/1806.03536, 2018.

[36] Chen Cai and Yusu Wang. A note on over-smoothing for graph neural networks. *arXiv preprint arXiv:2006.13318*, 2020.

[37] Yujia Li, Daniel Tarlow, Marc Brockschmidt, and Richard Zemel. Gated graph sequence neural networks. *arXiv preprint arXiv:1511.05493*, 2015.

[38] Xavier Bresson and Thomas Laurent. Residual gated graph convnets. *arXiv preprint arXiv:1711.07553*, 2017.

[39] Aliaksei Sandryhaila and José MF Moura. Discrete signal processing on graphs. *IEEE transactions on signal processing*, 61(7):1644–1656, 2013.

[40] Ajay Kumar Jaiswal, Haoyu Ma, Tianlong Chen, Ying Ding, and Zhangyang Wang. Spending your winning lottery better after drawing it. *arXiv preprint arXiv:2101.03255*, 2021.

[41] Haozhi Qi, Chong You, Xiaolong Wang, Yi Ma, and Jitendra Malik. Deep isometric learning for visual recognition. In *International Conference on Machine Learning*, pages 7824–7835. PMLR, 2020.

[42] Antonio Ortega, Pascal Frossard, Jelena Kovačević, José MF Moura, and Pierre Vandergheynst. Graph signal processing: Overview, challenges, and applications. *Proceedings of the IEEE*, 106(5):808–828, 2018.

[43] Kathrin Schacke. On the kronecker product. *Master's thesis, University of Waterloo*, 2004.

[44] Johannes Klicpera, Aleksandar Bojchevski, and Stephan Günnemann. Predict then propagate: Graph neural networks meet personalized pagerank. In *ICLR*, 2019.

[45] Prithviraj Sen, Galileo Namata, Mustafa Bilgic, Lise Getoor, Brian Gallagher, and Tina Eliassi-Rad. Collective classification in network data articles. *AI Magazine*, 29:93–106, 09 2008.

[46] Tianlong Chen, Kaixiong Zhou, Keyu Duan, Wenqing Zheng, Peihao Wang, Xia Hu, and Zhangyang Wang. Bag of tricks for training deeper graph neural networks: A comprehensive benchmark study. *arXiv preprint arXiv:2108.10521*, 2021.

[47] Keyulu Xu, Chengtao Li, Yonglong Tian, Tomohiro Sonobe, Ken-ichi Kawarabayashi, and Stefanie Jegelka. Representation learning on graphs with jumping knowledge networks. In *International Conference on Machine Learning*, pages 5453–5462. PMLR, 2018.

[48] Anees Kazi, Shayan Shekarforoush, S Arvind Krishna, Hendrik Burwinkel, Gerome Vivar, Karsten Kortüm, Seyed-Ahmad Ahmadi, Shadi Albarqouni, and Nassir Navab. Inceptiongcn: receptive field aware graph convolutional network for disease prediction. In *International Conference on Information Processing in Medical Imaging*, pages 73–85. Springer, 2019.

[49] Felix Wu, Amauri Souza, Tianyi Zhang, Christopher Fifty, Tao Yu, and Kilian Weinberger. Simplifying graph convolutional networks. In *International conference on machine learning*, pages 6861–6871. PMLR, 2019.

[50] Liqi Yang, Linhan Luo, Lifeng Xin, Xiaofeng Zhang, and Xinni Zhang. Dagnn: Demand-aware graph neural networks for session-based recommendation. *arXiv preprint arXiv:2105.14428*, 2021.

[51] Eli Chien, Jianhao Peng, Pan Li, and Olgica Milenkovic. Adaptive universal generalized pagerank graph neural network. *arXiv preprint arXiv:2006.07988*, 2020.

[52] Oleksandr Shchur, Maximilian Mumme, Aleksandar Bojchevski, and Stephan Günnemann. Pitfalls of graph neural network evaluation. *arXiv preprint arXiv:1811.05868*, 2018.

[53] Hongbin Pei, Bingzhe Wei, Kevin Chen-Chuan Chang, Yu Lei, and Bo Yang. Geom-gcn: Geometric graph convolutional networks. *arXiv preprint arXiv:2002.05287*, 2020.

[54] Pablo Barceló, Egor V. Kostylev, Mikaël Monet, Jorge Pérez, Juan L. Reutter, and Juan Pablo Silva. The logical expressiveness of graph neural networks. In *ICLR*, 2020.

[55] Sergey Ioffe and Christian Szegedy. Batch normalization: Accelerating deep network training by reducing internal covariate shift. In *International conference on machine learning*, pages 448–456. PMLR, 2015.

[56] Kaixiong Zhou, Xiao Huang, Yuening Li, Daochen Zha, Rui Chen, and Xia Hu. Towards deeper graph neural networks with differentiable group normalization. *Advances in Neural Information Processing Systems*, 33:4917–4928, 2020.

[57] Yuxin Wu and Kaiming He. Group normalization. In *Proceedings of the European conference on computer vision (ECCV)*, pages 3–19, 2018.

[58] Chaoqi Yang, Ruijie Wang, Shuochao Yao, Shengzhong Liu, and Tarek Abdelzaher. Revisiting over-smoothing in deep gcns. *arXiv preprint arXiv:2003.13663*, 2020.

[59] Nitish Srivastava, Geoffrey E. Hinton, Alex Krizhevsky, Ilya Sutskever, and Ruslan Salakhutdinov. Dropout: a simple way to prevent neural networks from overfitting. *J. Mach. Learn. Res.*, 15:1929–1958, 2014.

[60] Difan Zou, Ziniu Hu, Yewen Wang, Song Jiang, Yizhou Sun, and Quanquan Gu. Layer-dependent importance sampling for training deep and large graph convolutional networks. In *NeurIPS*, 2019.

Checklist

The checklist follows the references. Please read the checklist guidelines carefully for information on how to answer these questions. For each question, change the default **[TODO]** to [Yes] , [No] , or [N/A] . You are strongly encouraged to include a **justification to your answer**, either by referencing the appropriate section of your paper or providing a brief inline description. For example:

- Did you include the license to the code and datasets? [Yes] See Section **??**.
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1. For all authors...
   (a) Do the main claims made in the abstract and introduction accurately reflect the paper's contributions and scope? [Yes]
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   (a) Did you include the code, data, and instructions needed to reproduce the main experimental results (either in the supplemental material or as a URL)? [Yes]
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   (e) Did you discuss whether the data you are using/curating contains personally identifiable information or offensive content? [N/A]
5. If you used crowdsourcing or conducted research with human subjects...
   (a) Did you include the full text of instructions given to participants and screenshots, if applicable? [N/A]
   (b) Did you describe any potential participant risks, with links to Institutional Review Board (IRB) approvals, if applicable? [N/A]
   (c) Did you include the estimated hourly wage paid to participants and the total amount spent on participant compensation? [N/A]

A Derivation of Eq. 12 and 13

Recall that, in Section 2.3 our goal is to make initialized GCN isometric. Instead of exactly compute such initialization by solving an optimization, we propose the following objective, which has a closed-form solution:

$$\boldsymbol{W}^* = \underset{\boldsymbol{W}\in\mathbb{R}^{C'\times C}}{\operatorname{argmin}} \sum_{i,j\in[N]} \left\| \left(\boldsymbol{a}_i\boldsymbol{a}_j^\top\right) \otimes \left(\boldsymbol{W}^\top\boldsymbol{W}\right) - \left(\boldsymbol{e}_i\boldsymbol{e}_j^\top\right) \otimes \boldsymbol{I}_C \right\|_F^2 . \tag{17}$$

In this section, we give the detailed derivation to reach the closed-form solution. Notice that both $\boldsymbol{a}_i\boldsymbol{a}_j^\top$ and $\boldsymbol{e}_i\boldsymbol{e}_j^\top$ are sparsely structured. Let $\boldsymbol{A}^{(i,j)} = \boldsymbol{a}_i\boldsymbol{a}_j^\top$ and $\boldsymbol{E}^{(i,j)} = \boldsymbol{e}_i\boldsymbol{e}_j^\top$, then we have:

$$\boldsymbol{A}^{(i,j)}_{m,n} = \begin{cases} 1 & (i,m)\in\mathcal{E} \text{ and } (j,n)\in\mathcal{E} \\ 0 & \text{Otherwise} \end{cases}, \qquad \boldsymbol{E}^{(i,j)}_{m,n} = \begin{cases} 1 & m=i \text{ and } n=j \\ 0 & \text{Otherwise} \end{cases}, \tag{18}$$

where $\boldsymbol{A}^{(i,j)}_{m,n}$, $\boldsymbol{E}^{(i,j)}_{m,n}$ denotes the m-th row and the n-th column entry of $\boldsymbol{A}^{(i,j)}$, $\boldsymbol{E}^{(i,j)}$, respectively. Hence, for $m,n\in[N]$ such that $(i,m)\in\mathcal{E}$ and $(j,n)\in\mathcal{E}$ but $m\neq i$ or $n\neq j$, the block-wise difference at location (m,n) remains $\boldsymbol{W}^\top\boldsymbol{W}$. While for $m=i$ and $n=j$, the block difference is $\boldsymbol{W}^\top\boldsymbol{W}-\boldsymbol{I}_C$ as $\boldsymbol{A}$ contains self-loop (i.e., $(i,i)\in\mathcal{E}$ and $(j,j)\in\mathcal{E}$). Then we can simply our objective Eq. 17 as below:

$$\boldsymbol{W}^* = \underset{\boldsymbol{W}\in\mathbb{R}^{C'\times C}}{\operatorname{argmin}} \sum_{i,j\in[N]} \left(\sum_{\substack{m,n\in[N]\\(m,n)\neq(i,j)}} \left\|\boldsymbol{A}^{(i,j)}_{m,n}\boldsymbol{W}^\top\boldsymbol{W}\right\|_F^2 + \left\|\boldsymbol{W}^\top\boldsymbol{W}-\boldsymbol{I}_C\right\|_F^2 \right) \tag{19}$$

$$= \underset{\boldsymbol{W}\in\mathbb{R}^{C'\times C}}{\operatorname{argmin}} \sum_{i,j\in[N]} (d_i+d_j+d_id_j) \left\|\boldsymbol{W}^\top\boldsymbol{W}\right\|_F^2 + N^2\left\|\boldsymbol{W}^\top\boldsymbol{W}-\boldsymbol{I}_C\right\|_F^2 . \tag{20}$$

To solve Eq. 20, we decompose the diagonal and off-diagonal components from $\boldsymbol{W}^\top\boldsymbol{W}$. For off-diagonal terms, we can minimize them to zeros, thus we have $\boldsymbol{w}_k^\top\boldsymbol{w}_l = 0$, where we note that $\boldsymbol{w}_k$ denotes the k-th column of $\boldsymbol{W}$. For on-diagonal terms, it is equivalent to minimizing $\sum_{i,j\in[N]}(d_i+d_j+d_id_j)\|\boldsymbol{w}_k\|_2^4 + N^2(\|\boldsymbol{w}_k\|_2^2-1)^2$ for every $k\in[C]$. Then combining both arguments above, the optimal solution should satisfy:

$$\boxed{\begin{aligned} &\|\boldsymbol{w}_k\|_2^2 = \frac{N^2}{\sum_{i,j\in[N]}(d_i+1)(d_j+1)} && \forall k\in[C], \quad (21)\\ &\boldsymbol{w}_k^\top\boldsymbol{w}_l = 0 && \forall k,l\in[C], k\neq l, \quad (22)\end{aligned}}$$

which reaches our final conclusion in Eq. 14. To compute this magnitude given a graph, one needs to first compute the degree of each node, and plug them into our Eq. 12. The complexity to compute this value is as small as $\mathcal{O}(N^2)$, which is significantly lower than directly optimizing Eq. 10.

B Description of Methods in Table 3:

Table 3 presents comparison of our DE-Guided Rewiring dynamic rewiring with respect to various fancy methods to improve the trainability of deep vanilla-GCNs [1] such as as skip-connections, regularization. Our formal description of these methods are:

B.1 Skip Connections

Skip-connections [20, 21, 35, 17] helps in allevating the problem of over-smoothing and significantly improve the accuracy and training stability of deep GCNs. For a L layer GCN, we can apply various type of skip-connections after certain graph convolutional layers with the current and preceding embeddings X^l , $0\le l\le L$. We have compared our method with following four representative types of skip connections:

1. *Residual Connection:* $X^l = (1-\alpha)\cdot X^l + \alpha\cdot X^{l-1}$

2. *Initial Connection:* $X^l = (1-\alpha) \cdot X^l + \alpha \cdot X^0$
3. *Dense Connection:* $X^l = \text{COM}(\{X^k, 0 \leq k \leq l\})$
4. *Jumping Connection:* $X^L = \text{COM}(\{X^k, 0 \leq k \leq L\})$

where α is *residual* and *initial connections* is a hyperparameter to weight the contribution of a node features from the current layer l and previous layers. *Jumping connection* is a simplified case of *dense connection* and it is only applied at the end of whole forward propagation process to combine the node features from all previous layers.

B.2 Graph Normalization

Graph normalization [25, 55, 56, 57, 58] re-scale node embeddings over an input graph to constrain pairwise node distance and thus alleviate over-smoothing. Our investigated normalization mechanisms are formally depicted as follows:

1. *BatchNorm:* $x_{(k)} = \gamma \cdot \frac{x_{(k)} - \mathbf{E}[x_{(k)}]}{std(x_{(k)})} + \beta$
2. *PairNorm:* $\tilde{x}_i = x_i - \frac{1}{n}\sum_{i=1}^{n} x_i$, $\text{PairNorm}(x_i;\ s) = \frac{s \cdot \tilde{x}_i}{(\frac{1}{n}\sum_{i=1}^{n} \|\tilde{x}_i\|_2^2)^{1/2}}$
3. *NodeNorm:* $\text{NodeNorm}(x_i;\ p) = \frac{x_i}{std(x_i)^{1/2}}$

where s in *PairNorm* is a hperparameter controlling the average pair-wise variance and p in *NodeNorm* denotes the normalization order.

B.3 Dataset Details

Table 7 provided provides the detailed properties and download links for all adopted datasets. We adopt the following benchmark datasets since i) they are widely applied to develop and evaluate GNN models, especially for deep GNNs studied in this paper; ii) they contains diverse graphs from small-scale to large scale or from homogeneous to heterogeneous; iii) they are collected from different applications including citation network, social network, etc.

Dataset	Nodes	Edges	Features	Classes	Download Links
Cora	2,708	5,429	1,433	7	https://github.com/kimiyoung/planetoid/raw/master/data
Citeseer	3,327	4,732	3,703	6	https://github.com/kimiyoung/planetoid/raw/master/data
PubMed	19,717	44,338	500	3	https://github.com/kimiyoung/planetoid/raw/master/data
OGBN-ArXiv	169,343	1,166,243	128	40	https://ogb.stanford.edu/
CoauthorCS	18,333	81,894	6805	15	https://github.com/shchur/gnn-benchmark/raw/master/data/npz/
Computers	13,381	245,778	767	10	https://github.com/shchur/gnn-benchmark/raw/master/data/npz/
Photo	7,487	119,043	745	8	https://github.com/shchur/gnn-benchmark/raw/master/data/npz/
Texas	183	309	1,703	5	https://raw.githubusercontent.com/graphdml-uiuc-jlu/geom-gcn/master
Wisconsin	183	499	1,703	5	https://raw.githubusercontent.com/graphdml-uiuc-jlu/geom-gcn/master
Cornell	183	295	1,703	5	https://raw.githubusercontent.com/graphdml-uiuc-jlu/geom-gcn/master
Actor	7,600	33,544	931	5	https://raw.githubusercontent.com/graphdml-uiuc-jlu/geom-gcn/master

Table 7: Graph datasets statistics and download links.