DISENTANGLING DEEP NEURAL NETWORKS WITH RECTIFIED LINEAR UNITS USING DUALITY

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ABSTRACT

Despite their success deep neural networks (DNNs) are still largely considered as black boxes. The main issue is that the linear and non-linear operations are entangled in every layer, making it hard to interpret the hidden layer outputs. In this paper, we look at DNNs with rectified linear units (ReLUs), and focus on the gating property (‘on/off’ states) of the ReLUs. We extend the recently developed dual view in which the computation is broken path-wise to show that learning in the gates is more crucial, and learning the weights given the gates is characterised analytically via the so called neural path kernel (NPK) which depends on inputs and gates. In this paper, we present novel results to show that convolution with global pooling and skip connection provide respectively rotational invariance and ensemble structure to the NPK. To address ‘black box’-ness, we propose a novel interpretable counterpart of DNNs with ReLUs namely deep linearly gated networks (DLGN): the pre-activations to the gates are generated by a deep linear network, and the gates are then applied as external masks to learn the weights in a different network. The DLGN is not an alternative architecture per se, but a disentanglement and an interpretable re-arrangement of the computations in a DNN with ReLUs. The DLGN disentangles the computations into two ‘mathematically’ interpretable linearities (i) the ‘primal’ linearity between the input and the pre-activations in the gating network and (ii) the ‘dual’ linearity in the path space in the weights network characterised by the NPK. We compare the performance of DNN, DGN and DLGN on CIFAR-10 and CIFAR-100 to show that, the DLGN recovers more than 83.5% of the performance of state-of-the-art DNNs, i.e., while entanglement in the DNNs enable their improved performance, the ‘disentangled and interpretable’ computations in the DLGN recovers most part of the performance. This brings us to an interesting question: ‘Is DLGN a universal spectral approximator?’

1 Introduction

Despite their success deep neural networks (DNNs) are still largely considered as black boxes. The main issue is that in each layer of a DNN, the linear computation, i.e., multiplication by the weight matrix and the non-linear activations are entangled. Such entanglement has its pros and cons. The commonly held view is that such entanglement is the key to success of DNNs, in that, it allows DNNs to learn sophisticated structures in a layer-by-layer manner. However, in terms of interpretability, such entanglement has an adverse effect: only the final layer is linear and amenable to a feature/weight interpretation, and the hidden layers are non-interpretable due to the non-linearities.

Prior works [Jacot et al., 2018, Arora et al., 2019, Cao and Gu, 2019] showed that training an infinite width DNN with gradient descent is equivalent to a kernel method with the so called neural tangent kernel matrix. As a pure kernel method, the NTK matrix performed better than other pure kernel methods. However, in relation to ‘black box’-ness, there are two issues with NTK theory: (i) Issue I: Infinite width NTK matrix does not explain fully the success of DNNs because it was observed that finite width DNNs outperform their infinite width NTK counterparts, and it was an open question to understand this performance gap [Arora et al., 2019], and (ii) Issue II: Since the NTK is based on the gradients, it does not offer further insights about the inner workings of DNNs even for infinite width.
A dual view for DNNs with rectified linear units (ReLUs) was recently developed by Lakshminarayanan and Singh [2020] who exploited the gating property (i.e., “on/off” states) of the ReLUs. The dual view is essentially linearity in the path space, i.e., the output is the summation of path contributions. While the weights in a path are the same for each input, whether or not a path contributes to the output is entirely dictated by the gates in the path, which are ‘on/off’ based on the input. To understand the role of the gates, a deep gated network (DGN) (see Figure 1) was used to disentangle the learning in the gates from the learning in weights. In a DGN, the gates are generated (and learnt) in a ‘gating network’ which is a DNN with ReLUs and are applied as external signals and the weights are learnt in a ‘weight network’ consisting of gated linear units (GaLUs) [Fiat et al., 2019]. Each GaLU multiplies its pre-activation and the external gating signal. Using the DGN, two important insights were provided: (i) learning in the gates is the most crucial for finite width networks to outperform infinite width NTK; this addresses Issue I, and (ii) in the limit of infinite width, learning the weights with fixed gates, the NTK is equal to (but for a scalar) a so called neural path kernel (NPK) which is a kernel solely based on inputs and gates. This shifts Issue II on interpretability to that of interpretability of the gates as opposed to interpretability of the gradients.

Our Contribution. We extend the dual view to address ‘black box’-ness by completely disentangling the ‘gating network’ and the ‘weight network’. Our contributions are listed below.

- **Disentangling Gating Network.** For this, we propose a novel Deep Linearly Gated Network (DLGN) as a mathematically interpretable counterpart of a DNN with ReLUs (see Figure 1). In a DLGN, the gating network is a deep linear network, i.e., there is disentanglement because of the absence of non-linear activations. The gating network is mathematically interpretable, because, the transformations from input to the pre-activations are entirely linear; we call this primal linearity.

- **Dual View** (Section 3.1). We present an unnoticed insight in prior work on fully connected networks that the NPK is a product kernel and is invariant to layer permutations. We present new results to show that (i) the NPK is rotationally invariant for convolutional networks with global average pooling, and (ii) the NPK is an ensemble of many kernels in the presence of skip connections.

- **Disentangling Weight Network.** We then argue via theory and experiments that the weight network is disentangled in the path space, i.e., it learns path-by-path and not layer-by-layer. For this, in Section 3.2 we show via experiments that destroying the layer-by-layer structure by permuting the layers and providing a constant 1 as input (see DLGN in Figure 1) do not degrade performance. These counter intuitive results are difficult to reconcile using the commonly held ‘sophisticated structures are learnt in layers’ interpretation. However, these experimental results follow from the theory in Section 3.1. In other words, it is useful to think that the learning in the weight network happens path-by-path; we call this dual linearity, which (for infinite width) is interpreted via the NPK.

**Message.** The DLGN is not an alternative architecture per se, but a disentanglement and an interpretable rearrangement of the computations in a DNN with ReLUs. The DLGN disentangles the computations into two ‘mathematically’ interpretable lineairities (i) the ‘primal’ linearity and (ii) the ‘dual’ linearity interpreted via the NPK. Using the facts that the NPK is based on input and the gates, and in a DLGN, the pre-activations in the gating network are ‘primal’ linear, we have complete disentanglement. We compare the performance of DNN, DGN and DLGN on CIFAR-10 and CIFAR-100 to show that, the DLGN recovers more than 83.5% of the performance of state-of-the-art DNNs, i.e., while entanglement in the DNNs enable their improved performance, the ‘disentangled and interpretable’ computations in the DLGN recovers most part of the performance. This brings us to an interesting question: ‘Is DLGN a universal spectral approximator?’ (see Section 3.2).

### 1.1 Related Works.

We now compare our work with the related works.
• **Kernels.** Several works have examined theoretically as well as empirically two important kernels associated with a DNN namely its NTK based on the correlation of the gradients and the conjugate kernel based on the correlation of the outputs [Fan and Wang, 2020, Geifman et al., 2020, Liu et al., 2020, Chen et al., 2020, Xiao et al., 2020, Jacot et al., 2018, Arora et al., 2019, Novak et al., 2018, Lee et al., 2017, 2020a]. In contrast, the NPK is based on the correlation of the gates. We do not build pure-kernel method with NPK, but use it as an aid to disentangle finite width DNN with ReLUs.

• **ReLU, Gating, Dual Linearity.** A spline theory based on max-affine linearity was proposed in [Balestriero et al., 2018, Balestriero and Baraniuk, 2018] to show that a DNN with ReLUs performs hierarchical, greedy template matching. In contrast, the dual view exploits the gating property to simplify the NTK into the NPK. Gated linearity was studied in [Fiat et al., 2019] for single layered networks, along with a non-gradient algorithm to tune the gates. In contrast, we look at networks of any depth, and the gates are tuned via standard optimisers. The main novelty in our work in contrast to the above is that in DLGN the feature generation is linear. The gating in this paper refers to the gating property of the ReLU itself and has no connection to [Srivastava et al., 2015] where gating is a mechanism to regulate information flow. Also, the soft-gating used in our work and in [Lakshminarayanan and Singh, 2020] enables gradient flow via the gating network and is different from Swish [Ramachandran et al., 2018], which is the multiplication of pre-activation and sigmoid.

• **Finite vs Infinite Width.** [Lee et al., 2020b] perform an extensive comparison of finite versus infinite width DNNs. An aspect that is absent in their work, but present in the dual view is the disentanglement of gates and weights, and the fact that the learning in gates is crucial for finite width network to outperform infinite width DNNs. In our paper, we make use of theory developed for infinite width DNNs to provide empirical insights into inner workings of finite width networks.

• **Capacity.** Our experiments on destruction of layers, and providing constant 1 input are direct consequences of the insights from dual view theory. These are not explained by mere capacity based studies showing DNNs are powerful to fit even random labelling of datasets [Zhang et al., 2016].

## 2 Prior Work : Neural Tangent Kernel and Dual View

In this section, we will focus on the dual view [Lakshminarayanan and Singh, 2020] and how the dual view helps to address the open question in the NTK theory. We begin with a brief summary of NTK.

### 2.1 Infinite Width DNN = Kernel Method With Neural Tangent Kernel

An important kernel associated with a DNN is its *neural tangent kernel* (NTK), which, for a pair of input examples \( x, x' \in \mathbb{R}^d \), and network weights \( \Theta \in \mathbb{R}^{d_{\text{net}}} \), is given by:

\[
\text{NTK}(x, x') = \left\langle \nabla_{\Theta} \hat{y}(x), \nabla_{\Theta} \hat{y}(x') \right\rangle,
\]

where \( \hat{y}_\Theta(\cdot) \in \mathbb{R} \) is the DNN output. Prior works [Jacot et al., 2018, Arora et al., 2019, Cao and Gid, 2019] have shown that, as the width of the DNN goes to infinity, the NTK matrix converges to a limiting deterministic matrix \( \text{NTK}_\infty \), and training an infinitely wide DNN is equivalent to a kernel method with \( \text{NTK}_\infty \). While, as a pure kernel \( \text{NTK}_\infty \) performed better than prior kernels by more than 10%, [Arora et al., 2019] observed that on CIFAR-10:

\[
\text{CNTK-GAP: } 77.43\% \leq \text{CNN-GAP: } 83.30\%
\]

where, CNN-GAP is a convolutional neural network with global average pooling and CNTK-GAP is its corresponding \( \text{NTK}_\infty \) matrix. Due to this performance gap of about 5 − 6%, they concluded that \( \text{NTK}_\infty \) does not explain fully the success of DNNs, and explaining this gap was an open question.

### 2.2 Dual View For DNNs with ReLUs: Characterising the role of gates

In the dual view, the computations are broken down path-by-path. The input and the gates (in each path) are encoded in a neural path feature vector and the weights (in each path) are encoded in a neural path value vector, and the output is the inner product of these two vectors. The learning in the gates and the learning in the weights are separated in a deep gated network (DGN) setup, which leads to the two main results of dual view presented in Section 2.2.3 and Section 2.2.4, wherein, the neural path kernel, the Gram matrix of the neural path features will play a key role.
The ‘gating network’ has ReLUs which turn ‘on/off’ based on their pre-activation signals, and the ‘weight network’ has gated linear units (GaLUs) [Fiat et al., 2019, Lakshminarayanan and Singh, 2020], which multiply their respective outputs. In the case of fixed random gates, the feature network is initialised at random and frozen, and the value network is trained with \( \hat{y}_{DGN} \) as the output. In the case of fixed random gates, the feature network is initialised and the random flow is frozen, and the value network is trained with \( \hat{y}_{DGN} \) as the output. In the case of fixed gates, hard gating \( G(q) = \mathbb{1}_{\{q \neq 0\}} \) is used. Standalone Training: both feature and value network are initialised at random and trained together with \( \hat{y}_{DGN} \) as the output. Here, soft gating \( G(q) = \frac{1}{1 + \exp(-\beta q)} \) is used to allow gradient flow through feature network. On the right side is the CNN-GAP and its DGN used in [Lakshminarayanan and Singh, 2020]. In CNN-GAP, \( C_1, C_2, C_3, C_4 \) are convolutional layers, \( FC_1, FC_2 \) are fully connected layers. In CNN-GAP-DGN, \( G_{i}, i = 1, 2, 3, 4 \) are the gates of layers, the superscripts \( f \) and \( v \) stand for feature and value network respectively.

### 2.2.1 Neural Path Feature, Neural Path Value and Neural Path Kernel

Consider a fully connected DNN with ‘\( d \)’ layers and ‘\( w \)’ hidden units in each layer. Let the DNN accept input \( x \in \mathbb{R}^{d_n} \) and produce an output \( \hat{y}_{\Theta}(x) \in \mathbb{R} \).

**Definition 2.1.** A path starts from an input node, passes through a weight and a hidden unit in each layer and ends at the output node. We define the following quantities for a path \( p \):

- **Activity:** \( A_\Theta(x, p) \) is the product of the ‘\( d-1 \)’ gates in the path.
- **Value:** \( v_\Theta(p) \) is the product of the ‘\( d \)’ weights in the path. The neural path feature (NPF) given by \( \phi_\Theta(x) = (\phi_\Theta(x, p), p = 1, \ldots, P^{fc}) \in \mathbb{R}^{P^{fc}} \) and the neural path value (NPV) given by \( v_\Theta = (v_\Theta(p), p = 1, \ldots, P^{fc}) \in \mathbb{R}^{P^{fc}} \), where \( P^{fc} = d_{in}w^{(d-1)} \) is the total number of paths.

**Proposition 2.1.** The output of the DNN is then the inner product of the NPF and NPV:

\[
\hat{y}_\Theta(x) = \langle \phi_\Theta(x), v_\Theta \rangle = \sum_{p \in [P]} \phi_\Theta(x, p)v_\Theta(p) \tag{1}
\]

**Subnetwork Interpretation of DNNs with ReLUs.** A path is active only if all the gates in the path are active. This gives a subnetwork interpretation, i.e., for a given input \( x \in \mathbb{R}^{d_n} \), only a subset of the gates and consequently only a subset of the paths are active, and the input to output computation can be seen to be produced by this active subnetwork. The following matrix captures the correlation of the active subnetworks for a given pair of inputs \( x, x' \in \mathbb{R}^{d_n} \).

**Definition 2.2 (Overlap of active sub-networks).** The total number of ‘active’ paths for both \( x \) and \( x' \) that pass through input node \( i \) is defined to be:

\[
\text{overlap}_\Theta(i, x, x') \overset{\Delta}{=} |\{ p: p \text{ starts at node } i, A_\Theta(x, p) = A_\Theta(x', p) = 1 \}|
\]

**Lemma 2.1 (Neural Path Kernel (NPK)).** Let \( D \in \mathbb{R}^{d_n} \) be a vector of non-negative entries and for \( u, u' \in \mathbb{R}^{d_n} \), let \( \langle u, u' \rangle_D = \sum_{i=1}^{d_n} D(i)u(i)u'(i) \). Then the neural path kernel (NPK) is given by:

\[
\text{NPK}_\Theta(x, x') \overset{\Delta}{=} \langle \phi_\Theta(x), \phi_\Theta(x') \rangle = \langle x, x' \rangle_{\text{overlap}_\Theta(., x, x')}
\]

### 2.2.2 Deep Gated Network: Separating Gates and Weights

Deep Gated Network (DGN) is a setup to separate the gates from the weights. Consider a DNN with ReLUs with weights \( \Theta \in \mathbb{R}^{d_{out}} \). The DGN corresponding to this DNN (left diagram in Figure 2) has two networks of identical architecture (to the DNN) namely the ‘gating network’ and the ‘weight network’ with distinct weights \( \Theta^f \in \mathbb{R}^{d_{out}} \) and \( \Theta^v \in \mathbb{R}^{d_{out}} \).

The ‘gating network’ has ReLUs which turn ‘on/off’ based on their pre-activation signals, and the ‘weight network’ has gated linear units (GaLUs) [Fiat et al., 2019, Lakshminarayanan and Singh, 2020], which multiply their respective

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Figure 2: Shows the DGN on the left. **Training:** In the case of fixed learnt gates, the feature network is pre-trained using \( \hat{y}_f \) as the output, and then the feature network is frozen, and the value network is trained with \( \hat{y}_{DGN} \) as the output. In the case of fixed random gates, the feature network is initialised and the random flow is frozen, and the value network is trained with \( \hat{y}_{DGN} \) as the output. In the case of fixed gates, hard gating \( G(q) = \mathbb{1}_{\{q \neq 0\}} \) is used. **Standalone Training:** both feature and value network are initialised at random and trained together with \( \hat{y}_{DGN} \) as the output. Here, soft gating \( G(q) = \frac{1}{1 + \exp(-\beta q)} \) is used to allow gradient flow through feature network. On the right side is the CNN-GAP and its DGN used in [Lakshminarayanan and Singh, 2020]. In CNN-GAP, \( C_1, C_2, C_3, C_4 \) are convolutional layers, \( FC_1, FC_2 \) are fully connected layers. In CNN-GAP-DGN, \( G_{i}, i = 1, 2, 3, 4 \) are the gates of layers, the superscripts \( f \) and \( v \) stand for feature and value network respectively.
pre-activation inputs by the external gating signals provided by the ‘gating network’. Since both the networks have identical architecture, the ReLUs and GalUs in the respective networks have a one-to-one correspondence. Gating network realises \( \hat{\phi_{\Theta}}(x) \) by turning ‘on/off’ the corresponding GalUs in the weight network. The weight network realises \( v_{\Theta'} \) and computes the output \( \hat{y}_{\text{DGN}}(x) = \langle \hat{\phi_{\Theta}}(x), v_{\Theta'} \rangle \). The gating network is also called as the feature network since it realises the neural path features, and the weight network is also called as the value network since it realises the neural path value.

### 2.2.3 Learning Weights With Fixed Gates = Neural Path Kernel

During training, a DNN learns both \( \hat{\phi_{\Theta}}(x) \) as well as \( v_{\Theta} \) simultaneously, and a finite time characterisation of this learning in finite width DNNs is desirable. However, this is a hard problem. An easier problem is to understand in a DGN, how the weights in the value network are learnt when the gates are fixed in the feature network, i.e., how \( \hat{y}_{\text{DGN}}(x) = \langle \hat{\phi_{\Theta}}(x), v_{\Theta'} \rangle \) is learnt by learning \( v_{\Theta'} \) with fixed \( \hat{\phi_{\Theta}}(x) \). While \( \hat{y}_{\text{DGN}}(x) = \langle \hat{\phi_{\Theta}}(x), v_{\Theta'} \rangle \) is linear in the dual variables, it is still non-linear in the value network weights \( \Theta' \). However, [Lakshminarayanan and Singh, 2020] showed that the dual linearity is characterised by the NPK in the infinite width regime. We state the assumption followed by Theorem 5.1 in [Lakshminarayanan and Singh, 2020], wherein, the NTK \( (x, x') = (\nabla_{\Theta'} y_{\Theta_0}(x), \nabla_{\Theta'} y_{\Theta_0}(x')) \) is due to the gradient of \( \hat{y}_{\text{DGN}} \) with respect to the value network weights, and the NPK \( (x, x') = \langle \hat{\phi_{\Theta_0}}(x), \hat{\phi_{\Theta_0}}(x') \rangle \) is due to the feature network weights.

**Assumption 2.1.** \( \Theta_{\Theta_0} \sim \text{Bernoulli}(\frac{1}{2}) \) over \( \{-\sigma, +\sigma\} \) and statistically independent of \( \Theta_{\Theta_0'} \).

**Theorem 2.1** (Theorem 5.1 in [Lakshminarayanan and Singh, 2020]). Under Assumption 2.1 for a fully connected DGN:

\[
\text{NTK}_{\text{FC}}(x, x') \xrightarrow{(a) \ d \cdot \sigma^2 d^{-(d-1)}} \text{NPK}_{\text{FC}}(x, x'), \quad \text{as} \ w \to \infty
\]

\[
= d \cdot \sigma^2 d^{-(d-1)} \cdot \langle x, x' \rangle \cdot \text{overlap}(x, x')
\]

**Remark.** In the fully connected case, overlap \( (i, x, x') \) is identical for all \( i = 1, \ldots, d_{\Theta_0} \), and hence \( \langle x, x' \rangle \text{overlap}(i, x, x') \) in Lemma 2.1 becomes \( \langle x, x' \rangle \cdot \text{overlap}(x, x') \) in Theorem 2.1. It follows from NTK theory that an infinite width DGN with fixed gates is equivalent to kernel method with NPK.

### 2.2.4 Learning in Gates Key For Finite Width To Be Better Than Infinite Width

The fixed gates setting is an idealised setting, in that, it does not theoretically capture the learning of the gates, i.e., the neural path features \( \hat{\phi_{\Theta}}(x) \). However, the learning in the gates can be empirically characterised by comparing fixed learnt (FL) gates coming from a pre-trained DNN and fixed random (FR) gates coming from randomly initialised DNN, and the infinite width NTK. Using a CNN-GAP and its corresponding DGN, [Lakshminarayanan and Singh, 2020] showed on CIFAR-10 that (see Figure 2 for details on DGN training and the CNN-GAP architecture):

**FR Gates: 67.09% ≤ CNTK-GAP: 77.43% ≤ FL Gates: 79.68% ≈ CNN-GAP: 80.32%**

based on which it was concluded that learning in the gates (i.e., neural path features) is crucial for finite width CNN-GAP to outperform the infinite width CNTK-GAP. It was also shown that the DGN can be trained standalone (as shown in Figure 2) and is only marginally poor to the DNN.

### 3 Deep Linearly Gated Networks: Complete Disentanglement

\[
\text{DLGN: } x \rightarrow \text{Linear} \rightarrow \text{Pre-activations} \rightarrow \text{Gates} \rightarrow \hat{\phi_{\Theta}}(x) \rightarrow \text{Linear: } \hat{y}(x) = \langle \hat{\phi_{\Theta}}(x), v_{\Theta'} \rangle
\]

The deep linearly gated network (DLGN) has two ‘mathematically’ interpretable linearities, the ‘primal’ and the ‘dual’ linearities. The primal linearity is ensured in via construction and needs no theoretical justification. Once the pre-activations triggers gates, \( \hat{\phi_{\Theta}}(x) \) gets realised in the value network by activating the paths. Now, the value network itself is ‘dual’ linear, i.e., it simply computes/learns the inner product \( \hat{y}(x) = \langle \hat{\phi_{\Theta}}(x), v_{\Theta'} \rangle \). Gating lifts the ‘primal’ linear computations in the feature network to ‘dual’ linear computations in the value network. Dual linearity is characterised by the NPK (for infinite width) which in turn depends on the input and gates, and the fact that the pre-activations to the gates are primal linear implies complete disentanglement and interpretability.

Dual linearity is mathematically evident due to the inner product relationship, however, adopting it has the following conceptual issue: it is a commonly held view that ‘sophisticated features are learnt in the layers’, that is, given that the input \( x \in \mathbb{R}^{d_{\Theta_0}} \) is presented to the value network (as in Figure 2), it could be argued that the GalUs and linear operations
are entangled which in turn enable learning of sophisticated features in the layers. In what follows, we demystify this layer-by-layer view via theory (infinite width case) in Section 3.1, and experiments (on finite width networks) in Section 3.2 and then study the performance of DLGN in Section 3.2. The layer-by-layer view is demystified by showing that (i) a constant 1 input can be given to the value network, (ii) layer-by-layer structure can be destroyed. The constant 1 input is meant to show that if the input is not given to the value network then it is not possible to learn sophisticated structures from the input in a layer-by-layer manner. In terms of the dual linearity, providing a constant 1 input has only a minor impact, in that, the neural path feature becomes \( \phi(x, p) = 1 \cdot A(x, p) \), i.e., it still encodes the path activity which is still input dependent. Since \( \phi(x) \) depends only on gates, the NPK will depend only on the overlap matrix; results in Section 3.1 captures this in theory. Now, it could be argued that, despite a constant 1 input, the gates are still arranged layer-by-layer, due to which, the value network is still able to learn sophisticated structures in its layers. Section 3.1 has theory that points out that as long as the correlation of the gates is not lost, the layer-by-layer structure can be destroyed.

3.1 Dual Linearity: New Insights and New Results

We saw in Section 2.2.3 that dual linearity is characterised by the NPK for infinite width case. In this section, we: (i) cover standard architectural choices namely convolutions with global-average-pooling and skip connections in Theorems 3.2 and 3.3, the prior result Theorem 2.1 is only for the fully connected case, (ii) present new insights on Theorem 2.1 by restating it explicitly in terms of the gates in Theorem 3.1 and (iii) discuss how the NPK structure helps in demystifying the layer-by-layer view. Note: Results in this section are about the value network and hold for both DGN and DLGN.

3.1.1 Fully Connected: Product of LayerWise Base Kernels

**Theorem 3.1.** Let \( G_l(x) \in [0, 1]^w \) denote the gates in layer \( l \in \{1, \ldots, d - 1\} \) for input \( x \in \mathbb{R}^d_w \). Under Assumption 2.7 \((\sigma = \sqrt{\frac{d}{w}})\) as \( w \to \infty \), we have for fully connected DGN/DLGN:

\[
NTK_{FC}(x, x') \to d \cdot \sigma^{2d-1} \cdot NPK_{FC}(x, x') = d \cdot \sigma^{2d-1} \cdot \langle x, x' \rangle \cdot \prod_{l=1}^{d-1} \frac{\langle G_l(x), G_l(x') \rangle}{w},
\]

- **Product Kernel: Role of Depth and Width.** Theorem 3.1 is mathematically equivalent to Theorem 2.1 which follows from the observation that \( \text{overlap}(x, x') = \prod_{l=1}^{d-1} \langle G_l(x), G_l(x') \rangle \). While this observation is very elementary in itself, it is significant at the same time; Theorem 3.1 provides the most simplest kernel expression that characterises the information in the gates. From Theorem 3.1 it is evident that the role of width is averaging (due to the division by \( w \)). Each layer therefore corresponds to a base kernel \( \frac{\langle G_l(x), G_l(x') \rangle}{w} \) which measures the correlation of the gates. The role of depth is to provide the product of kernels. To elaborate, the feature network provides the gates \( G_l(x) \), and the value network realises the product kernel in Theorem 3.1 by laying out the GalUs depth-wise, and connecting them to form a deep network. The depth-wise layout is important: for instance, if we were to concatenate the gating features as \( \varphi(x) = \langle G_l(x), l = 1, \ldots, d - 1 \rangle \in \{0, 1\}^{(d - 1)w} \), it would have only resulted in the kernel \( \langle \varphi(x), \varphi(x') \rangle = \sum_{l=1}^{d-1} \langle G_l(x), G_l(x') \rangle \), i.e., a sum (not product) of kernels.

- **Constant 1 Input.** This has a minor impact, in that, the expression on right hand side of Theorem 3.1 becomes \( d \cdot \sigma^{2d-1} \cdot d_{\text{in}} \cdot \prod_{l=1}^{d-1} \frac{\langle G_l(x), G_l(x') \rangle}{w} \), i.e., the kernel still has information of the gates.

- **Destroying structure by permuting the layers.** \( \prod_{l=1}^{d-1} \frac{\langle G_l(x), G_l(x') \rangle}{w} \) is permutation invariant, and hence permuting the layers has no effect.

3.1.2 Convolution Global Average Pooling: Rotationally Invariant Kernel

We consider networks with circular convolution and global average pooling (architecture and notations are in the Appendix). In Theorem 3.2 let the circular rotation of vector \( x \in \mathbb{R}^d_w \) by \( r \) co-ordinates be defined as \( \text{rot}(x, r)(i) = x(i + r) \) if \( i + r \leq d_{\text{in}} \) and \( \text{rot}(x, r)(i) = x(i + r - d_{\text{in}}) \) if \( i + r > d_{\text{in}} \).

**Theorem 3.2.** Under Assumption 2.7 for a suitable \( \beta_{cv} \) (see Appendix for expansion of \( \beta_{cv} \)):

\[
NTK_{\text{CONV}}(x, x') \to \beta_{cv} \cdot \frac{\sum_{r=0}^{d_{\text{in}}-1} \langle x, \text{rot}(x', r) \rangle \cdot \text{overlap}(x, x, \text{rot}(x', r))}{d_{\text{in}}^2}, \text{ as } w \to \infty.
\]

\[
\sum_{r=0}^{d_{\text{in}}-1} \langle x, \text{rot}(x', r) \rangle \cdot \text{overlap}(x, x, \text{rot}(x', r)) = \sum_{r=0}^{d_{\text{in}}-1} \sum_{i=1}^{d_{\text{in}}} x(i) \cdot \text{rot}(x', r)(i) \cdot \text{overlap}(i, x, \text{rot}(x', r)), \text{ where the inner \`\Sigma\' is the inner product between } x \text{ and } \text{rot}(x', r) \text{ weighted by } \text{overlap} \text{ and the outer \`\Sigma\' covers all possible rotations,}
\]
which in addition to the fact that all the variables internal to the network rotate as the input rotates, results in the rotational invariance. It was observed by [Arora et al. 2019] that networks with global-average-pooling are better than vanilla convolutional networks. The rotational invariance holds for convolutional architectures only in the presence of global-pooling. So, this result explains why global-average-pooling helps. That said, rotational invariance is not a new observation; it was shown by [Li et al. 2019] that prediction using CNTK-GAP is equivalent to prediction using CNTK without GAP but with full translation data augmentation (same as rotational invariance) with wrap-around at the boundary (same as circular convolution). However, Theorem 3.2 is a necessary result, in that, it shows rotational invariance is recovered in the dual view as well.

- The expression in Theorem 3.2 becomes \( \frac{\partial}{\partial x} \sum_{r=0}^{d_{\text{in}}-1} \sum_{i=1}^{d_{\text{in}}} \text{overlap}(i, x, \text{rot}(x', r)) \) for a constant 1 input. The key novel insight is that the rotational invariance is not lost and overlap matrix measures the correlation of the paths which in turn depends on the correlation of the gates.

- Destroying structure by permuting the layers does not destroy the rotational invariance in Theorem 3.2. This is because, due to circular convolutions all the internal variables of the network rotate as the input rotates. Permuting the layers only affects the ordering of the layers, and does not affect the fact that the gates rotate if the input rotates, and cor relates in the gates is not lost.

3.1.3 Residual Networks With Skip Connections (ResNet): Ensemble Of Kernels

We consider a ResNet with \( \{(b+2)\} \) blocks and \( b \) skip connections between the blocks. Each block is a fully connected (FC) network of depth \( d_{\text{blk}} \) and width \( w \). There are \( 2^b \) many sub-FCNs within this ResNet (see Definition 3.1). Note that the blocks being fully connected is for expository purposes, and the result continue to hold for any kind of block.

Definition 3.1. [Sub FCNs] Let \( 2^b \) denote the power set of \( [b] \) and let \( \mathcal{J} \in 2^b \) denote any subset of \( [b] \). Define the \( \mathcal{J}^{th} \)-sub-FCN of the ResNet to be the fully connected network obtained by (i) including block \( j, \forall j \in \mathcal{J} \) and (ii) ignoring block \( j, \forall j \notin \mathcal{J} \).

Theorem 3.3. Let \( \text{NPK}^{FC} \) be the NPK of the \( \mathcal{J}^{th} \)-sub-FCN, and \( \beta_{\mathcal{J}}^{FC} \) (see Appendix for expansion of \( \beta_{\mathcal{J}}^{FC} \)) be the associated constant. Under Assumption 2.7 we have:

\[
\text{NTK}^{\text{RES}} \to \sum_{\mathcal{J} \in 2^b} \beta_{\mathcal{J}}^{FC} \text{NPK}^{FC}_{\mathcal{J}}, \text{ as } w \to \infty
\]

- Ensemble. To the best of our knowledge, this is the first theoretical result to show that ResNets have an ensemble structure, where each kernel in the ensemble, i.e., \( \text{NPK}^{FC}_{\mathcal{J}} \) corresponds to one of the \( 2^b \) sub-architectures (see Definition 3.1). The ensemble behaviour of ResNet and presence of \( 2^b \) architectures was observed by [Veit et al. 2016], however without any concrete theoretical formalism.

- Effect of constant 1 input is as before for kernels \( \text{NPK}^{FC} \) and translates to the ensemble \( \text{NTK}^{\text{RES}} \).

- Destroying structure. The ResNet inherits the invariances of the block level kernel. In addition, the ensemble structure allows to even remove layers. [Veit et al. 2016] showed empirically that removing single layers from ResNets at test time does not noticeably affect their performance, and yet removing a layer from architecture such as VGG leads to a dramatic loss in performance. Theorem 3.3 can be seen to provide a theoretical justification for this empirical result. In other words, due to the ensemble structure a ResNet is capable of dealing with failure of components. While failure of component itself does not occur unless one makes them fail purposefully as done in [Veit et al., 2016], the insight is that even if one or many of the kernels in the ensemble are corrupt and the good ones can compensate.

3.2 Numerical Experiments

Section 3.1 presented theoretical results which demystified the layer-by-layer view in value network, in this section we will verify these theoretical results in experiments. We then show that DLGN recovers major part of performance of state-of-the-art DNNs on CIFAR-10 and CIFAR-100.

Setup Details. We consider 3 DNN architectures, C4GAP, VGG-16 and Resnet-110, and their DGN and DLGN counterparts. Here C4GAP is a simple model (achieves about 80\% accuracy on CIFAR-10), mainly used to verify the theoretical insights in Section 3.1. VGG-16 and Resnet-110 are chosen for their state-of-the-art performance on CIFAR-10 and CIFAR-100. All models are trained using off-the-shelf optimisers (for more details, see Appendix C). The DGN and DLGN are trained from scratch, i.e., both the feature and value network are initialised at random and trained. In DGN and DLGN, we use soft gating (see Figure 3) so that gradient flows through the feature network and the gates are learnt (we chose \( \beta = 10 \)). In what follows, we use the notation \( \text{DGN}(x^f, x^v) \) and \( \text{DLGN}(x^f, x^v) \) where \( x^f \)
Disentangling Value Network. We show that destroying the layer-by-layer structure via permutations and providing a constant 1 input do not degrade performance. Since our aim here is not state-of-the-art performance, we use C4GAP with 4 convolutional layers which achieves only about 80% test accuracy on CIFAR-10, however, enables us to run all the $4! = 24$ layer permutations. The C4GAP, DGN and DLGN with layer permutations are shown in Figure 3. Once a permutation is chosen, it is fixed during both training and testing. The results in Table I of Figure 3 show that there is no significant difference in performance between DGN (a permutation is chosen, it is fixed during both training and testing). The results in Table I of Figure 3 show that there is no significant difference in performance between DGN vs DGN, and DLGN vs DLGN, i.e., constant 1 input does not hurt. Also, there is no significant difference between the models without permutations and the models with permutations. These counter intuitive and surprising results are difficult to explain using the commonly held ‘sophisticated features are learnt layer-by-layer’ view. However, neither the permutations or the constant 1 input destroys the correlation in the gates, and are not expected to degrade performance as per the insights in Section 3.1.

DLGN Performance. For this we choose VGG-16 and Resnet-110. The results in Table II of Figure 3 show that the DLGN recovers most part of the performance. For each we choose VGG-16 and Resnet-110. The results in Table II of Figure 3 show that the DLGN recovers most part of the performance. For each dataset, the top row has results for vanilla models without permutations (the results are averaged over 5 runs) and the bottom row has results of $4! - 1 = 23$ permutations (except the identity) for each model (the results are averaged over the 23 permutations).

3.3 Is DLGN a Universal Spectral Approximator

Motivated by the success of DLGN, we further break it down into DLGN-Shallow Features (DLGN-SF), wherein, the feature network is a collection of shallow single matrix multiplications. We compare a shallow DNN with ReLU
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Table III

| Dataset            | C1GAP (epochs) | C1GAP-16-ENS (epochs) | C1GAP (epochs) | C1GAP-16-ENS (epochs) | C4GAP-DLGN-SF (epochs) | VGG-16-DLGN-SF (epochs) |
|--------------------|----------------|-----------------------|----------------|-----------------------|------------------------|-------------------------|
| CIFAR-10           | 62.2±0.2       | 63.81                 | 70.0±0.2       | 72.2                  | 75.1±0.4               | 84.9±0.2                |
| CIFAR-100          | 36.3±0.2       | 38.23                 | 42.2±0.3       | 45.47                 | 47.3±0.6               | 56.3±0.2                |

Figure 4: C1GAP and C4GAP have width = 512 to make them comparable to VGG-16 whose maximum width is 512. The ensemble size is 16 to match the 16 layers of VGG-16. Note that C4GAP in Figure 3 has width=128. The last two columns show only DLGN-SF(x, 1). We observed the performance of DLGN-SF(x, x) to be ∼ 2% lesser and left it out from Table III for sake brevity. All results except for ENS are averaged over 5 runs.

called C1GAP against DLGN-SF of C4GAP (see Figure 4 and VGG-16 (see Appendix). The results are in Figure 4 based on which we observe the following:

- **Power of depth in value network and lifting to dual space.** Both C1GAP and C4GAP-DLGN-SF were trained with identical batch size, optimiser and learning rate (chosen to be the best for C1GAP). The performance of C1GAP at 200 epochs is ∼ 10% lower than that of C4GAP-DLGN-SF. After 2000 epochs of training and ensembling 16 such C1GAP’s as C1GAP-16-ENS closes the gap within ∼ 3% on C4GAP-DLGN-SF. Yet, a deeper architecture VGG-16-DLGN-SF is ∼ 10% better than C1GAP-16-ENS. Note that both VGG-16-DLGN-SF and C1GAP-16-ENS have gates for 16 layers produced in a shallow manner. While in a C1GAP-16-ENS, ‘16’ C1GAPs are ensembled, in VGG-16-DLGN-SF these gates for 16 layers are used as gating signals to turn ‘on/off’ the GalUs laid depth-wise as 16 layers of the value network, which helps to lift the computations to the dual space. Thus, using the gates to lift (instead of ensembling) the computations to the dual space in the value network is playing a critical role, investigating which is an important future work.

- **Power of depth in feature network.** By comparing CIFAR-100 performance of VGG-16-DLGN-SF in Figure 4 and that of VGG-16-DLGN in Figure 3, we see ∼ 6% improvement if we have a deep linear network instead of many shallow linear networks as the feature network. This implies depth helps even if the feature network is entirely linear, investigating which is an important future work.

- **DGN vs DLGN** In Table I and II of Figure 4, the difference between DGN and DNN is minimal (about 3%), however, the difference between DLGN and DNN is significantly large. Thus, it is important to understand the role of the ReLUs in the feature network of DGN. It is interesting to know whether this is simpler than understanding the DNN with ReLUs itself.

- **Is DLGN a Universal Spectral Approximator?** The value network realises the NPK which in general is an ensemble (assuming skip connections). The NPK is based on the gates whose pre-activations are generated linearly. It is interesting to ask whether the DLGN via its feature network learns the right linear transformations to extract the relevant spectral features (to trigger the gates) and via its value network learns the ensembling of kernels (based on gates) in a dataset dependent manner.

## 4 Conclusion

Entanglement of the non-linear and the linear operation in each layer of a DNN makes them uninterpretable. This paper proposed a novel DLGN which disentangled the computations in a DNN with ReLUs into two mathematically interpretable linearities, the ‘primal’ linearity from the input to the pre-activations that trigger the gates, and the ‘dual’ linearity in the path space. DLGN recovers more than 83.5% of performance of state-of-the-art DNNs on CIFAR-10 and CIFAR-100. Based on this success of DLGN, the paper concluded by asking ‘Is DLGN a universal spectral approximator?’.

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A Convolution With Global Average Pooling

In this section, we define NPFs and NPV in the presence of convolution with pooling. This requires three key steps (i) treating pooling layers like gates/masks (see Definition A.2) (ii) bundling together the paths that share the same path value (due to weight sharing in convolutions, see Definition A.3), and (iii) re-defining the NPF and NPV for bundles (see Definition A.4). Weight sharing due to convolutions and pooling makes the NPK rotationally invariant. We begin by describing the architecture.
Architecture: We consider (for sake of brevity) a 1-dimensional convolutional neural network with circular convolutions, with $d_{cv}$ convolutional layers ($l = 1, \ldots, d_{cv}$), followed by a global-average-pooling layer ($l = d_{cv} + 1$) and $d_{fc}$ ($l = d_{cv} + 2, \ldots, d_{cv} + d_{fc} + 1$) fully connected layers. The convolutional window size is $w_{cv} < d_{in}$, the number of filters per convolutional layer as well as the width of the FC is $w$.

Indexing: Here $i_{in}/i_{out}$ are the indices (taking values in $[w]$) of the input/output filters. $i_{cv}$ denotes the indices of the convolutional window taking values in $[w_{cv}]$. $i_{out}$ denotes the indices (taking values in $[d_{in}]$, the dimension of input features) of individual nodes in a given output filter. The weights of layers $l$ in $[d_{cv}]$ are denoted by $\Theta(i_{in}, i_{out}, l)$. For layers $l \in [d_{cv}] + d_{cv}$ are denoted by $\Theta(i_{in}, i_{out}, l)$. The pre-activations, gating and hidden unit outputs are denoted by $q_x, \Theta(i_{out}, i_{out}, l), G_x, \Theta(i_{out}, i_{out}, l)$, and $z_x, \Theta(i_{out}, i_{out}, l)$ for layers $l = 1, \ldots, d_{cv}$.

Definition A.1 (Circular Convolution). For $x \in \mathbb{R}^{d_{in}}$, $i \in [d_{in}]$ and $r \in \{0, \ldots, d_{in} - 1\}$, define:

1. $i + r = i + r, \text{ for } i + r \leq d_{in}$ and $i + r = i + r - d_{in}, \text{ for } i + r > d_{in}$.
2. $\Theta(x, r) = x(i + r), i \in [d_{in}]$.
3. $q_x, \Theta(i_{out}, i_{out}, l) = \sum_{i_{in} \in i_{out}, l} \Theta(i_{cv}, i_{in}, i_{out}, l) \cdot z_x, \Theta(i_{in}, i_{out}, l)$.

Definition A.2 (Pooling). Let $G^\text{pool}_x, \Theta(i_{out}, i_{out}, l) + 1$ denote the pooling mask, then we have $z_x, \Theta(i_{out}, i_{out}, l) + 1 = \sum_{i_{out}, l} x, \Theta(i_{in}, i_{out}, l) \cdot G^\text{pool}_x, \Theta(i_{out}, i_{out}, l) + 1$, where in the case of global-average-pooling $G^\text{pool}_x, \Theta(i_{out}, i_{out}, l) + 1 = \frac{1}{d_{in}}, \forall i_{out} \in [w], i_{out} \in [d_{in}]$.

Input Layer: $z_x, \Theta(\cdot, 1, 0) = x$

Convolutional Layers, $l \in [d_{cv}]$

- Pre-Activation: $q_x, \Theta(i_{out}, i_{out}, l) = \sum_{i_{cv}, i_{in}} \Theta(i_{cv}, i_{in}, i_{out}, l) \cdot z_x, \Theta(i_{in}, i_{out}, l)$
- Gating Values: $G_{x, \Theta}(i_{out}, i_{out}, l) = 1_{\{q_x, \Theta(i_{out}, l) > 0\}}$
- Hidden Unit Output: $z_x, \Theta(i_{out}, i_{out}, l) = q_x, \Theta(i_{out}, i_{out}, l) \cdot G_{x, \Theta}(i_{out}, i_{out}, l)$

GAP Layer, $l = d_{cv} + 1$

- Hidden Unit Output: $z_x, \Theta(i_{out}, d_{cv} + 1) = \sum_{i_{in}} z_x, \Theta(i_{in}, i_{out}, d_{cv}) \cdot G^\text{pool}_{x, \Theta}(i_{out}, i_{out}, d_{cv} + 1)$

- Fully Connected Layers, $l \in [d_{fc}] + (d_{cv} + 1)$

- Pre-Activation: $q_x, \Theta(i_{out}, l) = \sum_{i_{in}} \Theta(i_{in}, i_{out}, l) \cdot z_x, \Theta(i_{in}, l - 1)$
- Gating Values: $G_{x, \Theta}(i_{out}, l) = 1_{\{q_x, \Theta(i_{out}, l) > 0\}}$
- Hidden Unit Output: $z_x, \Theta(i_{out}, l) = q_x, \Theta(i_{out}, l) \cdot G_{x, \Theta}(i_{out}, l)$
- Final Output: $y_{\Theta}(x) = \sum_{i_{in}} \Theta(i_{in}, d_{fc} + 1) \cdot z_x, \Theta(i_{in}, d_{fc} - 1)$

Table 1: Shows the information flow in the convolutional architecture described at the beginning of Appendix A.

A.1 Neural Path Features, Neural Path Value

Proposition A.1. The total number of paths in a CNN is given by $P^{\text{conn}} = d_{in}(w_{cv}w)^{d_{cv}-w_{cv}+1}$.

Notation A.1 (Index Maps). The ranges of index maps $I_i, I_i^{\text{cv}}, I_i$ are $[d_{in}], [w_{cv}]$ and $[w]$ respectively.

Definition A.3 (Bundle Paths of Shifting Weights). Let $P^{\text{conn}} = \frac{P^{\text{conn}}}{d_{cv}}$, and $\{B_1, \ldots, B_{d_{con}}\}$ be a collection of sets such that $\forall i, j \in [P^{\text{conn}}], i \neq j$ we have $B_i \cap B_j = \emptyset$ and $\bigcup_{i=1}^{d_{con}} B_i = [P^{\text{conn}}]$. Further, if paths $p, p' \in B_i$, then $I_i^{\text{cv}}(p) = I_i^{\text{cv}}(p')$, $\forall l = 1, \ldots, d_{cv}$ and $I_i(p) = I_i(p')$, $\forall l = 0, \ldots, d_{cv}$.

Proposition A.2. There are exactly $d_{in}$ paths in a bundle.

Definition A.4. Let $x \in \mathbb{R}^{d_{in}}$ be the input to the CNN. For this input, $A_{\Theta}(x, p) \triangleq \left(\prod_{l=1}^{d_{cv}+1} G_{x, \Theta}(I_i^{\text{cv}}(p), I_i(p), l)\right) \cdot \left(\prod_{l=d_{cv}+1}^{d_{cv}+d_{fc}+1} G_{x, \Theta}(I_i(p), l)\right)$

$\phi_{x, \Theta}(p) \triangleq \sum_{p \in B_i} x(I_i(p)) \cdot A_{\Theta}(x, p)$

$\psi_{\Theta}(B_p) \triangleq \left(\prod_{l=1}^{d_{cv}} \Theta(I_i^{\text{cv}}(p), I_i-1(p), I_i(p), l)\right) \cdot \left(\prod_{l=d_{cv}+1}^{d_{cv}+d_{fc}+1} \Theta(I_i-1(p), I_i(p), l)\right)$

The results follow in a direct manner to any form of circular convolutions.
As a consequence of the skip connections, within the ResNet architecture there are $2^b$ sub-FC networks (see Definition 3.1). The total number of paths $P_{\text{res}}$ in the ResNet is equal to the summation of the paths in these $2^b$ sub-FC networks (see Proposition B.1). Now, The neural path features and the neural path value are $P_{\text{res}}$ dimensional quantities, obtained as the concatenation of the NPFs and NPV of the $2^b$ sub-FC networks.

**Proposition B.1.** The total number of paths in the ResNet is $P_{\text{res}} = d_m \cdot \sum_{i=0}^{b-1} (b)_i 2^{(i+2)d_m - 1}$.
Lemma B.1 (Sum of Product Kernel). Let $NPK_{\Theta}^{\text{RES}}$ be the NPK of the ResNet, and $NPK_{\Theta}^{\mathcal{J}}$ be the NPK of the sub-FCNs within the ResNet obtained by ignoring those skip connections in the set $\mathcal{J}$. Then,

$$NPK_{\Theta}^{\text{RES}} = \sum_{\mathcal{J} \in 2^{|\Theta|}} NPK_{\Theta}^{\mathcal{J}}$$

Proof. Proof is complete by noting that the NPF of the ResNet is a concatenation of the NPFs of the $2^b$ distinct sub-FC-DNNs within the ResNet architecture.

We re-state Theorem 3.3

Theorem B.1. Let $\sigma = \frac{c}{\sqrt{w}}$. Under Assumption 2.1, as $w \to \infty$, for $\beta_{\text{res}}^\mathcal{J} = (|\mathcal{J}| + 2) \cdot d_{\text{blk}} \cdot \sigma^2 (|\mathcal{J}| + 2) d_{\text{in}} - 1$,

$$NTK_{\Theta}^{\text{RES}} \to \sum_{\mathcal{J} \in 2^{|\Theta|}} \beta_{\text{res}}^\mathcal{J} NPK_{\Theta}^{\mathcal{J}}$$

Proof. Follows from Theorem 5.1 in [4].

C Numerical Experiments

We now list the details related to the numerical experiments which have been left out in the main body of the paper.

- **Computational Resource.** The numerical experiments were run in Nvidia-RTX 2080 TI GPUs and Tesla V100 GPUs.
  - All the models in Table I of Figure 3 we used Adam [Kingma and Ba, 2014] with learning rate of $3 \times 10^{-4}$, and batch size of 32.
  - In Section 3.2, the codes for experiments based on VGG-16 and ResNet-110 were refactored from following repository: “https://github.com/gahaalt/resnets-in-tensorflow2”.
  - For VGG-16-DLGN in Figure 3 and DLGN-SF in Figure 4 the max-pooling were replaced by average pooling so as to ensure that the feature network is entirely linear. For the comparison to be fair, we replaced the max pooling in VGG-16 reported in Figure 3 by average pooling. Batch normalisation layers were retained in VGG-16, VGG-16-DLGN and VGG-16-DLGN-SF (all three are shown in Figure 7).
  - For VGG-16-DLGN in Figure 3 and DLGN-SF in Figure 4 the max-pooling were replaced by average pooling so as to ensure that the feature network is entirely linear. For the comparison to be fair, we replaced the max pooling in VGG-16 reported in Figure 5 Batch normalisation layers were retained in VGG-16, VGG-16-DLGN and VGG-16-DLGN-SF.
  - All the VGG-16, Resnet-100 (and their DGN/DLGN) models in Table II of Figure 3 we used SGD optimiser with momentum 0.9 and the following learning rate schedule (as suggested in “https://github.com/gahaalt/resnets-in-tensorflow2”): for iterations $[0, 400)$ learning rate was 0.01, for iterations $[400, 32000)$ the learning rate was 0.1, for iterations $[32000, 48000)$ the learning rate was 0.01, for iterations $[48000, 64000)$ the learning rate was 0.001. The batch size was 128. The models were trained till 32 epochs.
  - The VGG-16-DLGN-SF in Table III of Figure 4 uses the same optimiser, batch size and learning rate schedule as the models in Table II of Figure 3 as explained in the previous point.
  - For C1GAP and C4GAP in Table III of Figure 4 we used Adam [Kingma and Ba, 2014] with learning rate of $10^{-3}$, and batch size of 32. This learning rate is best among the set $\{10^{-1}, 10^{-2}, 10^{-3}, 3 \times 10^{-4}\}$ for C1GAP.
  - Models Used in Figure 3 and Figure 4 are shown below.
Figure 5: Shows the permutations 1 – 12 C4GAP-DLGN in Table I of Figure 3. The top left is the identity permutation and is the vanilla model.
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Figure 6: Shows the permutations 13 – 24 C4GAP-DLGN in Table 1 of Figure 3.
Figure 7: Shows VGG-16 (left), VGG-16-DLGN (middle), VGG-16-DLGN-SF (right).