FIX YOUR CLASSIFIER: THE MARGINAL VALUE OF TRAINING THE LAST WEIGHT LAYER

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ABSTRACT

Neural networks are commonly used as models for classification for a wide variety of tasks. Typically, a learned affine transformation is placed at the end of such models, yielding a per-class value used for classification. This classifier can have a vast number of parameters, which grows linearly with the number of possible classes, thus requiring increasingly more resources.

In this work we argue that this classifier can be fixed, up to a global scale constant, with little or no loss of accuracy for most tasks, allowing memory and computational benefits. Moreover, we show that by initializing the classifier with a Hadamard matrix we can speed up inference as well. We discuss the implications for current understanding of neural network models.

1 INTRODUCTION

Deep neural network have become a widely used model for machine learning, achieving state-of-the-art results on many tasks. The most common task these models are used for is to perform classification, as in the case of convolutional neural networks (CNNs) used to classify images to a semantic category. CNN models are currently considered the standard for visual tasks, allowing far better accuracy than preceding approaches (Krizhevsky et al., 2012; He et al., 2016; Szegedy et al., 2015).

Neural networks (NN) are known to be a universal approximation model (Cybenko, 1989; Hornik, 1991), even with a small number of non-linear transformations (layers). They were recently shown to even allow arbitrarily chosen data points to be memorized (Zhang et al., 2017a). Training NN models and using them for inference requires large amounts of memory and computational resources, thus, extensive amount of research has been done lately to reduce the size of networks. Han et al. (2015) used weight sharing and specification, Micikevicius et al. (2017) used mixed precision to reduce the size of the neural networks by half. Tai et al. (2015) and Jaderberg et al. (2014) used low rank approximations to speed up NNs.

Hubara et al. (2016b), Li et al. (2016) and Zhou et al. (2016), used a more aggressive approach, in which weights, activations and gradients were quantized to further reduce computation during training. Although aggressive quantization benefits from smaller model size, the extreme compression rate comes with a loss of accuracy.

Past work noted the fact that predefined (Park & Sandberg, 1991) and random (Huang et al., 2006) projections can be used together with a learned affine transformation to achieve competitive results on several tasks. In this study suggest the reversed proposal - that common NN models used can learn useful representation even without modifying the final output layer, which often holds a large number of parameters that grows linearly with number of classes.

1.1 CLASSIFIERS IN CONVOLUTIONAL NEURAL NETWORKS

Convolutional neural networks (CNNs) are commonly used to solve a variety of spatial and temporal tasks. CNNs are usually composed of a stack of convolutional parameterized layers, spatial pooling layers and fully connected layers, separated by non-linear activation functions. Earlier architectures of CNNs (LeCun et al., 1998; Krizhevsky et al., 2012) used a set of fully-connected layers at later stage of the network, presumably to allow classification based on global features of an image.

Despite the enormous number of trainable parameters these layers added to the model, they are known to have a rather marginal impact on the final performance of the network (Zeiler & Fergus, 2014) and are easily compressed and reduced after a model was trained by simple means such as matrix decomposition and sparsification (Han et al., 2015). Further more, modern architecture choices are characterized with the removal of most of the fully connected layers (Lin et al., 2013; Szegedy et al., 2015; He et al., 2016), which was found to lead to better generalization and overall accuracy, together with a huge decrease in the number of trainable parameters.

Additionally, numerous works showed that CNNs can be trained in a metric learning regime (Bromley et al., 1994; Schroff et al., 2015; Hoffer & Ailon, 2015), where no explicit classification layer was introduced and the objective regarded only distance measures between intermediate representations. All of these properties provide evidence that fully-connected layers are in fact redundant and play a small role in learning and generalization.

Despite the apparent minor role they play, fully-connected layers are still commonly used as classification layers, transforming from the dimension of network features N to the number of required class categories C. Therefore, each classification model must hold $N \cdot C$ number of trainable parameters that grows in a linear manner with the number of classes.

In this work we claim that for common use-cases of convolutional network, the parameters used for the final classification transform are completely redundant, and can be replaced with a predetermined linear transform. The use of a fixed transform can, in many cases, allow a huge decrease in model parameters, and a possible computational benefit. We suggest that these can help the deployment of those models in devices with low computation ability and smaller memory capacity. Moreover, as we keep the classifier fixed, less parameters need to be updated, reducing the communication cost for models deployed in distributed systems. The use of a fixed transform which doesn't dependence on the number classes can allow models to scale to a huge number of possible outputs, without a linear cost in the number of parameters. We also suggest that these finding can shed a light on the importance of the preceding non-linear layers to learn and generalize.

2 Using a fixed classifier

2.1 Fully-connected classifiers

We focus our attention on the final representation obtained by the network, before the classifier. We denote these representation as $x = F(z; \theta)$ where F is assumed to be a deep neural network with input z and parameters θ, e.g. - a convolutional network, trained by back-propagation and stochastic-gradient-decent (SGD).

In common NN models, this representation is followed by an additional affine transformation

$$y = W^T x + b$$

where W and b are also trained by back-propagation and SGD.

For x of N length, and C different possible outputs, W is required to be a matrix of $N \times C$. Training is done using cross-entropy loss, by feeding the network outputs through a softmax activation

$$v_i = \frac{e^{y_i}}{\sum_j^C e^{y_j}}, \; i \in \{1, \ldots, C\}$$

and reducing the expected negative log likelihood with respect to ground-truth target $t \in \{1, \ldots, C\}$, by minimizing

$$\mathcal{L}(x, t) = -\log v_t = -w_t \cdot x - b_t + \log \left(\sum_j^C e^{w_j \cdot x + b_j} \right)$$

where w_i is the ith column of W.

2.2 Choosing the projection matrix

To evaluate our conjecture regarding the importance of the final classification transformation, we replaced the trainable parameter matrix W with a fixed orthonormal projection $Q \in \mathbb{R}^{N \times C}$, such

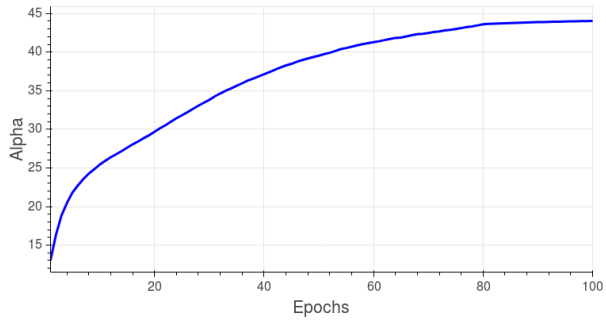


Figure 1: The softmax scale coefficient α was observed to follow a logarithmic growth over the course of training.

that $\forall i \neq j : q_i \cdot q_j = 0$ and $\|q_i\|_2 = 1$, where q_i is the ith column of Q. This can be ensured by a simple random sampling and singular-value decomposition

As the rows of classifier weight matrix are fixed with an equally valued L_2 norm, we find it beneficial to also restrict the representation of x by normalizing it to reside on the n-dimensional sphere

$$\hat{x} = \frac{x}{\|x\|_2} \tag{1}$$

This allows faster training and convergence, as the network does not need to account for changes in the scale of its weights.

We now face the problem that $w_i \cdot \hat{x}$ is bounded between -1 and 1. This causes convergence issues, as the softmax function is scale sensitive, and the network is affected by the inability to re-scale its input. This is similar to the phenomenon described by Vaswani et al. (2017) with respect to softmax function used for attention mechanisms. In the same spirit, we can amend this issue with a fixed scale T applied to softmax inputs $f(y) = \mathrm{softmax}(\frac{1}{T}y)$, also known as a softmax temperature. However, this introduces an additional hyper-parameter which may differ between networks and datasets. Instead, we suggest to introduce a single scalar parameter α to learn the softmax scale, effectively functioning as an inverse of the softmax temperature $\frac{1}{T}$.
Using normalized weights and an additional scale coefficient is similar in spirit to weight-normalization (Salimans & Kingma, 2016), with the difference that we use a single scale for all entries in the weight matrix, in contrast to a scale for each row that Salimans & Kingma (2016) uses.

We keep the additional vector of bias parameters $b \in \mathbb{R}^C$, and train using the same negative-log-likelihood criterion. More explicitly, our classifier output is now

$$v_i = \frac{e^{\alpha q_i \cdot \hat{x} + b_i}}{\sum_j^C e^{\alpha q_j \cdot \hat{x} + b_j}}, \; i \in \{1, \ldots, C\}$$

and we minimize the loss:

$$\mathcal{L}(x, t) = -\alpha q_t \cdot \frac{x}{\|x\|_2} + b_t + \log\left(\sum_{i=1}^{C} \exp\left(\alpha q_i \cdot \frac{x}{\|x\|_2} + b_i\right)\right)$$

where x is the network output for a specific sample, and $t \in \{1, \ldots, C\}$ is the ground-truth label for that sample.

Observing the behavior of the α parameter over time revealed a logarithmic growth depicted in graph 1. Interestingly, this is the same behavior exhibited by the norm of a learned classifier, first described by Hoffer et al. (2017) and linked to the generalization of the network. This was recently explained by Anonymous (2018) as convergence to a max margin classifier. We suggest that using a single parameter will enable simpler examination and possible further exploration of this phenomenon and its implications.

2.3 USING A FIXED HADMARD MATRIX

We further suggest the use of a Hadamard matrix (Hedayat et al., 1978) as the final classification transform. Hadamard matrix H is a $n \times n$ matrix, where all of its entries are either $+1$ or -1. Further more, H is orthogonal, such that $HH^T = nI_n$ where I_n is the identity matrix.
We can use a truncated Hadamard matrix $\hat{H} \in \{-1, 1\}^{C \times N}$ where all C rows are orthogonal as our final classification layer such that

$$y = \hat{H}\hat{x} + b$$

This usage allows two main benefits:

- A deterministic, low-memory and easily generated matrix that can be used to classify.
- Removal of the need to perform a full matrix-matrix multiplication - as multiplying by a Hadamard matrix can be done by simple sign manipulation and addition.

We note that n must be a multiple of 4, but it can be easily truncated to fit normally defined networks.

We also note the similarity of using a Hadamard matrix as a final classifier to methods of weight binarization such as the one suggested by Courbariaux et al. (2015). As the classifier weights are fixed to need only 1-bit precision, it is now possible to focus our attention on the features preceding it.

3 EXPERIMENTAL RESULTS

Table 1: Validation accuracy results on learned vs. fixed classifier

Network	Dataset	Learned	Fixed	# Params	% Fixed params
Resnet56 (He et al., 2016)	Cifar10	93.03%	93.14%	855,770	0.07%
DenseNet(k=12)(Huang et al., 2017)	Cifar100	77.73%	77.67%	800,032	4.2%
Resnet50 (He et al., 2016)	ImageNet	75.3%	75.3%	25,557,032	8.01%
DenseNet169(Huang et al., 2017)	ImageNet	76.2%	76%	14,149,480	11.76%
ShuffleNet(Zhang et al., 2017b)	ImageNet	65.9%	65.4%	1,826,555	52.56%

3.1 CIFAR10/100

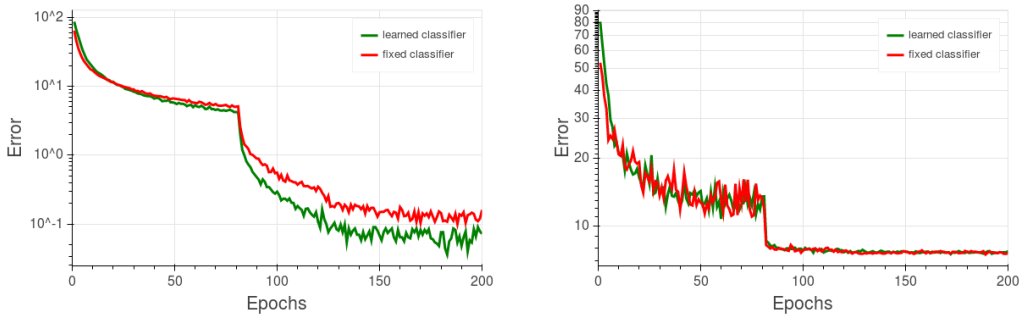
We used the well known Cifar10 and Cifar100 datasets by Krizhevsky (2009) as an initial test-bed to explore the idea of a fixed classifier. Cifar10 is an image classification benchmark dataset containing $50,000$ training images and $10,000$ test images. The images are in color and contain 32×32 pixels. There are 10 possible classes of various animals and vehicles. Cifar100 holds the same number of images of same size, but contains 100 different classes.

We trained a residual network of He et al. (2016) on the Cifar10 dataset. We used a network of depth 56 and the same hyper-parameters used in the original work. We compared two variants: the original model with a learned classifier, and our version - where a fixed transformation is used. The results shown in figure 2 demonstrate that although training error is considerably lower for the network with learned classifier, both models achieve the same classification accuracy on the validation set.

We then followed to train a model on the Cifar100 dataset. We used the DenseNet-BC model of Huang et al. (2017) with depth of 100 layers and $k = 12$. We continued to train according to the original regime and setting described for this network and dataset. Naturally, the higher number of classes caused the number of parameters to grow and encompass about 4% of the whole model. Validation accuracy for the fixed-classifier model remained equally good as the original model, and we continued to observe the same training curve.

3.2 IMAGENET

In order to validate our results on a more challenging dataset, we used the Imagenet dataset introduced by Deng et al. (2009). The Imagenet dataset spans over 1000 visual classes, and over 1.2



(a) Training error

(b) Validation error

Figure 2: Comparing training and validation error of fixed and learned classifier

million samples. CNNs used to classify Imagenet such as Krizhevsky et al. (2012), He et al. (2016), Szegedy et al. (2016) usually have a hidden representation leading to the final classifier of at least 1024 dimensions. This architectural choice, together with the large number of classes, causes the size of classifier to exceed millions of parameters and taking a sizable share from the entire model size.

We evaluated our fixed classifier method on Imagenet using Resnet50 by He et al. (2016) with the same training regime and hyper-parameters. By using a fixed classifier, approximately 2-million parameters were removed from the model, accounting for about 8% of the model parameters. Following the same procedure, we trained a Densenet169 model (Huang et al., 2017) for which a fixed classifier reduced about 12% of the parameters. Similarly to results on Cifar10 dataset, we observed the same convergence speed and approximately the same final accuracy on both the validation and training sets.

Furthermore, we were interested in evaluating more challenging models where the classifier parameters constitutes the majority amount. For this reason we chose the Shufflenet architecture (Zhang et al., 2017b), which was designed to be used in low memory and limited computing platforms. The Shufflenet network contains about 1.8 million parameters, out of which 0.96 million are part of the final classifier. Fixing the classifier resulted with a model with only 0.86 million parameters. This model was trained and found, again, to converge to similar validation accuracy as the original.

Interestingly, this method allowed Imagenet training in an under-specified regime, where there are more training samples than number of parameters. This is an unconventional regime for modern deep networks, which are usually over-specified to have many more parameters than training samples (Zhang et al., 2017a). Moreover, many recent theoretical results related to neural network training (Soudry & Hoffer, 2017; Xie et al., 2016; Safran & Shamir, 2016; Soltanolkotabi et al., 2017; Soudry & Carmon, 2016) and even generalization (Gunasekar et al., 2017; Advani & Saxe, 2017; Wilson et al., 2017) usually assume over-specification.

Table 1 summarizes our fixed-classifier results on convolutional networks, comparing to originally reported results. We offer our drop-in replacement for learned classifier that can be used to train models with fixed classifiers and replicate our results [1].

3.3 Language modeling

As language modeling requires classification of all possible tokens available in the task vocabulary, we were interested to see if a fixed classifier can be used, possible saving a very large number of trainable parameters (vocabulary size can have tens or even hundreds of thousands of different words). Recent works have already found empirically that using the same weights for both word embedding and classifier can yield equal or better results than using a separate pair of weights (Inan et al., 2016; Press & Wolf, 2017; Vaswani et al., 2017). This is compliant with our findings that the linear classifier is largely redundant. To examine further reduction in the number of parameters, we removed both classifier and embedding weights and replaced them with a fixed transform.

[1] Code is available at `https://github.com/paper-submissions/NoLinear`

We trained a language model on the WikiText2 dataset described in Merity et al. (2016), using the same setting in Merity et al. (2017). We used a recurrent model with 2-layers of LSTM (Hochreiter & Schmidhuber, 1997) and embedding + hidden size of 512. As the vocabulary of WikiText2 holds about $33K$ different words, the expected number of parameters in embedding and classifier is about 34-million. This number makes for about 89% from the $38M$ parameters used for the whole model.

We found that using a random orthogonal transform yielded poor results compared to learned embedding. We suspect that, in oppose to image classification benchmarks, the embedding layer in language models holds information of the words similarities and relations, thus requiring a fine initialization. To test our intuition, we opted to use pre-trained embeddings using word2vec algorithm by Mikolov et al. (2013) or PMI factorization as suggested by Levy & Goldberg (2014). We find that using fixed word2vec embeddings, we achieve much better results. Specifically, we use 89% less parameters than the fully learned model, and obtain only somewhat worse perplexity.

We argue that this implies a required structure in word embedding that stems from semantic relatedness between words and the natural imbalance between classes. However, we suggest that with a much more cost effective ways to train word embeddings (e.g Mikolov et al. (2013)), we can narrow the gap and avoid their cost when training bigger models.

Table 2: Validation perplexity results

Network	Dataset	Learned	Fixed	# Params	% Fixed params
2-layer LSTM (h=512)	WikiText-2	74.1	81.2	38,312,446	88.94%

4 DISCUSSION

4.1 IMPLICATIONS TO FUTURE DNN MODELS AND USE CASES

In the last couple of years a we observe a rapid growth in the number of classes benchmark datasets contain - e.g. Cifar100 (Krizhevsky, 2009), ImageNet1K, ImageNet22k (Deng et al., 2009) and language modeling (Merity et al., 2016). Therefore the computational demands of the final classifier will increase as well and should be considered no less than the architecture chosen. We use the work by Sun et al. (2017) as our use case, which introduced JFT-300M - an internal Google dataset with over 18K different classes. Using a Resnet50 (He et al., 2016), with a 2048 sized representation, this led to a model with over 36M parameters. This means that over 60% of the model parameters reside in the final classification layer.

Sun et al. (2017) further describes the difficulty in distributing this amount of parameters between the training servers, and the need to split them between 50 sub-layers. We also note the fact that the training procedure needs to account for synchronization after each parameter update - which must incur a non-trivial overhead.

Our work can help considerably in this kind of scenario - where using a fixed classifier removes the need to do any gradient synchronization for the final layer. Furthermore, using a Hadamard matrix, we can remove the need to save the transformation altogether, and make it more efficient, allowing considerable memory and computational savings.

4.2 FUTURE WORK

Understanding that linear classifiers used in NN models are largely redundant allows us to consider new approaches in training and understanding these models.

Recent works (Neyshabur et al., 2017; Bartlett et al., 2017) suggested a connection between generalization capabilities of models and various norm-related quantities of their weights. Such results might be potentially simplified in our model, since we have a single scalar variable, which seems to be the only relevant scale in the model (since we normalize the last hidden layer, and fix the last weight layer).

The use of fixed classifiers might be further simplified in Binarized Neural Networks (Hubara et al., 2016a), where the activations and weights are restricted to ± 1 during propagations. In this case

the norm of the last hidden layer is constant for all samples (equal to the square root of the layer width). This constant can be absorbed into the scale constant α, and there is no need in a per-sample normalization as in eq. 1.

We also plan to further explore more efficient ways to learn word embeddings, where similar redundancy in classifier weights may suggest simpler forms of token representations - such as low-rank or sparse versions, allowing similar benefits to the fixed transformations we suggested.

5 CONCLUSION

In this work we suggested removing the parameters from the classification layer used in deep neural networks. We showed empirical results suggesting that keeping a fixed classifier does not cause little or no decline in classification performance for common balanced datasets such as Cifar and Imagenet, while allowing a noticeable reduction in trainable parameters. We argue that fixing the last layer can reduce the computational complexity for training as well as the communication cost in distributed learning. Further more, using a Hadamard matrix as classifier might lead to some computational benefits when properly implemented, and save memory otherwise spent on large amount of transformation coefficients. As datasets tend to become more complex by time (e.g., Cifar100, ImageNet1K ImageNet22k, JFT-300M, language modeling) we believe that resource hungry affine transformation should not be used.

We also found that new efficient methods to create pre-defined word embeddings should be explored, as they require huge amount of parameters that can possibly be avoided when learning a new task. Based on these findings, we recommend future research to focus on representations learned by the non-linear part of neural networks - up to the final classifier, as it seems to be highly redundant.

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