
ANTI-DISTILLATION: KNOWLEDGE TRANSFER FROM SIMPLE MODEL TO A COMPLEX ONE

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ABSTRACT

This paper considers the problem of adapting the model to new data with a large amount of information. We propose to build a more complex model using the parameters of a simple one. It is necessary to take into account not only the accuracy of the prediction on the original samples but also the adaptability to new data and the stability of the obtained solution. The novelty of the work lies in the fact that our method allows adapting the pre-trained model to a more heterogeneous dataset. This study uses both probabilistic and algebraic methods for obtaining a student model. In the computational experiment, we analyse the quality of predictions on FashionMnist and CIFAR10 datasets.

Keywords Distillation · Knowledge Transfer · Weight Initialization · Machine Learning

1 Introduction

Training a model from scratch can lead to poor results or take a long time. To get better results faster, researchers have been developing various methods allowing to use existing trained models to solve new problems. For instance, there are knowledge distillation [2, 3], transfer learning [5], fine-tuning, low-rank model approximation [4]. Moreover, there are methods for initializing model weights for faster convergence [1]. These approaches help to decrease the time needed for training or inference and achieve higher quality.

Consider the distillation method. Statement of the initial problem is the transfer of knowledge from a cumbersome neural network or ensemble of ones to a smaller model in the classification problem. Hinton and others [2] were able to achieve this by training the student model to reproduce the probability distribution of the classes produced by the teacher model. The use of such soft targets helped to carry more information, so the student models generalization ability is comparable to the teachers. However, this research focuses on model compression under conditions of input data persistence. We want solve the inverse problem: to keep the model under conditions of increasing sample complexity.

This work proposes a method for increasing the complexity of the model based on a pre-trained one. This is done by growing the dimension of the weight space and initializing part of the student neural network with teacher model weights. Our approach allows to speed up neural network training and obtain a more robust model. As the learning is supposed to start close to the optimum point. In this way, we can adapt the pre-trained model to more variable data and reuse previously learned information.

This paper presents computational experiments on various ways of complicating the model. We consider fully connected, convolutional layers and LSTM modules. The experiment

compares uniform initialization with one based on a previously trained model and analyse differences in convergence rate, prediction variance, and achieved quality.

2 Anti-Distillation problem statement

Consider c -class classification. There are two sets

$$\begin{aligned} D_1 &= \{(\mathbf{x}_i, y_i)\}_{i=1}^{m_1}, \mathbf{x}_i \in \mathbb{R}^{n_1}, y_i \in C_1 = \{1, \dots, c_1\}, \\ D_2 &= \{(\mathbf{x}'_i, y'_i)\}_{i=1}^{m_2}, \mathbf{x}'_i \in \mathbb{R}^{n_2}, y'_i \in C_2 = \{1, \dots, c_2\}. \end{aligned}$$

Set D_2 is *more complex* than D_1 , i.e., $n_2 > n_1$ or $c_2 > c_1$.

Having the teacher model

$$g_{\text{tr}} : \mathbb{R}^{n_1} \rightarrow \Delta^{c_1}, g_{\text{tr}}(\mathbf{x}) = g(\mathbf{x}, \hat{\mathbf{u}})$$

where optimal model parameters $\hat{\mathbf{u}} \in \mathbb{R}^{N_{\text{tr}}}$ are defined as follows:

$$\hat{\mathbf{u}} = \arg \min_{\mathbf{u}} \mathcal{L}_g(\mathbf{u}, D_1) = \arg \min_{\mathbf{u}} \sum_{i=1}^{m_1} l(y_i, g(\mathbf{x}_i, \mathbf{u})),$$

here, l is a cross-entropy loss

$$l(y, \hat{y}) = - \sum_{k=1}^c [y = k] \log \hat{y}_k, y \in C, \hat{y} \in \Delta^c,$$

where Δ^c is the set of c -dimensional probability vectors.

Our approach proposes constructing student model

$$f_{\text{st}} : \mathbb{R}^{n_2} \rightarrow \Delta^{c_2}, f_{\text{st}}(\mathbf{x}) = f(\mathbf{x}, \hat{\mathbf{w}}),$$

$$\hat{\mathbf{w}} = \arg \min_{\mathbf{w}} \mathcal{L}_f(\mathbf{w}, D_2).$$

We find the solution to the above optimization problem using gradient optimization methods. The model weights $\mathbf{w} \in \mathbb{R}^{N_{\text{st}}}$ update as

$$\begin{aligned} \mathbf{w}_{t+1} &= T(\mathbf{w}_t | \mathcal{L}_f, D_2), \\ T : \mathbb{R}^{N_{\text{st}}} &\rightarrow \mathbb{R}^{N_{\text{st}}} \end{aligned}$$

is the optimization operator and $t \in \mathbb{N}$ is the gradient step number.

The function

$$\varphi : \mathbb{R}^{N_{\text{tr}}} \rightarrow \mathbb{R}^{N_{\text{st}}}$$

determines the student model parameters initialization $\mathbf{w}_1 = \varphi(\hat{\mathbf{u}})$.

3 Our setup

We construct a more complex model by extending fully connected layers and increasing number of feature maps in convolutional layers. There are different initialization methods for extended weights:

1. Zero initialization.
2. Uniform initialization.

4 Computational experiment

The goal of computational experiment is to compare the performance of models depending on the initialization of parameters.

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