OPTIMIZATION AS A MODEL FOR FEW-SHOT LEARNING

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ABSTRACT

Though deep neural networks have shown great success in the large data domain, they generally perform poorly on few-shot learning tasks, where a model has to quickly generalize after seeing very few examples from each class. The general belief is that gradient-based optimization in high capacity models requires many iterative steps over many examples to perform well. Here, we propose an LSTM-based *meta-learner* model to learn the exact optimization algorithm used to train another *learner* neural network in the few-shot regime. The parametrization of our model allows it to learn appropriate parameter updates specifically for the scenario where a set amount of updates will be made, while also learning a general initialization of the learner network that allows for quick convergence of training. We demonstrate that this meta-learning model is competitive with deep metric-learning techniques for few-shot learning.

1 Introduction

Deep learning has shown great success in a variety of tasks with large amounts of labeled data in image classification (He et al., 2015), machine translation (Wu et al., 2016), and speech modeling (Oord et al., 2016). These achievements have relied on the fact that optimization of these deep, high-capacity models requires many iterative updates across many labeled examples. This type of optimization breaks down in the small data regime where we want to learn from very few labeled examples. In this setting, rather than have one large dataset, we have a set of datasets, each with few annotated examples per class. The motivation for this task lies not only in the fact that humans, even children, can usually generalize after just one example of a given object, but also because models excelling at this task would have many useful applications. Firstly, they would help alleviate data collection as we would not require millions of labeled examples to attain reasonable performance. Furthermore, in many fields, data exhibits the characteristic of having many different classes but few examples per class. Models that are able to generalize from few examples would be able to capture this type of data effectively.

There seem to be two main reasons why gradient-based optimization fails in the face of few labeled examples. Firstly, the variants of gradient-based optimization algorithms, such as momentum (Nesterov, 1983), Adagrad (Duchi et al., 2011), Adadelta (Zeiler, 2012), and ADAM (Kingma & Ba, 2014), weren't designed specifically to perform well under the constraint of a set number of updates. Specifically when applied to non-convex optimization problems, with a reasonable choice of hyperparameters these algorithms don't have very strong guarantees of speed of convergence, beyond that they will eventually converge to a good solution after what could be many millions of iterations. Secondly, for each separate dataset considered, the network would have to start from a random initialization of its parameters, which considerably hurts its ability to converge to a good solution after a few updates. Transfer learning (Caruana, 1995; Bengio et al., 2012; Donahue et al., 2013) can be applied to alleviate this problem, by fine-tuning a pre-trained network from another task which has more labelled data; however, it has been observed that the benefit of a pre-trained network greatly decreases as the task the network was trained on diverges from the target task (Yosinski et al., 2014). What is needed is a systematic way to learn a beneficial common initialization that would serve as

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a good point to start training for the set of datasets being considered. This would provide the same benefits as transfer learning, but with the guarantee that the initialization is an optimal starting point for fine-tuning.

Previous work has suggested one manner in which to acquire quick knowledge from few examples, through the idea of *meta-learning* (Thrun, 1998; Schmidhuber et al., 1997). Meta-learning suggests framing the learning problem at two levels. The first is quick acquisition of knowledge *within* each separate task presented. This process is guided by the second, which involves slower extraction of information learned *across* all the tasks.

We present a method here that addresses the weakness of neutral networks trained with gradient-based optimization on the few-shot learning problem by framing the problem within a meta-learning setting. We propose an LSTM-based *meta-learner* optimizer that is trained to optimize a *learner* neural network. The meta-learner captures both short-term knowledge within a task and long-term knowledge common among all the tasks. By using an objective that directly captures an optimization algorithm's ability to have good generalization performance given only a set number of updates, the model is trained to converge to a good solution quickly on each task. Additionally, the formulation of our meta-learner model allows it to learn a task-common initialization for the learner, which captures fundamental knowledge shared among all the tasks.

2 TASK DESCRIPTION

We first begin by detailing the meta-learning formulation we use. In the typical machine learning setting, we are interested in a dataset D and usually split D so that we optimize parameters θ on a training set D_{train} and evaluate its generalization on the test set D_{test} . In meta-learning, however, we are dealing with meta-sets $\mathscr D$ containing multiple regular datasets, where each $D \in \mathscr D$ has a split of D_{train} and D_{test} .

We consider the k-shot, N-class classification task, where for each dataset D, the training set consists of k labelled examples for each of N classes, meaning that D_{train} consists of $k \cdot N$ examples, and D_{test} has a set number of examples for evaluation.

In meta-learning, we thus have different meta-sets for meta-training, meta-validation, and meta-testing ($\mathcal{D}_{meta-train}$, $\mathcal{D}_{meta-validation}$, and $\mathcal{D}_{meta-test}$, respectively). On $\mathcal{D}_{meta-train}$, we are interested in training a learning procedure (the meta-learning model) that can take as input one of its training sets D_{train} and produce a model that achieves high average classification performance on its corresponding test set D_{test} . Using $\mathcal{D}_{meta-validation}$ we can perform hyper-parameter selection of the meta-learning model and evaluate its generalization performance on $\mathcal{D}_{meta-test}$.

For this formulation to correspond to the few-shot learning setting, each training set in datasets $D \in \mathcal{D}$ will contain few labeled examples (we consider k=1 or k=5), that must be used to generalize to good performance on the corresponding test set.

3 Model

We now move to the description of our proposed model for meta-learning.

3.1 Model Description

Consider a single dataset $D \in \mathscr{D}_{meta-train}$. Suppose we have a *learner* neural net model with parameters θ that we want to train on D_{train} . The standard optimization algorithms used to train deep neural networks are some variant of gradient descent, which uses updates of the form

$$\theta_t = \theta_{t-1} - \alpha_t \nabla_{\theta_{t-1}} \mathcal{L}_t, \tag{1}$$

where θ_{t-1} are the parameters of the learner after t-1 updates, α_t is the learning rate at time t, \mathcal{L}_t is the loss optimized by the learner for its t^{th} update, $\nabla_{\theta_{t-1}}\mathcal{L}_t$ is the gradient of that loss with respect to parameters θ_{t-1} , and θ_t is the updated parameters of the learner.

Our key observation that we leverage here is that this update resembles the update for the cell state in an LSTM

$$c_t = f_t \odot c_{t-1} + i_t \odot \tilde{c}_t, \tag{2}$$

if $f_t = 1, c_{t-1} = \theta_{t-1}, i_t = -\alpha_t$, and $\tilde{c}_t = \nabla_{\theta_{t-1}} \mathcal{L}_t$.

Thus, we propose training a meta-learner LSTM to learn an update rule for training a neural network. We set the cell state of the LSTM to be the parameters of the learner, or $c_t = \theta_t$, and the candidate cell state $\tilde{c}_t = \nabla_{\theta_{t-1}} \mathcal{L}_t$, given how valuable information about the gradient is for optimization. We define parametric forms for i_t and f_t so that the meta-learner can determine optimal values through the course of the updates.

Let us start with i_t , which corresponds to the learning rate for the updates. We let

$$i_t = \sigma \left(\mathbf{W}_I \cdot \left[\nabla_{\theta_{t-1}} \mathcal{L}_t, \mathcal{L}_t, \theta_{t-1}, i_{t-1} \right] + \mathbf{b}_I \right),$$

meaning that the learning rate is a function of the current parameter value θ_t , the current gradient $\nabla_{\theta_t} \mathcal{L}_t$, the current loss \mathcal{L}_t , and the previous learning rate i_{t-1} . With this information, the meta-learner should be able to finely control the learning rate so as to train the learner quickly while avoiding divergence.

As for f_t , it seems possible that the optimal choice isn't the constant 1. Intuitively, what would justify shrinking the parameters of the learner and forgetting part of its previous value would be if the learner is currently in a bad local optima and needs a large change to escape. This would correspond to a situation where the loss is high but the gradient is close to zero. Thus, one proposal for the forget gate is to have it be a function of that information, as well as the previous value of the forget gate:

$$f_t = \sigma \left(\mathbf{W}_F \cdot \left[\nabla_{\theta_{t-1}} \mathcal{L}_t, \mathcal{L}_t, \theta_{t-1}, f_{t-1} \right] + \mathbf{b}_F \right).$$

Additionally, notice that we can also learn the initial value of the cell state c_0 for the LSTM, treating it as a parameter of the meta-learning. This corresponds to the initial weights of the learner model (that the meta-learner is training). Learning this initial value allows the meta-learner to determine the optimal initial weights of the learner so that training begins from a beneficial starting point, that allows optimization to proceed rapidly.

3.2 PARAMETER SHARING & PREPROCESSING

Because we want our meta-learner to produce updates for deep neural networks, which consist of tens of thousands of parameters, to prevent an explosion of meta-learner parameters we need to employ some sort of parameter sharing. Thus as in Andrychowicz et al. (2016), we share parameters across the coordinates of the learner gradient. This means each coordinate has it own hidden and cell state values but the LSTM parameters are the same across all coordinates. This allows us to use a compact LSTM model and additionally has the nice property that the same update rule is used for each coordinate, but one that is dependent on the respective history of each coordinate during optimization. We can easily implement parameter sharing by having the input be a batch of gradient coordinates and loss inputs $(\nabla_{\theta_{t,i}} \mathcal{L}_t, \mathcal{L}_t)$ for each dimension i.

Because the different coordinates of the gradients and the losses can be of very different magnitudes, we need to be careful in normalizing the values so that the meta-learner is able to use them properly during training. Thus, we also found that the preprocessing method of Andrychowicz et al. (2016) worked well when applied to both the dimensions of the gradients and the losses at each time step:

$$x \to \begin{cases} \left(\frac{\log(|x|)}{p}, \operatorname{sgn}(x)\right) & \text{if } |x| \ge e^{-p} \\ (-1, e^p x) & \text{otherwise} \end{cases}$$

This preprocessing adjusts the scaling of gradients and losses, while also separating the information about their magnitude and their sign (the later being mostly useful for gradients). We found that the suggested value of p=10 in the above formula worked well in our experiments.

3.3 Training

The question now is how do we train the LSTM meta-learner model to be effective at few-shot learning tasks? As observed in Vinyals et al. (2016), in order to perform well at this task, it is key

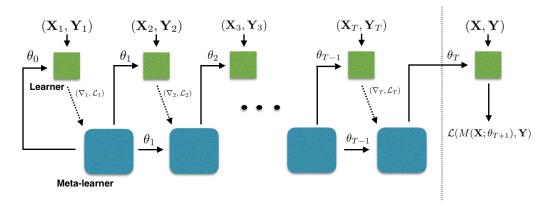


Figure 1: Computational graph for the forward pass of the meta-learner. The dashed line divides examples from the training set D_{train} and test set D_{test} . Each $(\mathbf{X}_i, \mathbf{Y}_i)$ is the i^{th} batch from the training set whereas (\mathbf{X}, \mathbf{Y}) is all the elements from the test set. The dashed arrows indicate that we do not back-propagate through that step when training the meta-learner. We refer to the learner as M, where $M(\mathbf{X}; \theta)$ is the output of learner M using parameters θ for inputs \mathbf{X} . We also use ∇_t as a shorthand for $\nabla_{\theta_{t-1}} \mathcal{L}_t$.

to have training conditions match those of test time. During evaluation of the meta-learning, for each dataset $D = (D_{train}, D_{test}) \in \mathscr{D}_{meta-test}$, a good meta-learner model will, given a series of learner gradients and losses on the training set D_{train} , suggest a series of updates for the learner model that trains it towards good performance on the test set D_{test} .

Thus to match test time, when considering each dataset $D \in \mathscr{D}_{meta-train}$, the training objective we use is the loss \mathcal{L}_{test} of the final learner model on D's test set D_{test} . While iterating over the examples in D's training set D_{train} , at each time step t the LSTM meta-learner receives $(\nabla_{\theta_{t-1}}\mathcal{L}_t,\mathcal{L}_t)$ from the learner and proposes the new set of parameters θ_t . The process repeats for T steps, after which the learner and its final parameters are evaluated on the test set to produce the loss that is then used to train the meta-learner. The training algorithm is described in Algorithm 1 and the corresponding computational graph is shown in Figure 1.

3.3.1 Gradient independence assumption

Notice that our formulation would imply that the losses \mathcal{L}_t and gradients $\nabla_{\theta_{t-1}}\mathcal{L}_t$ of the learner are dependent on the parameters of the meta-learner. Gradients on the meta-learner's parameters should normally take this dependency into account. However, as discussed by Andrychowicz et al. (2016), this complicates the computation of the meta-learner's gradients. Thus, following Andrychowicz et al. (2016), we make the simplifying assumption that these contributions to the gradients aren't important and can be ignored, which allows us to avoid taking second derivatives, a considerably expensive operation. We were still able to train the meta-learner effectively in spite of this simplifying assumption.

3.3.2 Initialization of Meta-Learner LSTM

When training LSTMs, it is advised to initialize the LSTM with small random weights and to set the forget gate bias to a large value so that the forget gate is initialized to be close to 1, thus enabling gradient flow (Zaremba, 2015). In addition to the forget gate bias setting, we found that we needed to initialize the input gate bias to be small so that the input gate value (and thus the learning rate) used by the meta-learner LSTM starts out being small. With this combined initialization, the meta-learner starts close to normal gradient descent with a small learning rate, which helps initial stability of training.

Algorithm 1 Train Meta-Learner

Input: Meta-training set $\mathscr{D}_{meta-train}$, Learner M with parameters θ , Meta-Learner R with parameters Θ .

```
1: \Theta_0 \leftarrow random initialization
 3: for d = 1, n do
 4:
           D_{train}, D_{test} \leftarrow \text{random dataset from } \mathcal{D}_{meta-train}
 5:
                                                                                                        6:
 7:
           for t = 1, T do
 8:
                 \mathbf{X}_t, \mathbf{Y}_t \leftarrow \text{random batch from } D_{train}
                 \mathcal{L}_t \leftarrow \mathcal{L}(M(\mathbf{X}_t; \theta_{t-1}), \mathbf{Y}_t)
 9:
                                                                                                ⊳ Get loss of learner on train batch
                 c_t \leftarrow R((\nabla_{\theta_{t-1}} \mathcal{L}_t, \mathcal{L}_t); \Theta_{d-1})
10:
                                                                              ▶ Get output of meta-learner using Equation 2
                 \theta_t \leftarrow c_t
                                                                                                         ▶ Update learner parameters
11:
12:
           end for
13:
           \mathbf{X}, \mathbf{Y} \leftarrow D_{test} \\ \mathcal{L}_{test} \leftarrow \mathcal{L}(M(\mathbf{X}; \underline{\theta}_T), \mathbf{Y})
14:
15:
                                                                                                 Update \Theta_d using \nabla_{\Theta_{d-1}} \mathcal{L}_{test}
16:
                                                                                                 ▶ Update meta-learner parameters
17:
18: end for
```

3.4 BATCH NORMALIZATION

Batch Normalization (Ioffe & Szegedy, 2015) is a recently proposed method to stabilize and thus speed up learning of deep neural networks by reducing internal covariate shift within the learner's hidden layers. This reduction is achieved by normalizing each layer's pre-activation, by subtracting by the mean and dividing by the standard deviation. During training, the mean and standard deviation are estimated using the current batch being trained on, whereas during evaluation a running average of both statistics calculated on the training set is used. We need to be careful with batch normalization for the learner network in the meta-learning setting, because we do not want to collect mean and standard deviation statistics during meta-testing in a way that allows information to leak between different datasets being considered. One easy way to prevent this issue is to not collect statistics at all during the meta-testing phase, but just use our running averages from metatraining. This, however, has a bad impact on performance, because we have changed meta-training and meta-testing conditions, causing the meta-learner to learn a method of optimization that relies on collection of these statistics that does not translate well to evaluation. In order to keep the two phases as similar as possible, we found that a better strategy was to collect statistics for each dataset $D \in \mathcal{D}$ for all of $\mathscr{D}_{meta-train}$, $\mathscr{D}_{meta-validation}$, and $\mathscr{D}_{meta-test}$, but then erase the running statistics when we consider the next dataset. This does not cause any information to leak between different datasets, but also allows the meta-learner to be trained on conditions that are matched between training and testing.

4 Related Work

While this work falls within the broad literature of transfer learning in general, we focus here on positioning it relative to previous work on meta-learning and few-shot learning.

4.1 META-LEARNING

Meta-learning has a long history, but has grown to prominence recently as many have advocated for it as a key to achieving human-level intelligence in the future (Lake et al., 2016). The ability to learn at two levels (learning within each task presented, while accumulating knowledge about the similarities and differences between tasks) is seen as being crucial to improving AI. Previous work has used a variety of techniques in the meta-learning setting. In Santoro et al. (2016), a memory-augmented neural network is trained to learn how to store and retrieve memories to use for

each classification task. The work of Andrychowicz et al. (2016) uses an LSTM to train a neural network model; however, they are interested in learning a general optimization algorithm to train neural networks for large-scale classification, whereas we are interested in the few-shot learning problem. This work also builds upon Hochreiter et al. (2001) and Bosc, both of which used LSTMs to train multi-layer perceptrons to learn on binary classification and time-series prediction tasks. Our work also bears similarity to Maclaurin et al. (2015), who tune the hyperparameters of momentum gradient descent by backpropagating through the chain of gradient steps to optimize the validation performance.

4.2 FEW-SHOT LEARNING

The best performing methods for few-shot learning have been mainly metric learning methods. Deep siamese networks (Koch, 2015) train a convolutional network to embed examples so that items in the same class are close while items in different classes are far away, according to some distance metric. Matching networks (Vinyals et al., 2016) refine this idea so that training and testing conditions match, by defining a differentiable nearest neighbor loss involving the cosine similarities of embeddings produced by a convolutional network.

5 EVALUATION

In this section, we describe the results of experiments, examining the properties of our model and comparing our method's performance against different approaches. Following Vinyals et al. (2016), we consider the k-shot, N-class classification setting where a meta-learner trains on many related but small training sets of k examples for each of N classes. Here, a meta-learning dataset is defined by splits of all possible classes into disjoint sets of training, validation and testing classes. To generate each instance of a k-shot, N-class task dataset $D = (D_{train}, D_{test}) \in \mathscr{D}$, we do the following: we first sample N classes from the list of classes corresponding to the meta-learning split we consider. We then sample k examples from each of those classes. These k examples together compose the training set D_{train} . Then, an additional fixed amount of the rest of the examples are sampled to yield a test set D_{test} . We generally have 15 examples per class in the test sets. When training the meta-learner, we iterate by sampling these datasets repeatedly. For meta-validation and meta-testing however, we produce a fixed number of these datasets to evaluate each method. We produce enough datasets to ensure that the confidence interval of the mean accuracy is small.

For the learner, we use a simple CNN containing 4 convolutional layers, each of which is a 3×3 convolution with 32 filters, followed by batch normalization, a ReLU non-linearity, and lastly a 2×2 max-pooling. The network then has a final linear layer followed by a softmax for the number of classes being considered. The loss function $\mathcal L$ is the average negative log-probability assigned by the learner to the correct class. For the meta-learner, we use a 2-layer LSTM, where the first layer is a normal LSTM and the second layer is our modified LSTM meta-learner. The gradients and losses are preprocessed and fed into the first layer LSTM, and the regular gradient coordinates are also used by the second layer LSTM to implement the state update rule shown in (1). At each time step, the learner's loss and gradient is computed on a batch consisting of the entire training set D_{train} , because we consider training sets with only a total of 5 or 25 examples. We train our LSTM with ADAM using a learning rate of 0.001 and with gradient clipping using a value of 0.25. We plan on releasing the code used in our evaluation experiments.

5.1 EXPERIMENT RESULTS

The Mini-ImageNet dataset was proposed by Vinyals et al. (2016) as a benchmark offering the challenges of the complexity of ImageNet images, without requiring the resources and infrastructure necessary to run on the full ImageNet dataset. Because the exact splits used in Vinyals et al. (2016) were not released, we create our own version of the Mini-Imagenet dataset by selecting a random 100 classes from ImageNet and picking 600 examples of each class. We use 64, 16, and 20 classes for training, validation and testing, respectively. We consider 1-shot and 5-shot classification for 5 classes. We use 15 examples per class for evaluation in each test set. We compare against two baselines and a recent metric-learning technique, Matching Networks (Vinyals et al., 2016), which has achieved state-of-the-art results in few-shot learning. The results are shown in Table 1.

Model	5-class	
	1-shot	5-shot
Baseline-nearest-neighbor	38.9%	48.1%
Baseline-finetune	37.3%	51.9%
Matching Network	42.4%	50.9%
Matching Network FCE	45.1%	52.2%
Meta-Learner LSTM (OURS)	44.3%	60.0%

Table 1: Average classification accuracies on Mini-ImageNet. Marked in bold are the best results for each scenario, as well as other results with an overlapping confidence interval.

The first baseline we use is a nearest-neighbor baseline (Baseline-nearest-neighbor), where we first train a network to classify between all the classes jointly in the original meta-training set. At meta-test time, for each dataset D, we embed all the items in the training set using our trained network and then use nearest-neighbor matching among the embedded training examples to classify each test example. The second baseline we use (Baseline-finetune) represents a coarser version of our meta-learner model. As in the first baseline, we start by training a network to classify jointly between all classes in the meta-training set. We then use the meta-validation set to search over SGD hyperparameters, where each training set is used to fine-tune the pre-trained network before evaluating on the test set. We use a fixed number of updates for fine tuning and search over the learning rate and learning rate decay used during the course of these updates.

For Matching Networks, we implemented our own version of both the basic and the fully-conditional embedding (FCE) versions. In the basic version, a convolutional network is trained to learn independent embeddings for examples in the training and test set. In the FCE version, a bidirectional-LSTM is used to learn an embedding for the training set such that each training example's embedding is also a function of all the other training examples. Additionally, an attention-LSTM is used so that a test example embedding is also a function of all the embeddings of the training set.

For our meta-learner, we train different models for the 1-shot and 5-shot tasks, that make 12 and 5 updates, respectively. We noticed that better performance for each task was attained if the meta-learner is explicitly trained to do the set number of updates during meta-training that will be used during meta-testing.

We attain results that are much better than the baselines discussed and competitive with Matching Networks. For 5-shot, we are able to do much better than Matching Networks, whereas for 1-shot, the confidence interval for our performance intersects the interval for Matching Network FCE. Again, we note that the numbers do not match the ones provided by Vinyals et al. (2016) simply because we created our version of the dataset and implemented our own versions of their model.

5.2 VISUALIZATION OF META-LEARNER

We also visualize the optimization strategy learned by the meta-learner, in Figure 2. We can look at the i_t and f_t gate values in Equation 2 at each update step, to try to get an understanding of how the meta-learner updates the learner during training. We visualize the gate values while training on different datasets D_{train} , to observe whether there are variations between training sets. We consider both 1-shot and 5-shot classification settings, where the meta-learner is making 10 and 5 updates, respectively. For the forget gate values for both tasks, the meta-learner seems to adopt a simple weight decay strategy that seems consistent across different layers. The input gate values are harder to interpret to glean the meta-learner's strategy. However, there seems to a be a lot of variability between different datasets, indicating that the meta-learner isn't simply learning a fixed optimization strategy. Additionally, there seem to be differences between the two tasks, suggesting that the meta-learner has adopted different methods to deal with the different conditions of each setting.

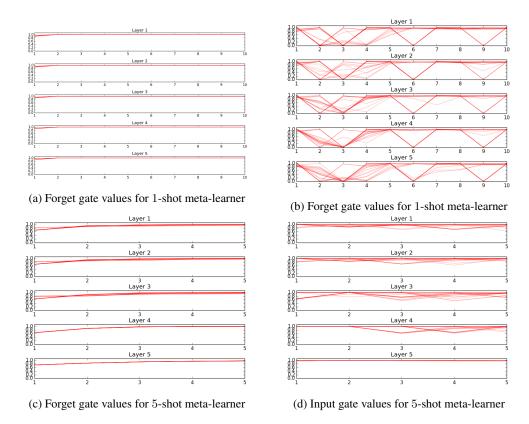


Figure 2: Visualization of the input and forget values output by the meta-learner during the course of its updates. Layers 1-4 represent the values for a randomly selected parameter from the 4 convolutional layers and layer 5 represents the values for a random parameter from fully-connected layer. The different curves represent training steps on different datasets.

6 Conclusion

We described an LSTM-based model for meta-learning, which is inspired from the parameter updates suggested by gradient descent optimization algorithms. Our LSTM meta-learner uses its state to represent the learning updates of the parameters of a learner model. It is trained to discover both a good initialization for the learner's parameters, as well as a successful mechanism for updating its parameters to a given small training set for some new classification task. Our experiments demonstrate that our approach outperforms natural baselines and is competitive to the state-of-the-art in metric learning for few-shot learning.

In this work, we focused our study to the few-shot and few-classes setting. However, it would be more valuable to train meta-learners that can perform well across a full spectrum of settings, i.e. for few or lots of training examples and for few or lots of possible classes. Our future will thus consider moving towards this more challenging scenario.

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