



# Approximate Second Order Optimization for Fast Few-Shot Learning

Master's Thesis

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# Abstract

Few-shot object detection refers to the problem of detecting objects with only a few examples. It is a challenging but real setting as many datasets arising from nature have long-tailed distributions. Currently, mainstream research in few-shot object detection employs traditional gradient descent to find a solution to optimization problems. Unfortunately gradient descent's non-competitive convergence rate results in high latency and consequently restricts real-time applications. Another issue is catastrophic forgetting where previously learned knowledge is lost as the model overfits on few-shot data.

In this thesis, we first revisited the optimization problem in few-shot object detection and designed a steep gradient descent strategy based on Conjugate Gradients (CG). In addition, we applied a meta learned weight generator and a feature-wise regularizer to make CG even faster and more stable on existing benchmark. Next, we proposed a simple yet effective hierarchical detection approach (HDA) to address catastrophic forgetting. We ran extensive experiments on the COCO dataset to measure model performance and converging speed. Experiments show that our optimization strategy CG is about 30 times faster than stochastic gradient descent when applied to the same two-stage tuning approach (TFA). Consequently, HDA achieves state-of-the-art in base class performance without any knowledge forgetting and has a competitive performance on novel classes. HDA converges 200 to 250 times faster than SGD in TFA, and as such shows potential for real-time applications.

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# Contents

1	Introduction       1         1.1       Focus of this Work       2         1.2       Thesis Organization       2
2	Related Work       4         2.1       Few-shot Learning       4         2.2       Optimization in Few-shot Learning       6         2.3       Object Detection       6         2.4       Few-shot Object Detection       6
3	Methods       9         3.1       Fast Convergence with Conjugate Gradient       9         3.1.1       Baseline Approach: TFA       9         3.1.2       Conjugate Gradient-based Optimization       10         3.2       Refined CG in TFA with Meta-Learning       12         3.3       Hierarchical Detection Approach       13
4	Experiments and Results174.1Dataset174.2Evaluation Metrics174.3Implementation Details174.4Experiments184.4.1Naive CG in TFA184.4.2Meta-Learning to Refine CG in TFA194.4.3HDA with CG194.4.4Running Time20
5	Discussion       23         5.1       Benchmarks       23         5.2       Model Performance in Average Precision       23         5.3       Model Performance in Convergence Speed       24
6	Conclusion         29           6.1         Summary         29           6.2         Outlook         29

CONTENTS

# **List of Figures**

3.1	Work flow of meta learning weight generator and regularizer	13
3.2	Workflow of hierarchical detection approach	15
5.1	Influence of meta learned modules on TFA+CG for 1-shot COCO dataset (seed 0)	24
5.2	Convergence of different models on 1-shot COCO dataset (seed 0)	26

LIST OF FIGURES

# **List of Tables**

4.1	Average precision of the model TFA+SGD	18
4.2	Average precision of the model TFA+CG	19
4.3	Average precision of the model TFA+CG+meta	19
4.4	Average precision of the model HDA+CG	20
4.5	Average precision of the model HDA+CG+aug	20
4.6	Time per iteration and number of iterations in optimization	20
4.7	Running time for feature extraction and optimization	21
	Performance of different models in generalized few-shot object detection	
5.2	Overall running time of different models	27

LIST OF TABLES

# **Chapter 1**

# Introduction

Humans have the ability to quickly learn new concepts after experiencing similar tasks. This ability is commonly demonstrated as they go about their daily lives. For example, a pianist can readily learn to play the violin while exerting less effort compared to a novice with no background in music. If this could be simply be explained by age, consider how a toddler can learn new words after being given only a few instructions. Such an ability to effectively adapt to a new task with only a few examples is an ability also desired in machine learning. In technical terms, we define this challenge as the few-shot learning problem. Few-shot learning is significant for real-world applications of machine learning because on the one hand, real-world data from many domains naturally have a long-tailed distribution with imbalances in sample sizes across different classes. On the other hand, constraints such as computational and memory resources, physical labor costs for human annotations, as well as data privacy restricts the ability to apply the traditional machine learning paradigm of training a model on an abundance data. Real world applications require making the most of limited data. The general paradigm of few-shot learning is to have a data-abundant set of base classes and a data-scarce set of novel classes and try to generalize the knowledge learned from the abundant dataset of base classes to a few-shot dataset of novel classes [13, 26, 20, 22, 37, 33, 6, 40].

Currently, in few-shot learning, the instance-level task of object detection is less well studied compared to the image-level task of classification. As one of the fundamental tasks in computer vision, object detection aims at localizing object instances in an image, which is usually represented by bounding box, and determining the class label for each instance. Few-shot object detection is challenging given that it requires both classification and localization of the detected objects but with only a few examples. Besides, it is important for few-shot object detection to retain the performance on base classes while learning novel classes because in real applications, a scene could contain instances of both base and novel classes. The requirement to achieve high performance on both base and novel classes is also known as generalized few-shot object detection [38, 23, 12]. Currently, there are two major paradigms for few-shot object detection including meta-learning and transfer learning. The meta-learning-based approaches first learns task-level knowledge from the data-abundant base set and then applies it to a few-shot task. Benchmarks [21, 27, 11] for few-shot object detection were first developed with meta-learning as its idea to first learn task-level knowledge fits the setting of few-shot learning well. Subsequent to the meta learning-based approach, the two-stage fine-tuning approach TFA [38] and its following works [23, 12] developed the transfer learning-based approach, where the model is first trained on the data-abundant base dataset and then only a few layers are fine-tuned while the other layers remain fixed. Besides, for few-shot object detection, researchers noticed the necessity of keeping the performance on base classes while learning novel classes, as in real-world object detection one scene could contain the instances of both base and novel classes [12].

### CHAPTER 1. INTRODUCTION

While existing literature have developed a few methods to tackle few-shot object detection, there is still room for improvement in the space of "fast adaptation" and the avoidance of "catastrophic forgetting". Fast adaptation: Currently, Stochastic Gradient Descent (SGD) is commonly used for model updates in fewshot object detection. On the one hand, SGD is computationally cheap and stable at every single iteration and it has been well implemented in popular machine learning libraries. On the other hand, it endures slow convergence with a vast number of iterations, which fails to meet the demand of fast adaption and therefore hinders the potential of few-shot object detection in real-world applications. Catastrophic forgetting: As aforementioned, retaining the performance on base classes is also important. However, in few-shot learning, it is often the case that the knowledge learned on old classes is largely forgotten when the model adapts to new classes, which is referred as catastrophic forgetting. Even in the transfer learning-based approach where the model performance on base classes is largely refined than its previous meta learning-based benchmarks, forgetting still happens compared to the pre-trained base model. To mitigate the forgetting, various techniques such as knowledge distillation have been applied [12]. However, even in the work [12] that the model performance on base class has been successfully retained, complexity of the model is increased and loss in knowledge distillation is still inevitable. If we want a robot to quickly detect objects in a novel scene but with a few examples, current approaches are unsatisfactory and more works need to be done to ensure a real-time performance without catastrophic forgetting.

### **1.1 Focus of this Work**

Given the issues analyzed above, the focus of the thesis can be captured with three words: **Faster**, **Higher**, **Stronger**. We propose a steep gradient descent technique and a novel hierarchical detection framework to achieve faster convergence in adaptation, higher performance on base classes without catastrophic forget-ting, and stronger performance in real-world settings.

To achieve fast adaptation, we tailored an optimization strategy based on Conjugate Gradient (CG) [8]. We started by interpreting the optimization problem in the transfer learning-based approach as a shallow learning problem, where only a few layers are optimized. We then analyzed how to apply Newton's method properly for such a problem by approximating the Hessian matrix at every step using the Gauss-Newton's method. Then we applied CG to efficiently approximate the second-order optimization at each step. We first show that our optimization method is significantly faster than the traditional SGD by applying CG to the popular transfer learning approach TFA [38]. During subsequent analysis, we investigated how task-level knowledge could benefit transfer learning approaches with CG, wherein we determined that a good initial guess of weights can speed up CG while knowledge of the sensitivity of weights to different classes can regularize the CG to minimize overfitting to few-shot data. Therefore, we meta-learned a weight generator with feature embedding and a feature-wise regularizer to boost the performance of the current transfer learning approach (HDA) that completely retains base class performance, thus negating catastrophic forgetting. The faster convergence and higher performance of HDA together make it stronger in real world challenges such as online learning of robots.

To summarize, the contributions of the thesis are as follows:

- We designed a CG-based steepest gradient descent strategy for few-shot object detection, which achieves mush faster convergence than traditional gradient descent and thus ensures a seamless performance in real-time applications.
- We integrated the meta-learning strategy with the transfer learning approach and utilized task-level

knowledge to make CG even faster and minimize overfitting when applied to current transfer learning approach.

• We proposed a simple yet effective hierarchical detection approach to completely alleviate the catastrophic forgetting on base classes without any increase of model complexity.

## **1.2** Thesis Organization

The remaining chapters of the thesis are organized as follows:

Chapter 2 will introduce the recent literature that are related to our work, including researches in fewshot learning, object detection, and few-shot object detection.

Chapter 3 will detail the methodology used in the work, including Conjugate Gradient-based optimization strategy, meta learning of weight generator and feature-wise regularization, and hierarchical detection approach.

Chapter 4 will present the dataset, evaluation metrics, implementation details of our experiments and results.

Chapter 5 will compare the results with existing benchmarks and discuss the implications.

Chapter 6 will summarize the work done in the thesis and give an outlook for future researches.

CHAPTER 1. INTRODUCTION

# Chapter 2

# **Related Work**

## 2.1 Few-shot Learning

Few-shot learning aims at learning a model that can generalize well to new knowledge provided considerably small datasets. Currently, several paradigms have been developed for the few-shot learning issues: metalearning [13, 26, 20, 32, 14, 39, 34], metric learning [22, 37, 33, 35, 7], and transfer learning [7, 9]. The goal of meta-learning is to first learn task-level knowledge before learning a specific task. This task-level knowledge helps the model quickly adapt to new tasks with limited training data, which makes it suitable for the problem of few-shot learning. Optimization-based meta-learning is a prevalent paradigm and one way is to learn a good weight initialization. MAML [13] and its variants [26, 20] try to learn a good initial weight by optimizing it over simulated tasks in the meta-learning phase, and with the meta-learned weights, the model can generalize to the few-shot data with only a few steps of gradient update. Instead of learning the initial weights directly, another way is to meta-learn weight generation. Rusu et al. [32] meta learned a low dimensional embedding space to sample the model weights by conditioning on the few-shot data, which is easier to optimize compared to the high dimension space of model parameters. Gidaris et al. [14] adopted an attention-based weight generator for few-shot classification. Wang et al. [39] applied a meta-learned network module to generate the model weights by feature embedding. Sun et al. [34] metalearned the scaling and shifting of pre-trained weights of a deep neural network to leverage the advantages of both meta-learning and transfer learning. Compared to learning the model weights directly, generating the weights conditioned on inputs is usually more task-specific, which was also applied in the thesis work.

Another paradigm to tackle the few-shot learning issue is metric learning, which also achieved competitive results. Metric learning aims at measuring the distance of the input data in a learned space. There are different variants to measure the distance. Siamese Neural Network [22] compare the similarity of the inputs for one-shot learning. Matching Nets [37] maps a query example to its label by matching it with the labeled support set. The Prototypical Network [33] computes the distance based on the prototype representation of input classes. Relation Network [35], which is also meta-learned, computes the relation scores between the query image and the few-shot examples of novel classes. Researchers also found that distance-based classifiers, which compute the cosine similarities between the feature and weight vectors, can reduce the intra-class variation and thus benefit the joint prediction of base and novel classes [14, 7, 38].

On another note regarding transfer learning, Chen *et al.* [7] fixed the pre-trained weights of base representation and fine-tuned the classifier for few-shot novel classes, which achieved results comparable with state-of-the-art. Dhillon *et al.* propose a transfer learning-based approach for image classification and consider it as a strong baseline for few-shot learning problem.

# CHAPTER 2. RELATED WORK 2.2 Optimization in Few-shot Learning

Bertinetto *et al.* [1] noticed that in the few-shot learning setting, updating only the parameters sensitive to specific classes actually formulate a shallow learning problem, and adaptation strategies that are more efficient than gradient descent exist for such a problem. They tested ridge and sigmoid regressions with closed-form solutions to achieve fast convergence for the meta-learning-based few-shot learning. Similarly, FIML [36] proposed a steepest gradient optimization algorithm that can be generalized to various learning objectives. A series of works [8, 2] applied the conjugate gradient method to update the few-shot learner in online tracking, where only the annotation of the initial frame is available and the model needs to be able to quickly adapt the following frames. A similar approach with conjugate gradient was also applied to other online settings such as video object segmentation [3] and has been proved to be efficient. While in current researches of few-shot object detection the gradient descent method is still dominant, we applied the conjugate gradient method for fast convergence that is preferable in real-time applications.

## 2.3 Object Detection

Currently, the paradigms of object detection are dominated by the one-stage [28, 25, 29, 30] and two-stage detection approaches [16, 15, 31, 18, 17]. YOLO series [28, 29, 30] is one of the representative works in one-stage detection, which applied a single neural network to predicting both the bounding box and classification score based on the anchors densely placed on an image. For two-stage detection, R-CNN [16] and its following variants (e.g. Fast R-CNN [15], Faster-RCNN [31], Mask R-CNN [17]) are popular, which firstly samples class-agnostic proposals from image features in the first stage, and makes region-wise predictions given the sampled proposals in the second stage. Compared to the one-stage approach, the two-stage approach is usually slower but more accurate. Recently, the attention-based transformer was also introduced to vision task [10] and Carion *et al.* [5] proposed DETR with encoder-decoder architecture to detect objects directly from global context with less heuristic design, which also achieve competitive results. Currently, the two-stage approach Faster-RCNN is widely used as the base detector to build benchmarks in existing researches for few-shot object detection [38, 19, 23, 12]. In the thesis, we applied the two-stage Faster RCNN as our basic model architecture.

## 2.4 Few-shot Object Detection

Few-shot object detection aims to detect the novel objects with only a few training examples, which is less studied compared to few-shot classification but is more complicated as an instance-level task that requires both localization and classification. Existing literature mainly adopted two paradigms to tackle the issue: meta-learning-based approach [21, 27, 11, 41, 19] and transfer learning based approach [6, 40, 38, 23, 12]. Similar to the idea of meta-learning introduced before, researchers leverage the meta-learned task-level knowledge to the detection task with limited training data. MetaYOLO [21] meta learned a feature learner module to extract the generic features of novel objects and a reweighting module to make predictions provided these features. ONCE [27] studied the few-shot object detection in the incremental setting by extracting class-agnostic features and meta-learning a code generator that is class-specific. Fan *et al.* [11] proposed Attention-RPN and Multi-Relation Detector to learn a metric space to measure the similarity of object pairs for detection. Following the DETR [41] introduced before, Meta-DETR meta leaned an encoder-decoder transformer for the few-shot object detection. DCNet [19] meta learned a Dense Relation Distillation module and Context-aware Aggregation module to fully leverage the features of novel objects.

### CHAPTER 2. RELATED WORK

For the transfer learning approach, LSTD [6] is one of the early works that adapted the detector learned on data-abundant objects to the target domain of few-shot novel objects. Wang et al. [38] proposed the two-stage fine-tuning approach TFA, where in the first stage a base predictor was trained in data-abundant base objects, and in the second stage the final layers for classification and bounding box localization were tuned on a balanced few-shot dataset containing both base and novel classes. This tuning-based approach is simple yet effective, which outperforms previous benchmarks using meta-learning. Additionally, they measured the model performance on both base and novel classes, as in real case one scenario can contain objects of both base and novel classes. Following TFA, LEAST [23] attributed the layers of the base model into the class agnostic layers extracting features at the image level and the class sensitive layers extracting features at the object level. Compared to TFA, more layers considered to be class sensitive were tuned, which boosts the performance on novel classes accordingly, however, deteriorates the performance on base classes. To mitigate this knowledge forgetting, they further applied knowledge distillation and the clustered exemplars of base objects. Fan et al. [12] noticed that even though TFA has achieved competitive results, it is still inferior to the performance of the pre-trained detector on base classes. They proposed Retentive R-CNN, which inherited the tuning approach of TFA with an auxiliary consistency loss to distill the knowledge of the base detector. Retentive R-CNN achieved performance on base classes that are on par with the pretrained detector. However, even with distillation, loss of pre-learned knowledge is still inevitable. In this thesis, we proposed the hierarchical detection approach, which is simpler, more intuitive, and achieved the performance on base classes exactly as the pre-trained detector without any losses.

CHAPTER 2. RELATED WORK

# **Chapter 3**

# Methods

Our work follows the transfer-learning approach [6, 40, 38, 23, 12] for few-shot object detection, which has a problem setting as follows: given data-abundant base categories  $C_b$  and data-scarce novel categories  $C_n$ with corresponding datasets  $D_b$  and  $D_n$ , we first train a base detector  $M_b$  on  $D_b$ , which is supposed to have strong ability to detect objects in  $C_b$ . Then we freeze the modules in the base detector  $M_b$  that are considered as class-agnostic and fine-tune only final layers on  $D_n$  or a combination of  $D_n$  on a small subset of  $D_b$  to obtain the novel detector  $M_n$  that is able to detect the novel categories  $C_n$ .

In the following sections, we will first revisit the optimization problem in one of the representative transfer-learning approaches TFA [38], and introduce our faster convergence algorithm using Conjugate Gradient (CG). Next, we will describe a dedicated strategy of meta-learning to make CG even faster and more stable based on the TFA framework. Then, we will introduce a novel hierarchical detection framework to address the catastrophic forgetting issues on base classes  $C_b$  with the ability of fast adaptation, and a new setting for few-shot object detection.

### **3.1** Fast Convergence with Conjugate Gradient

### 3.1.1 Baseline Approach: TFA

The two-stage fine-tuning approach TFA [38] deployed the two-stage detector Faster-RCNN as the base model architecture. Faster-RCNN [31] consists of a convolutional neural network (CNN) module for extracting generic image features, a Regional Proposal Network (RPN) to generate proposals for potential objects and a Region of Interest (ROI) head for making detections. Specifically for the ROI head, it first extracts the box features from the sampled proposals with two fully connected linear layers, and then the extracted box features will be passed two separate linear layers: one as a classifier C to classify each proposal and another as bounding box regressor  $\mathcal{R}$  to localize each proposal, wherein the classification and the localization are made simultaneously.

TFA proposes to first train Faster-RCNN on the data-abundant  $D_b$  to have a base detector  $M_b$ , and in the second stage, it freezes all the modules of  $M_b$  except for the two final linear layers, namely the classifier C and  $\mathcal{R}$ , which will be fine-tuned on a balanced few-shot dataset containing both base categories  $C_b$  and novel categories  $C_n$ . Therefore, the loss  $\mathcal{L}$  being optimized is actually composed of two components:

$$\mathcal{L} = \mathcal{L}_{cls} + \mathcal{L}_{loc} \tag{3.1}$$

where  $\mathcal{L}_{cls}$  is a cross entropy loss for classification in  $\mathcal{C}$  and  $\mathcal{L}_{loc}$  is a smoothed  $L_1$  loss for bounding box regression in  $\mathcal{R}$ . Stochastic Gradient Descent (SGD) [4], as a widely used optimization technique in the

#### CHAPTER 3. METHODS

machine learning, was also applied to optimize  $\mathcal{L}$  in TFA and many other researches that tackle few-shot object detection. SGD updates the model weights w by simply computing the first-order gradient of  $\mathcal{L}$ :

$$w^{i+1} = w^i - \Delta w^i \tag{3.2}$$

$$\Delta w^i = \alpha \cdot \Delta \mathcal{L}(w^i) \tag{3.3}$$

where  $\Delta w^i$  is the gradient update at iteration *i*.  $\alpha$  is the step length, and  $\Delta \mathcal{L}$  is the first-order gradient of  $\mathcal{L}$ .

SGD is computationally cheap for each iteration and is stable to reach convergence in various learning settings. However, optimization with SGD can usually suffer a slow convergence, which harms the potential of few-shot object detection in an online setting. To address this issue, we propose an optimization strategy based on Conjugate Gradient to ensure the seamless performance of the few-shot detector in real-time applications, which we will introduce in the following part.

### 3.1.2 Conjugate Gradient-based Optimization

. In the second stage of TFA, only the parameters of two separate linear layers each for C and  $\mathcal{R}$  are optimized. This actually formulates a shallow learning problem. As mentioned before, besides gradient descent, other optimization strategies could be used to optimize such a non-deep model for a faster convergence [1, 36]. To make the loss  $\mathcal{L}$  fully convex and differentiable, we replace the smoothed  $L_1$  loss with a  $L_2$  loss for  $\mathcal{L}_{loc}$ . Thus each term of  $\mathcal{L}$  is now formulated as:

$$\mathcal{L}_{cls}(pred_{cls}, gt_{cls}) = -\log \frac{\exp(pred_{cls}[gt_{cls}])}{\sum_{j} \exp(pred_{cls}[j])}$$
(3.4)

$$\mathcal{L}_{loc}(pred_{loc}, gt_{cls}, gt_{loc}) = \frac{1}{n}(pred_{loc}[gt_{cls}] - gt_{loc})^2$$
(3.5)

where  $pred_{loc}$  is the classification score from C, and  $gt_{cls}$  is the ground truth label. j indexes the other predictions scores other than  $gt_{cls}$ .  $pred_{loc}$  is the prediction of bounding box regression from  $\mathcal{R}$ , and  $gt_{cls}$  is the ground truth localization. n of the number of detected proposals.

These two terms together formulate a convex, differentiable and positive-definite loss function for  $\mathcal{L}$ . One well-known steep gradient technique to optimize such a formulation is Newton's method, which leverages the second-order information to search for a steep converging direction. To find the steepest gradient  $\Delta w^i$  for every iteration *i*, Newton's method approximate the loss  $\mathcal{L}(w^{i+1})$  at next parameter estimation  $w^{i+1}$  using second-order Taylor expansion at current estimation  $w^i$ :

$$\mathcal{L}(w^{i+1}) = \mathcal{L}(w^i + \Delta w^i) \approx \mathcal{L}(w^i) + (\Delta w^i)^T \Delta \mathcal{L}(w^i) + \frac{1}{2} (\Delta w^i)^T H_{w^i} \Delta w^i$$
(3.6)

where  $H_{w^i}$  is the Hessian matrix of second-order derivatives.

A solution can be found by setting the the derivative of  $\mathcal{L}(w^i + \Delta w^i)$  to  $\Delta w^i$  as zero:

$$\frac{\partial \mathcal{L}(w^i + \Delta w^i)}{\partial \Delta w^i} = \frac{\partial}{\partial \Delta w^i} [\mathcal{L}(w^i) + (\Delta w^i)^T \Delta \mathcal{L}(w^i) + \frac{1}{2} (\Delta w^i)^T H_{w^i} \Delta w^i] = 0$$
(3.7)

$$\Delta w^{i} = -H_{w^{i}}^{-1} \cdot \Delta \mathcal{L}(w^{i}) \tag{3.8}$$

However, calculating the inverse of Hessian matrix  $H^{-1}$  is computationally expensive and can be numerically unstable. As analyzed before that we are actually optimizing a shallow learning problem where only the parameters of two separate linear layers are tuned, it is viable to approximate  $\mathcal{L}$  at every step as a

simpler positive-definite quadratic loss. Gauss-Newton algorithm can be applied to problems with such a formulation, which avoids the calculation of second-order Hessian by approximating it with the first-order Jacobian. Gauss-Newton algorithm defines a residual r and optimizes  $\mathcal{L}$  by formulating it as a non-linear least square of r:

$$r(w) = w(x) - y \tag{3.9}$$

$$\mathcal{L}(w) = \|r(w)\|^2 = \|(w(x) - y)\|^2$$
(3.10)

where r(w) is the residual vector of at weight estimate w, (x, y) is a pair of input and ground truth data, w(x) is the model prediction given the input x.

Then it approximates the residual  $r(w^{i+1})$  at next parameter estimation  $w^{i+1}$  with the first-order Taylor expansion of  $r(w^i)$  at current estimation  $w^i$ , and then approximate the loss  $\mathcal{L}(w^{i+1})$  by replacing  $r(w^{i+1})$  with its first-order expansion:

$$r(w^{i+1}) = r(w^i + \Delta w^i) = r(w^i) + J_{w^i} \Delta w^i$$
(3.11)

$$\mathcal{L}(w^{i+1}) = ||r(w^{i+1})||^2 \approx (\Delta w^i)^T J_{w^i}^T J_{w^i} \Delta w^i + 2(\Delta w^i)^T J_{w^i}^T r(w^i) + r(w^i)^T r(w^i)$$
(3.12)

where  $J_w = \frac{\partial r(w^i)}{\partial w^i}$  is the Jacobian of first-order derivative of r at w.

A solution can be found by setting the the derivative of  $\mathcal{L}(w^{i+1})$  to  $\Delta w^i$  as zero:

$$\frac{\partial \mathcal{L}(w^{i+1})}{\partial \Delta w^i} = \frac{\partial}{\partial \Delta w^i} [(\Delta w^i)^T J_{w^i}^T J_{w^i} \Delta w^i + 2(\Delta w^i)^T J_{w^i}^T r(w^i) + r(w^i)^T r(w^i)] = 0$$
(3.13)

$$\Delta w^{i} = -(J_{w^{i}}^{T}J_{w^{i}})^{-1}J_{w^{i}}r(w^{i})$$
(3.14)

If we set  $H_{w^i} \approx J_{w^i}^T J_{w^i}$ , and  $B_{w^i} = J_{w^i} r(w^i)$ , a numerical solution of  $\Delta w^i$  can then be found by solving the following system of linear equations:

$$H_{w^i}\Delta w^i = B_{w^i} \tag{3.15}$$

Provided that  $H_{w^i}$  is positive-definite, one efficient way to find an approximated solution of  $\Delta w^i$  is the iterative Conjugate Gradient (CG) algorithm, which has been introduced and successfully applied to the online tracking [8, 2] that shares a similar setting of few-shot learning. CG approximate the solution of  $\Delta w^i$  at every iteration *i* by iteratively calculating a series of conjugate gradients  $\beta_j$  and corresponding coefficients  $\Delta c_j$ :

$$\Delta w^i = \Sigma_j \beta_j \Delta c_j \tag{3.16}$$

Intuitively, it finds the optimal step length and direction for each step of Newton iteration, which results in a steep gradient for fast convergence. To summarize, we adopt Newton's method as the foundation to derive the steepest gradient for every iteration. Given the non-deep model to optimize in the transferlearning-based approach, we further approximate the loss at every Newton iteration as a quadratic loss where the Gauss-Newton algorithm can be applied to find the gradient update by solving a linear system. Then the solution of this linear system, namely the steep gradient, can be efficiently approximated by a series of CG iterations.

### CHAPTER 3. METHODS

## **3.2** Refined CG in TFA with Meta-Learning

**Task-level knowledge to learn**. While previously the transfer-learning and meta-learning based approaches are two explicit streamlines, we argue that the task-level knowledge can also benefit the transfer-learning based approach if it is learned properly. We consider two kinds of task-level knowledge that can especially benefit the TFA framework to which CG is deployed. One of task-level knowledge is a good initialization of novel weights  $w_n^0 = \left\{ w_{n,i}^0, i = 1, \cdots, N \right\}$  for detecting  $C_n$  (N is the number of weight matrix or bias vector in each linear layer of C and B), which can help CG achieve even faster convergence. For this goal, we meta learn a network module  $\mathcal{N} = \{\mathcal{N}_i, i = 1, \cdots, N\}$  to generate the initial weights by embedding the task-specific box features, where  $\mathcal{N}_i$  is the sub-network to predict specific  $w_{n,i}^0$ . Given the input features  $\overline{x}_n \in \mathbb{R}^{c_n \times d}$  (each row of  $\overline{x}_n$  is the average of box features belonging to the same class in  $C_n$ .  $c_n$  is the number of novel classes. d is the dimension of flattened box features), each sub-network  $\mathcal{N}_i$  generates the corresponding weight or bias  $w_{n,i}^0 \in \mathbb{R}^{out_{n,i} \times feat_i}$  for each linear layer of classifier C and box regressor  $\mathcal{B}$  $(out_{n,i})$  is the output dimension of  $w_{n,i}^0$ .  $feat_i$  is the dimension of input feature.  $feat_i = d$  for weight and  $feat_i = 1$  for bias). The predicted novel weights  $w_n^0$  will be concatenated with the pretrained base weights  $w_b^0$  to have the initial weights  $w^0 = \{w_b^0, w_n^0\}$  for fine-tuning.  $\mathcal{N}_i$  consists of two fully connected linear layers and a function  $\mathcal{F}$  to rearrange the dimension of the outputs and scale the predicted weights so that it is at the same magnitude as  $w_h^0$ :

$$w_{n,i}^0 = \mathcal{N}_i(\overline{x}) = \mathcal{F}(\text{Linear2}_i(\text{ReLU}(\text{Linear1}_i(\overline{x}_n)))) \tag{3.17}$$

Another is the feature-wise sensitivity of base weights  $w_b$  when making detection of different classes. As in the fine-tuning stage, model performance on the base classes  $C_b$  tends to drop when the well pretrained weights  $w_b^0$  get updated, which is referred as 'catastrophic forgetting' [42, 12]. This is especially an issue for CG, as it can easily overfit the few-shot dataset due to fast convergence, which leads to the drop of model performance on  $C_b$ . Task-level knowledge of the sensitivity of the base weights  $w_b$  for different classes can be used to regularize the update of  $w_b$  so that the decline in model performance on  $C_b$  can be mitigated during fine-tuning. For this goal, we meta learn a set of regularizing vectors  $\lambda = \{\lambda_i, i = 1, \dots, N\}$  to regularize the base weights  $w_b = \{w_{b,i}, i = 1, \dots, N\}$  feature-wise by comparing it with the pretrained base weights  $w_b^0$ , where  $w_{b,i} \in \mathbb{R}^{out_{b,i} \times feat_i}$  and  $\lambda_i \in \mathbb{R}^{feat_i}$ . This feature-wise regularization term  $\mathcal{L}_{reg}$ will added to  $\mathcal{L}$ :

$$\mathcal{L} = \mathcal{L}_{cls} + \mathcal{L}_{loc} + \mathcal{L}_{reg} \tag{3.18}$$

$$\mathcal{L}_{reg} = \Sigma_i^N \|\lambda_i \odot (w_{b,i} - w_{b,i}^0)\|^2$$
(3.19)

where  $\odot$  is element-wise multiplication between  $\lambda_i$  and each row of the difference matrix between  $(w_{b,i}$  and  $w_{b,i}^0$ .

Meta learning phase. We follow the idea of optimization-based mata learning [13] to meta learn the weight generator  $\mathcal{N}$  and feature-wise regularizing factor  $\lambda$ . The core of the meta learning phase is to simulate the few-shot object detection task and optimize the modules over sampled tasks to learn the task-level knowledge. To simulate the few-shot dataset, we randomly sample a stream of tasks from the base dataset  $D_b$ . Each task contains a support set  $D_s$ , a query set  $D_q$ , and a random split of pesudo base classes  $C_{meta,b}$  and pesudo novel classes  $C_{meta,n}$  from base classes  $C_b$ . For each sampled task, the meta learning phase jointly optimize  $\mathcal{N}$  and  $\lambda$  with an inner loop and an outer loop. The inner loop is used to optimize the model weights  $w_{meta}$  is initialized with  $w_{meta}^0 = \left\{ w_{meta,b}^0, w_{meta,n}^0 \right\}$ , where  $w_{meta,b}^0$  is a subset of pretrained weights  $w_b^0$  corresponding to  $C_{meta,b}$ , and  $w_{pseudo,n}^0$  is generated by  $\mathcal{N}$  via embedding the box features of

 $C_{meta,n}$  in  $D_s$ . Then  $w_{meta}$  get updated over the support set  $D_s$  with a few iterations, where the feature-wise regularizing factor  $\lambda$  is applied to the loss. With the updated model weights  $w_{meta}$  passed from the inner loop, the outer loop will update  $\mathcal{N}$  and  $\lambda$  with a step of gradient descent by evaluating  $w_{meta}$  on the query set  $D_q$ . Note that in this phase the parameters of  $\mathcal{N}$  and  $\lambda$  are set as learnable, and the computational process though out both loops are tacked so that the loss evaluated on  $D_q$  can be back propagated to update  $\mathcal{N}$  and  $\lambda$ .

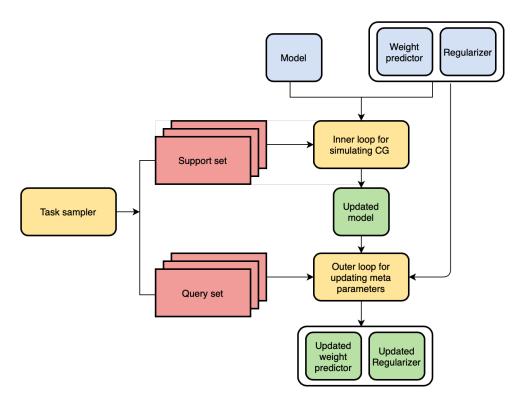


Figure 3.1: Work flow of meta learning weight generator and regularizer

Meta testing phase. In this phase, we fix the meta learned  $\mathcal{N}$  and  $\lambda$  and applied it to refine the CG optimization deployed in the TFA framework.  $\mathcal{N}$  predicts the novel weight  $w_n^0$  for  $C_n$  by embedding  $\overline{x}_n$ , which will be concatenated with  $w_b^0$  to have the initial weights  $w^0$ .  $\lambda$  will be applied to regularize the update of base weights  $w_b$  feature-wise to alleviate the knowledge forgetting on  $C_b$ .

## 3.3 Hierarchical Detection Approach

In this section we proposed a novel Hierarchical Detection Approach (HDA) for few-shot object detection to address the catastrophic forgetting on base classes  $C_b$ . It is based on transfer learning similar as TFA, but different from TFA and its variants that make joint detection on  $C_b$  and  $C_n$  by fine-tuning both  $w_b$ and  $w_n$ , HDA detect the objects according to their label hierarchy. Given the sampled box proposals, at the first hierarhical level it applied the pretrained base detector  $M_b$  to extracting the box proposals and detecting the object classes  $C = \{C_b, C_{bg,1}\}$ , where  $C_{bg,1}$  is a super-class of all proposals that are considered as background by  $M_b$ . Then at the second hierarchical level, we train a non-deep novel detector  $M_n = \{C_n, \mathcal{B}_n\}$ , which only contains a classifier  $\mathcal{C}_n$  and a box regressor  $\mathcal{B}_n$ . From the background proposals

### CHAPTER 3. METHODS

corresponding to  $C_{bg,1}$  and applied it to detecting  $C_{bg,1} = \{C_n, C_{bg,2}\}$ , where  $C_{bg,2}$  is another super class of background proposals detected by  $M_n$ . HDA was proposed based on two assumptions:

- Assumption 1. The pretrained detector  $M_b$  is good enough to detect  $C_b$ .
- Assumption 2.  $M_b$  is robust enough to distinguish  $C_{bq,1}$  from  $C_b$ .

The first assumption is based on the fact that  $M_b$  was pre-trained on a large dataset  $D_b$  with abundant annotations of classes  $C_b$ . Further fine-tuning the well-trained parameters  $w_b$  on a few-dataset usually does not improve the model performance on  $C_b$ . On the contrary, it results in 'catastrophic forgetting' of prel-learned knowledge, which requires additional tricks such as knowledge distillation to alleviate this forgetting. Therefore, to make sure the performance on  $C_b$ , one intuitive approach would be to fix  $M_b$  and apply it directly to detect  $C_b$ . However, in the few-shot object detection task, the model needs to detect not only  $C_b$ , but also  $C_n$ . This requires the second assumption regarding the robustness  $M_b$  against background proposals, which is supposed to contain most of the box proposals of  $C_n$  so that it is reliable for the further detection on  $C_n$ . In the following sections, we will detail the workflow of the hierarchical detection.

Detection on base classes. Same as TFA, HDA applied Faster-RCNN as base detector and the pretrained detector  $M_b$  was decomposed into three parts  $M_b = \{\mathcal{M}_{base}, \mathcal{C}_b, \mathcal{B}_b\}$ , where  $\mathcal{M}_{base}$  is used to extract the class-agnostic proposals and box features,  $C_b$  and  $B_b$  are two separate linear layers for classification and localization on  $C_b$ . We first applied the class-agnostic  $\mathcal{M}_{base}$  to sample a set of box proposals P = $\{P_i, i = 1, \dots, K\}$  for detection of all classes  $C = \{C_b, C_n\}$ . Each  $P_i$  here represents the bounding box, box features and associated data for training or inference. To make sure that the performance on base classes can be completely retained while not hindering the detection on novel classes, we proposed a dedicated workflow to filter the proposals  $P_b$  and  $P_{bg,1}$  for detecting  $C_b$  and  $C_n$ . Provided the box proposals P and the classification score  $S \in \mathbb{R}^{K \times (\|C_b\|_0 + 1)}$  from  $C_b$  (K is the number of box proposals,  $\|C_b\|_0 + 1$  is the number of output channels of  $C_b$  including background class), the filtering of  $P_b$  is actually to simulate how inference on S will be done to output the detection for  $C_b$ . As shown in Algorithm 1, step 1 and step 2 filter the indices  $I_b$  of proposals which have a classification score in any of the output channels for  $C_b$  greater than the score\_thresh. This makes sure that the proposals with very likely objectness are selected. However, the selected proposals with high objectness could also contain many proposals belonging to  $C_n$ . Thus step 3 to step 4 further filter the proposal indices using the Non-maximum Suppression (NMS) with the given threshold NMS\_thresh} and selects the top-k (top\_k) proposals from the NMS outputs according to the scores in descending order. Step 1 to 5 together select a subset of proposals  $P_b$  that are most likely to contain positive objects in  $C_b$  and will leave the remaining proposals  $P_{rest}$  for further detection.

**Detection on novel classes.** From the remaining proposals  $P_{rest}$  and scores  $S_{rest}$  filtered out by step 1 to step 5 in the Algorithm 1, we filter the proposals  $P_{bg,1}$  belonging to the super-class of background  $C_{bg,1}$  at the first hierarchical level, which is done by selecting proposals that have a background score greater than scores of any other class in  $C_b$  (step 6 to step 8). Then the background proposals  $P_{bg,1}$  will be passed to the novel detector  $M_n = \{C_n, \mathcal{B}_n\}$  for detecting objects in  $C = \{C_n, C_{bg,2}\}$ , where  $C_n$  is the novel class and  $C_{bg,2}$  is another super-class of background at the second hierarchical level.

**Training and inference**. For training, we fix the pretrained  $M_b$  and only need to train the novel detector  $M_n$ . The dataset is same as the one used in TFA, namely a balanced few-shot dataset  $D_{train}$  containing both base and novel classes.  $D_{train}$  will be first passed to the base detector to generate proposals P, which will be split into  $P_b$  and  $P_{bg,1}$  according to the procedure described before. Then  $P_{bg,1}$  will be used to train the classifier  $C_n$  and box regressor  $\mathcal{B}_n$  of  $M_n$ . The inference process follows a similar logic, where we passed the test set  $D_{test}$  first to  $M_b$  to extract the proposals P and then filter it into  $P_b$  and  $P_{bg,1}$ . Then we passed  $P_b$  to the  $C_b$  and  $\mathcal{B}_b$  of  $M_b$  for detecting  $C_b$ , and passed  $P_{bg,1}$  to the trained  $M_n$  for detecting  $C_n$ .

#### Algorithm 1 Procedure of Proposal Filtering **Input:** Extracted proposals $P = \{P_i, i = 1, \cdots, K\}$ , classification scores $S \in \mathbb{R}^{K \times (\|C_b\|_0 + 1)}$ **Require:** score\_thresh, nms\_thresh, top\_k **Output:** Proposals $P_b$ for detecting $C_b$ , proposals $P_{bq,1}$ for detecting $C_n$ 1: $S_b \leftarrow S[:,:-1]$ $\triangleright S_b$ is the classification score for $C_b$ 2: $I_b \leftarrow S_b > \text{score\_thresh}$ $\triangleright$ *I*<sup>*b*</sup> are indices of selected base proposals 3: $I_b \leftarrow \text{NMS}(P, I_b, \text{nms\_thresh})$ 4: $I_b \leftarrow I_b[:topk]$ 5: $P_b \leftarrow P[I_b]$ $\triangleright P_b$ is the set of filtered proposals for detecting $C_b$ 6: $P_{rest} \leftarrow P \setminus P_b, \ S_{rest} \leftarrow S \setminus S[I_b]$ 7: $I_{bg,1} \leftarrow S_{rest}[:,-1] > \operatorname{Max}(S_{rest}[:,:-1], \dim = 1)$ $\triangleright I_{bg,1}$ are indices of background proposals 8: $P_{bg,1} \leftarrow P_{rest}[I_{bg,1}]$ $\triangleright P_{bg,1}$ is the set of proposals for detecting $C_n$

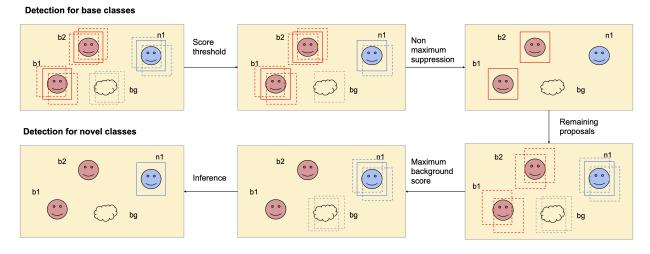


Figure 3.2: Workflow of hierarchical detection approach

15

CHAPTER 3. METHODS

# Chapter 4

# **Experiments and Results**

### 4.1 Dataset

We used the Microsoft COCO dataset [24] for our experiments, which has abundant images and annotations and is widely used for researches in few-shot object detection. COCO dataset has 80 categories that can be used for training. Following previous works [38, 23, 12], we split the 80 categories into 60 base categories and 20 novel categories. The test set has 5000 images that contain all categories, and the remaining images were used for training. A base dataset containing 60 base categories is used for training the base model and also for meta-learning. For few-shot learning of novel categories, we have a  $K \in \{1, 2, 3, 5, 10\}$  -shot datasets sampled over 10 random seeds following [38], which is a balanced small dataset containing all categories.

## 4.2 Evaluation Metrics

**Precision aspect**. Following the setting of generalized few-shot object detection [38, 23, 12], we apply the average precision to measure the model performance in all classes (AP), base classes (bAP), and novel classes (nAP), respectively. Average precision is defined as the area under the precision-recall curve at a certain IOU threshold. Same as previous benchmarks [38], here AP refers to the mean of average precisions of all classes with an IOU threshold from 0.5 to 0.95. The definition for bAP and nAP is similar to AP, but they are calculated for base and novel classes, respectively. Besides, AP at a certain IOU threshold is also used for measuring the model performance. For instance, AP50 is the AP calculated at an IOU threshold of 0.5.

**Time aspect**. To measure the speed of convergence with different optimization strategies, we measure the following metrics: time used for extracting features and proposals from the frozen layers T', time per iteration t, number of measured iterations n, and time needed for reaching convergence  $T = t \cdot n$ . The unit for time measurement is seconds (s).

### **4.3 Implementation Details**

We use Faster-RCNN as the basic model architecture and ResNet-101 as the backbone. The base detector  $M_b$  pre-trained by Wang *et al.* [38] is directly applied in the thesis, wherein  $M_b$  was trained using SGD with 110000 iterations, momentum 0.9, weight decay 0.0001, and base learning rate of 0.02 on the dataset

### CHAPTER 4. EXPERIMENTS AND RESULTS

containing 60 base categories. After the base detector being trained, all its layers except for the layers of C and B are fixed, and C and B will be tuned in TFA or adapted for hierarchical detection.

When applying CG, we first extract all proposals and associated box features from the fixed layers of the detector, and then optimize CG on all the extracted proposals and features. To have a fair comparison regarding the running speed of SGD with CG, we adapt the implementation of SGD in TFA, where the proposals and features for each image are first extracted, and then the extracted proposals and features are batched to fine-tune the model with SGD.

All experiments are run on the  $K \in \{1, 2, 3, 5, 10\}$  -shot datasets sampled over 10 random seeds, and the results of average precision are aggregated over the 10 groups of different random seeds to measure the model performance more accurately [38]. For the time aspect, we report the time metrics from experiments that were run on seed 0.

To measure the time metrics, we run all experiments on a single GPU with the same type 'TITXAN Xp'.

## 4.4 Experiments

### 4.4.1 Naive CG in TFA

**Baseline**. To have a fair comparison for the running time with CG, we adapt the original implementation of SGD in TFA w/cos to run it on the extracted proposals and features. The remaining setting for running SGD is as same as the one in TFA. Specifically, we first train the weights corresponding to detecting  $C_n$  on the novel dataset and then concatenate the weights with the pre-trained weights corresponding to  $C_b$ , and then fine-tune the model weights on the few-shot dataset. Same as the setting in TFA, we apply the SGD with momentum 0.9 and weight decay 0.0001. The learning rate for training novel weights is 0.01 and the one for fine-tuning is 0.001. The number of iterations increases with the size of the few-shot dataset. The model run in this experiment is noted as TFA+SGD for short. Table 4.1 displays the model performance of TFA+SGD.

Shots		All			Base		Novel			
Shots	AP	AP50	AP75	AP	AP50	AP75	AP	AP50	AP75	
1	24.1	40.3	25.4	31.5	52.3	33.2	2.3	4.4	2.1	
2	24.7	40.5	26.4	31.7	51.5	34.1	3.8	7.5	3.5	
3	25.0	40.6	27.0	31.6	51.0	34.5	4.9	9.4	4.7	
5	25.7	41.3	27.6	31.8	50.9	34.8	6.4	12.6	6.0	
10	26.1	42.3	27.8	31.6	51.0	34.5	8.3	16.4	7.7	

Table 4.1: Average precision of the model TFA+SGD

Naive CG. We replace the SGD with CG in the two-stage tuning approach TFA with cosine output layers (TFA w/cos) to observe how the model performance and converging speed will change. As aforementioned, for the few-shot dataset, we first extract the proposals and box features from the fixed layers of the pre-trained detector and then run CG directly on these extracted proposals and features. For weights in C and B, we initialize the parameters corresponding to  $C_n$  with randomly initialized weights. For all the datasets with different shots, we run CG with 100 Newton iterations and each Newton iteration is approximated with 2 CG iterations. The model run in this experiment is noted as TFA+CG for short. Table 4.2 displays the model performance of TFA+CG.

Shots		All			Base		Novel			
bilots	AP	AP50	AP75	AP	AP50	AP75	AP	AP50	AP75	
1	24.2	40.5	25.3	31.4	52.2	33.0	2.6	5.4	2.3	
2	25.3	41.8	27.1	32.4	53.0	34.9	4.0	8.1	3.5	
3	26.2	42.7	28.3	33.3	53.7	36.5	4.9	9.8	4.5	
5	27.3	44.1	29.7	34.3	54.6	37.6	6.4	12.5	6.0	
10	28.5	45.6	31.0	35.2	55.4	38.6	8.5	16.1	8.1	

CHAPTER 4. EXPERIMENTS AND RESULTS Table 4.2: Average precision of the model TFA+CG

### 4.4.2 Meta-Learning to Refine CG in TFA

Each element of the regularizer  $\lambda$  is randomly initialized from the Gaussian distribution  $\mathcal{G} \sim (0, 0.1)$ , and the predicted weights from  $\mathcal{N}$  are scaled with the factor 0.1 to make sure that the magnitude of the predicted weights is same as the one of the pre-trained weights. For each sampled task, we randomly split the 60 categories in  $C_b$  into 40 pseudo base categories  $C_{meta,b}$  and 20 pseudo novel categories  $C_{meta,n}$ . Both the support set and the query set are 60-way-1-shot datasets. Each meta-learning loop contains an inner loop and an outer loop. For the inner loop of meta-learning, we run 5 Newton iterations and 2 CG iterations for each Newton iteration to update the model weights. For the outer loop of meta-learning, we run a single SGD step to update  $\mathcal{N}$  and  $\lambda$  with the learning rate 0.01, momentum 0.09, and weight decay 0.0001. We run 1000 meta-learning loops to meta learn  $\mathcal{N}$  and  $\lambda$ . When applying these meta-learned modules to TFA with CG, we run 30 Newton iterations, each with 2 CG iterations. The model run in this experiment is noted as TFA+CG+meta for short. Table 4.3 displays the model performance of TFA+CG+meta.

Shots		All			Base			Novel			
bilots	AP	AP50	AP75	AP	AP50	AP75	AP	AP50	AP75		
1	24.3	40.6	25.7	31.6	52.8	33.4	2.3	4.2	2.3		
2	25.3	41.4	27.4	32.6	53.3	35.4	3.3	6.0	3.3		
3	25.9	41.9	28.3	33.3	53.5	36.4	3.8	6.9	3.9		
5	26.4	42.2	28.9	33.8	53.6	37.0	4.5	8.0	4.6		
10	26.9	42.3	29.4	34.2	53.6	37.5	4.8	8.5	5.0		

Table 4.3: Average precision of the model TFA+CG+meta

### 4.4.3 HDA with CG

Naive CG. For the proposal filtering in HDA, we set score\_thresh = 0.05, nms\_thresh = 0.5 and top\_k = 100 following the common setting of Faster R-RCNN during inference in previous benchmarks [38, 23, 12]. Given the filtered proposals, we run 30 Newton iterations and each with 2 CG iterations to train the novel detector head, which is randomly initialized. The model is noted as HDA+CG for short. Table 4.4 displays the model performance of HDA+CG.

**HDA with feature augmentation**. To boost model performance on novel classes and make CG less overfitting, we further augment the filtered proposals that have the ground truth label belonging to the novel

Shots		All			Base			Novel			
bilots	AP	AP50	AP75	AP	AP50	AP75	AP	AP50	AP75		
1	30.0	45.8	32.6	39.2	59.3	42.8	2.4	5.2	1.9		
2	30.4	46.6	32.9	39.2	59.3	42.8	4.1	8.7	3.4		
3	30.7	47.1	33.3	39.2	59.3	42.8	5.1	10.3	4.5		
5	31.0	47.7	33.5	39.2	59.3	42.8	6.4	12.8	5.8		
10	31.4	48.3	34.0	39.2	59.3	42.8	8.0	15.2	7.5		

Table 4.4: Average precision of the model HDA+CG

CHAPTER 4. EXPERIMENTS AND RESULTS

categories  $C_n$ . For each Newton iteration, we first duplicate every single shot of the proposal and corresponding box features into 5 shots, and apply Gaussian noise sampled from  $\mathcal{G} \sim (0, 0.1)$  to all the box features and then apply dropout with a rate of 0.5. The model is noted as HDA+CG+aug for short. Table 4.5 displays the model performance of HDA+CG+aug.

All Base Novel Shots AP75 AP75 AP50 AP **AP50** AP75 AP AP AP50 1 30.2 46.2 32.7 39.2 59.3 42.8 3.0 6.8 2.3 2 42.8 30.6 47.1 33.0 39.2 59.3 4.7 10.4 3.6 3 31.0 47.5 33.2 39.2 59.3 42.8 5.6 12.0 4.6 5 39.2 31.2 48.2 33.6 59.3 42.8 7.1 14.9 6.0

59.3

42.8

9.1

18.4

7.9

39.2

Table 4.5: Average precision of the model HDA+CG+aug

### 4.4.4 Running Time

10

31.8

49.4

34.1

For each model, we measure the converging speed by counting the time per iteration t and number of iterations n, based on which we calculate the time T needed for optimization. Besides, we measure the time needed for extracting proposals and box features before the optimization starts, in order to have an overall view regarding the time needed for model training. The time measurement is based on the experiments run on datasets sampled from random seed 0. As TFA+SGD has a stage for pre-training the novel weights and a stage for fine-tuning, we counted the running time for each stage, which are noted as TFA+SGD (novel) and TFA+SGD (tuning), respectively. The results can be referred in Table 4.6 and Table 4.7.

Table 4.6: Time per iteration and number of iterations in optimization

Model			$t\left(s ight)$				n					$T\left(s ight)$			
	1	2	3	5	10	1	2	3	5	10	1	2	3	5	10
TFA+SGD (novel)	0.0055	0.0059	0.0053	0.0055	0.0058	500	1500	1500	1500	2000	2.75	8.85	7.95	8.25	11.60
TFA+SGD (tuning)	0.0067	0.0067	0.0068	0.0070	0.0071	16000	32000	48000	80000	160000	107.20	214.40	326.40	560.00	1136.00
TFA+CG	0.0378	0.0687	0.1024	0.1683	0.3339	100	100	100	100	100	3.78	6.87	10.24	16.83	33.39
TFA+CG+meta	0.0464	0.0787	0.1129	0.1808	0.3401	100	100	100	100	100	4.64	7.87	11.29	18.08	34.01
HDA+CG	0.0191	0.0316	0.0458	0.0737	0.1407	30	30	30	30	30	0.57	0.95	1.37	2.21	4.22
HDA+CG+aug	0.0211	0.0359	0.0513	0.0822	0.1589	30	30	30	30	30	0.63	1.08	1.54	2.47	4.77

CHAPTER 4. EXPERIMENTS AND RESULTS

Model			$T'\left(s ight)$					$T\left(s ight)$			T' + T(s)				
model	1	2	3	5	10	1	2	3	5	10	1	2	3	5	10
TFA+SGD (novel)	3.03	5.51	6.75	10.38	19.58	2.75	8.85	7.95	8.25	11.60	5.78	14.36	14.70	18.63	31.18
TFA+SGD (tuning)	8.55	16.70	23.73	39.20	78.76	107.20	214.40	326.40	560.00	1136.00	115.75	231.10	350.13	599.20	1214.76
TFA+CG	9.36	17.43	23.20	37.53	74.42	3.78	6.87	10.24	16.83	33.39	13.14	24.30	33.44	54.36	107.81
TFA+CG+meta	8.82	18.77	24.84	37.72	76.87	4.64	7.87	11.29	18.08	34.01	13.46	26.64	36.13	55.80	110.88
HDA+CG	16.22	26.59	38.00	61.80	122.55	0.57	0.95	1.37	2.21	4.22	16.79	27.54	39.37	64.01	126.77
HDA+CG+aug	13.79	26.84	40.74	59.69	132.79	0.63	1.08	1.54	2.47	4.77	14.42	27.92	42.28	62.16	137.56

Table 4.7: Running time for feature extraction and optimization

# Chapter 5

# Discussion

## 5.1 Benchmarks

For comparison, we mainly cite the previous benchmarks [38, 23, 12] that apply the setting of generalized few-shot object detection (Table 5.1), where AP and bAP are given besides nAP to have a comprehensive view of model performance and the results are run on multiple groups of random seeds were given to reduce the variation of results on single seed. Specifically, we compare the two TFA models developed by Wang *et al.* [38]: TFA w/fc and TFA w/cos, where the former applies the fully connected layer in the classification layer C and the later applies the cosine similarity in C. Besides, they provide the benchmark that fully fine-tune the Faster-RCNN detector, noted as FRCN+ft-full. We also compare with the metrics reported in LEAST [23] and Retentive R-CNN [12]. For converging speed, we mainly compare our models with TFA+SGD, as both LEAST and Retentive R-CNN follow the tuning approach of TFA.

## 5.2 Model Performance in Average Precision

For models in TFA family, our implementation of SGD in TFA (TFA+SGD) is roughly on par with the original implementation of TFA  $w/\cos S$ . Its metrics are slightly lower than the one of TFA  $w/\cos S$  with less than 1 point, probably due to the fact that we extracted all the proposals at one time while in the original implementation, different groups of images were batched for every iteration to extract the proposals, the randomness of which could make the model less overfitting. Our naive CG implementation in TFA, TFA+CG, achieves performance on par with TFA  $w/\cos S$  with much less iterations. Especially for the fewer-shot case (e.g. 1-shot), TFA+CG outperforms the original TFA  $w/\cos S$  much (2.6 vs. 1.9 on 1-shot dataset).

For TFA+CG+meta where the meta learned modules are applied, we observe that for the 1-shot case, it has a performance on par with TFA+CG. To further verify that whether the meta learned modules are effective in the 1-shot case, we run the TFA+CG+meta for 200 iterations on the dataset sampled in seed 0 and compared it with TFA+CG initialized with random weights (Figure 5.1). Compared to the TFA+CG with randomly initialized weights, TFA+CG+meta with the generated weights from  $\mathcal{N}$  reaches higher nAP in the beginning iterations (highlighted by the green box), which suggests that the  $\mathcal{N}$  has learned a good guess of the initial weights that can speed up the convergence. With more iterations (e.g. from 100 to 200 iterations), we observe that bAP of TFA+CG without regularization has a drop of about 2 points while the bAP of TFA+CG+meta is steady (highlighted by the orange box). This indicates that the regularizer  $\lambda$  has effectively learned the sensitivity of features to different classes and regularizes the weight updates

#### CHAPTER 5. DISCUSSION

accordingly. However, in the case with more shots, the meta-learned modules harm the model performance on novel classes although the bAPs are well retained. As the weight generator and regularizer were meta learned in the 1-shot case, the results imply that these modules are not feasible for the case of higher shots, for which we might need to meta learn the modules again.

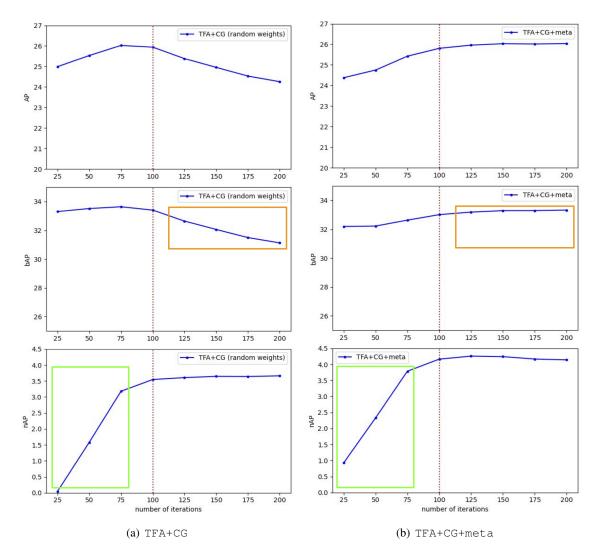


Figure 5.1: Influence of meta learned modules on TFA+CG for 1-shot COCO dataset (seed 0)

Figure 5.2 displays the evaluation of convergence of other models (TFA+SGD, HDA+CG, HDA+CG+aug) on COCO 1-shot dataset sampled from seed 0. Compared to the baseline TFA+SGD, it can be observed that the HDA models achieve significantly higher nAP with much less iterations. To further evaluate the influence the feature augmentation on HDA, we run the experiments of HDA with 40 iterations, which have 10 additional iterations than the one we use to run the benchmarks. With more iterations, we observe that there is a slight decrease on nAP in HDA+CG, while the nAP of HDA+CG+aug converges to a higher value and is more steady after convergence, indicating that feature augmentation can not only boost the performance of HDA on novel classes, but also make it less overfitting.

Our HDA model with feature augmentation, HDA+CG+aug, achieves the overall AP at the top among all

benchmarks (Table 5.1). With the base detector being fixed, it has a bAP exactly the same as the pre-trained model without any forgetting and it also achieves competitive metrics on nAP. Specifically, its performance on base classes is on par with the competitive Retentive R-CNN. Compared to Retentive R-CNN, our HDA approach, however, is much simpler without any auxiliary modules for knowledge distillation. It also has much fewer parameters to optimize, where only a detector head for novel classes needs to be trained. This also makes the HDA extremely fast in optimization, which we will discuss in the following section. For performance on novel classes, the benchmark LEAST is the best and nAP of our HDA+CG+aug is comparable with the second-place Retentive R-CNN. But in LEAST, more layers are fine-tuned, which can be expected to boost the nAP, however, results in more overfitting and have the bAP dropped accordingly. As can be observed, even with the examplar being applied to retain the performance on base classes, bAP of LEAST is just on par with TFA models and is still inferior to HDA models.

Model			AP					bAP					nAP		
Woder	1	2	3	5	10	1	2	3	5	10	1	2	3	5	10
FRCN+ft-full	16.2	15.8	15.0	14.4	13.4	21.0	20.0	18.8	17.6	16.1	1.7	3.1	3.7	4.6	5.5
TFA w/fc	24.0	24.5	24.9	25.6	26.2	31.5	31.4	31.5	31.8	32.0	1.6	3.8	5.0	6.9	9.1
TFA w/cos	24.4	24.9	25.3	25.9	26.6	31.9	31.9	32.0	32.3	32.4	1.9	3.9	5.1	7.0	9.1
TFA+SGD (Our Impl.)	24.1	24.7	25.0	25.7	26.1	31.5	31.7	31.6	31.8	31.6	2.3	3.8	4.9	6.4	8.3
LEAST	-	-	-	-	-	29.5	-	-	31.3	31.3	4.2	-	-	9.3	12.8
Retentive R-CNN	-	-	-	31.4	31.8	-	-	-	39.3	39.2	-	-	-	7.7	9.5
TFA+CG (Ours)	24.2	25.3	26.2	27.3	28.5	31.4	32.4	33.3	34.3	35.2	2.6	4.0	4.9	6.4	8.5
TFA+CG+meta (Ours)	24.3	25.3	25.9	26.4	26.9	31.6	32.6	33.3	33.8	34.2	2.3	3.3	3.8	4.5	4.8
HDA+CG (Ours)	30.0	30.4	30.7	31.0	31.4	39.2	39.2	39.2	39.2	39.2	2.4	4.1	5.1	6.4	8.0
HDA+CG+aug (Ours)	30.2	30.6	31.0	31.2	31.8	39.2	39.2	39.2	39.2	39.2	3.0	4.7	5.6	7.1	9.1

Table 5.1: Performance of different models in generalized few-shot object detection

## 5.3 Model Performance in Convergence Speed

Regarding the time needed for convergence (Table 4.6 and Table 4.7), compared to TFA+SGD, the running time for a single iteration of TFA+CG is longer than SGD, as CG iteratively approximates the second-order optimization at every step. However, benefiting from the steep gradient, TFA+CG can reach convergence for all datasets within 100 iterations, while for SGD the number of iterations is much larger and it significantly increases with the size of the dataset. Overall, TFA+CG is about 30 times faster compared to TFA+SGD regarding time needed for convergence (T), and the time needed for feature extraction are close (T'). Additionally, the original TFA with SGD first pre-trains the weights for novel classes, which brings additional time costs, while CG with the steepest gradient optimizes the model weights from scratch and can delineate comparable results with much less time.

Overall, the models in the HDA family are fastest, compared to TFA+SGD, HDA+CG is about 200 to 250 times faster regarding the time needed for convergence (T). Compared to TFA, HDA has much fewer parameters to optimize as aforementioned, which also helps to speed up the optimization. For instance, the time per iteration of HDA+CG is almost halved compared to the one of TFA+CG, and it can reach convergence for datasets with different shots with only 30 iterations. The converging speed of HDA+CG+aug is slightly slower than HDA+CG due to the feature augmentation at every iteration. Regarding the time needed for feature extraction, HDA takes a longer time than TFA due to the process of proposal filtering. Table 5.2

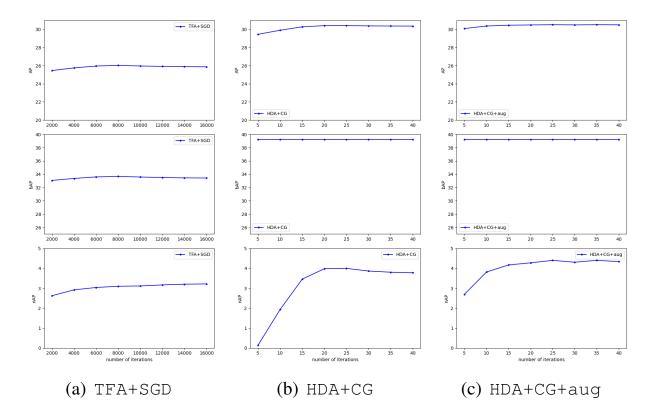


Figure 5.2: Convergence of different models on 1-shot COCO dataset (seed 0)

compares the overall time needed by HDA and TFA. For TFA+SGD, we summarize the time in both stages of pretraining novel weights and fine-tuning. With the time for feature extraction taken into account, HDA+CG is about 7 to 10 times faster than TFA+SGD.

Model	T' + T(s)									
	1	2	3	5	10					
TFA+SGD	121.53	245.46	364.83	617.83	1245.94					
HDA+CG	16.79	27.54	39.37	64.01	126.77					
HDA+CG+aug	14.42	27.92	42.28	62.16	137.56					

Table 5.2: Overall running time of different models

CHAPTER 5. DISCUSSION

# Chapter 6

# Conclusion

## 6.1 Summary

In this work, we tailor a steep gradient optimization for few-shot object detection based on Conjugate Gradient. Compared to traditional gradient descent, CG performs 30 times faster in optimization when applied to the previous benchmark of the two-stage tuning approach TFA. Additionally, we consider task-level knowledge that can refine the CG in TFA to make it even faster while minimizing overfitting. Experiment shows that in the 1-shot case for which the meta-learning was conducted, the predicted weight can make CG faster in beginning iterations and the feature-wise regularizer can effectively retain model performance on base classes. Furthermore, we propose a novel hierarchical detection approach (HDA) to address the issue of catastrophic forgetting. We observe that HDA can completely retain the model performance on base classes while achieving competitive results on novel classes. HDA's optimization is 200 to 250 times faster than SGD in TFA and overall run-time is 7 to 10 times faster than the previous benchmarks. We conducted extensive experiments on datasets sampled from different groups of random seeds and measure the model performance comprehensively. Our approach helps release the potential of few-shot object detection in real-time applications while without the worries of catastrophic forgetting.

## 6.2 Outlook

Future works that are well worth exploring include:

**Hierarchical detection setting**. Following the hierarchical detection approach proposed in the thesis, we observe a novel experimental setting of hierarchical detection that has not been explored in the existing literature. Perceiving objects hierarchically is natural for humans. For instance, when annotating the images in the COCO dataset, the researchers asked the labors to first decide the super-category of the object to annotate as it is more time-efficient [24]. With the HDA, we could further experiment the setting to first detect the super-category of an object and then decide its child class.

**Incremental learning**. Incremental learning is a common setting in robotics, where only the novel data is available and the robot needs to learn it without seriously forgetting learned knowledge [42]. Given the fast convergence of CG and the merit of HDA without catastrophic forgetting, it would be interesting to test our model in this setting.

**Efficient feature extraction**. In TFA+SGD, the time needed for convergence is dominant compared to the time needed for feature extraction. But with CG being applied, in HDA+CG, the time needed for convergence is largely shortened while the time needed for feature extraction becomes dominant. Researchers can

CHAPTER 6. CONCLUSION

look into a more efficient way to extract the proposals and box features to further speed up the overall time needed to run the model.

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