INTERLEAVING MULTI-TASK NEURAL ARCHITECTURE SEARCH

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ABSTRACT

Multi-task neural architecture search (MTNAS), which searches for a shared architecture for multiple tasks, has been broadly investigated. In these methods, multiple tasks are learned simultaneously by minimizing the weighted sum of their losses. How to balance these losses by finding the optimal loss weights requires a lot of tuning, which is time-consuming and labor intensive. To address this problem, we propose an interleaving MTNAS framework, where no tuning of loss weights is needed. In our method, a set of tasks (e.g., A, B, C) are performed in an interleaving loop (e.g., ABCABCABC...) where each task transfers its knowledge to the next task. Each task is learned by minimizing its loss function alone, without intervening with losses of other tasks. Loss functions of individual tasks are organized into a multi-level optimization framework which enables all tasks performed end-to-end. The effectiveness of our method is demonstrated in a variety of experiments.

1 Introduction

Neural architecture search (Zoph & Le, 2017; Liu et al., 2019a; Real et al., 2019), which aims at automatically searching for high-performance neural architectures with minimal human intervention, has attracted much attention recently and demonstrated promising effectiveness in a variety of machine learning tasks, including classification, segmentation, object detection, image captioning, etc. When multiple tasks are closely related, it is beneficial to search for a shared architecture for all tasks via multi-task learning. A number of methods have been developed for multi-task NAS (MTNAS) (Liang et al., 2018; Bruggemann et al., 2020; Gao et al., 2020). Existing MTNAS methods learn a shared architecture by minimizing the weighted sum of losses of all tasks. How to balance these loss terms is challenging: typically, more decrease of one loss renders less decrease of other losses; as a result, improving one task renders performance degradation of other tasks. To alleviate this problem, it is needed to carefully tune the tradeoff weights between loss terms (Mao et al., 2022), which is time-consuming and labor-intensive.

We aim to address this problem, by proposing an interleaving multi-task NAS framework. Different from existing MTNAS methods which perform multiple tasks by minimizing a combined loss function, our method is featured with multiple loss functions where each loss focuses on learning one task. This can avoid tuning the tradeoff weights between losses of different tasks. Different tasks are organized into a loop where each task can transfer its knowledge to any other task along the loop. Loss functions of individual tasks are unified into a multi-level optimization framework (Franceschi et al., 2018) which enables different tasks to be performed end-to-end.

In our framework, there are K tasks whose learnable parameters include network weights and a shared architecture. Each task has a data encoder and a task-specific head. Data encoders of all tasks share the same architecture, but have different network weights. The K tasks perform M rounds of interleaving learning. In the first round, we first learn task t_1 , then learn task t_2 , and so on. At the end of the first round, t_K is learned. Then we move to the second round, which starts with learning t_1 , then learns t_2 , and so on. This pattern repeats until the M rounds of learning are finished. Between two consecutive tasks t_k and t_{k+1} , knowledge transfer is performed based on distribution matching. Since all tasks are in a loop, the knowledge of one task can be propagated to any other task. After M rounds of learning, each task uses its model trained in the final round to make predictions on a validation dataset and updates their shared architecture by minimizing validation losses.

The major contributions of this paper are as follows:

- We propose an interleaving multi-task neural architecture search method, which searches for a shared architecture of multiple tasks in an interleaving fashion. Our framework solves multiple optimization problems sequentially, each focusing on training one task, which therefore can circumvent the difficulty of balancing different loss terms.
- We formulate interleaving MTNAS as a multi-level optimization problem which enables all tasks to be performed end-to-end.
- We propose a new knowledge transfer method based on distribution matching.
- Experiments on image classification and object detection demonstrate the effectiveness of our method.

2 Related works

2.1 MULTI-TASK LEARNING (MTL)

MTL (Ruder, 2017; Zhang & Yang, 2021) aims to improve multiple tasks simultaneously by learning them jointly and transferring knowledge across tasks. Various MTL approaches have been proposed, based on 1) hard parameter sharing (Caruana, 1998; Long et al., 2015; Doersch & Zisserman, 2017; Kokkinos, 2017; Sener & Koltun, 2018; Leang et al., 2020), where multiple tasks share the same weight parameters, such as encoder weights; 2) soft parameter sharing (Duong et al., 2015; Dai et al., 2016; Misra et al., 2016; Yang & Hospedales, 2016; Lu et al., 2017; Liu et al., 2019b; Maninis et al., 2019), where parameters of different tasks are constrained to be similar; 3) task similarity learning (Williams et al., 2007; Zhang & Yeung, 2014; Bingel & Søgaard, 2017; Standley et al., 2020; Zamir et al., 2020), which identifies similarity between tasks and encourages similar tasks to share more commonalities; 4) loss weighting (Chen et al., 2018; Kendall et al., 2018; Sener & Koltun, 2018; Gong et al., 2019; Leang et al., 2020), which weighs each task's loss. Existing methods mostly learn multiple tasks by minimizing the weighted sum of their losses, which requires careful tuning of tradeoff weights for balancing different loss terms. Different from these methods, our method performs MTL in an interleaving way. Zeng et al. (2019) proposed a cyclic MTL approach for neural simile recognition where several subtasks are learned in a loop. Tian et al. (2021) proposed to cyclically co-learn sounding object visual grounding and audio-visual sound separation. In these two approaches, subtasks are learned by minimizing the weighted sum of their training losses, which requires extensive tuning of tradeoff weights as well. In an audio source separation task, Doire & Okubadejo (2019) minimize the sum of losses of different sources by alternating mini-batches of data from different sources. This work suffers the difficulty of balancing loss terms since the mini-batch update of one source may lead to loss increase of other sources.

2.2 NEURAL ARCHITECTURE SEARCH (NAS)

The goal of neural architecture search (NAS) is to automatically identify highly-performing neural architectures that can potentially surpass human-designed ones. NAS research has made considerable progress in the past few years. Early NAS approaches (Zoph & Le, 2017; Pham et al., 2018; Zoph et al., 2018) are based on reinforcement learning, where a policy network learns to generate highquality architectures by maximizing the validation accuracy (as reward). These approaches are conceptually simple and can flexibly perform search in any search space. In differentiable search methods (Cai et al., 2019; Liu et al., 2019a; Xie et al., 2019), each candidate architecture is a combination of many building blocks. The combination coefficients represent the importance of building blocks. Architecture search amounts to learning these differentiable coefficients, which can be done using differentiable optimization algorithms such as gradient descent. Differentiable NAS methods started with DARTS (Liu et al., 2019a) and have been improved rapidly since then. For example, P-DARTS (Chen et al., 2019) allows the architecture depth to increase progressively during search. It also performs search space regularization and approximation to improve stability of search algorithms and reduce search cost. In PC-DARTS (Xu et al., 2020), the redundancy of search space exploration is reduced by sampling sub-networks from a super network. It also performs operation search in a subset of channels via bypassing the held-out subset in a shortcut. Another paradigm of NAS methods (Liu et al., 2018; Real et al., 2019) are based on evolutionary algorithms. In these approaches, architectures are considered as individuals in a population. Each architecture is associated with a fitness score representing how good this architecture is. Architectures with higher fitness scores have higher odds of generating offspring (new architectures), which replace architectures that have low-fitness scores. Our framework is orthogonal to existing NAS approaches and is applicable to most differentiable methods.

There are several works proposed for multi-task neural architecture search. Liang et al. (2018) proposed an evolutionary algorithm based multi-task NAS method which searches for a routing of modules for each task where candidate modules are shared by different tasks. Pasunuru & Bansal (2019) developed a framework for multi-task and continual NAS. Bruggemann et al. (2020) proposed branched multi-task architecture search, to automatically determine encoder branching in multi-task architectures. Gao et al. (2020) proposed a task-agnostic NAS framework for general-purpose multi-task learning. Silvestri (2020) proposed a one-shot NAS method for multi-task learning. Cai & Luo (2021) proposed a multi-task learning framework for conducting multi-objective evolutionary NAS. Similar to the MTL methods discussed in Section 2.1, these multi-task NAS methods require extensive tuning of tradeoff weights as well.

3 METHOD

In this section, we present details of the interleaving MTNAS framework. There are K tasks. These tasks could be the same (e.g., image classification on CIFAR-10, on CIFAR-100) or different (e.g., image classification on CIFAR-10, object detection on MS-COCO (Lin et al., 2014)). Each task k has a deep neural network model, consisting of a data encoder and a task-specific head. Given an input data example, it is fed into the encoder to generate a latent representation; the latent representation is fed into the head to predict an output label. Each task k has a training dataset $D_k^{\rm (tr)}$ and a validation dataset $D_k^{\rm (val)}$. The weight parameters of an encoder and head can be different in different rounds. Let $W_k^{\rm (m)}$ and $H_k^{\rm (m)}$ denote the encoder weights and head weights of task k in round k. Besides weight parameters, these tasks have a shared architecture k to learn. These tasks perform interleaving learning (with k rounds) in the following order:

$$\underbrace{t_1, t_2, \cdots, t_K}_{\text{Round 1}} \underbrace{t_1, t_2, \cdots, t_K}_{\text{Round } 2} \cdots \underbrace{t_1, t_2, \cdots, t_K}_{\text{Round } m} \cdots \underbrace{t_1, t_2, \cdots, t_K}_{\text{Round } M}$$
(1)

where t_k denotes that the k-th task performs learning. In the first round, t_1 is learned first, then t_2 is learned, then t_3 , etc. When learning t_{k+1} , we transfer knowledge from t_k to t_{k+1} . After all K tasks are learned in round 1, we move to the next round. When learning t_1 in the second round, we transfer knowledge from t_K learned in the first round to t_1 . Then t_2 is learned using knowledge transferred from t_1 , and so on. This interleaving procedure iterates for M rounds. Table 9 shows the notations of our method.

3.1 A MULTI-LEVEL OPTIMIZATION FRAMEWORK

There are $M \times K + 1$ learning stages: in each of the M rounds, each of the K tasks is learned at a stage; additionally, there is a final validation stage.

3.1.1 The first stage

At the very first stage, we train the model weights of task 1 in round 1, including encoder weights $W_1^{(1)}$ and task head $H_1^{(1)}$, by minimizing a training loss $L_{1,1}^{(tr)}$. The loss is defined on the model of task 1 in round 1 (including $W_1^{(1)}$, $H_1^{(1)}$, architecture A) and the training data $D_1^{(\text{tr})}$ of task 1. $L_{1,1}^{(tr)}$ is application specific. For example, in classification, $L_{1,1}^{(tr)}$ is a cross-entropy loss: given a data-label pair (x,y), the input data x is first fed into $W_1^{(1)}$ then $H_1^{(1)}$ to yield a prediction \hat{y} , and cross-entropy loss is defined on \hat{y} and ground-truth label y. A is needed to define $L_{1,1}^{(tr)}$, but is not optimized at this stage. Otherwise, a trivial solution of A will be yielded where A can perfectly overfit $D_1^{(\text{tr})}$ but has poor generalization capability. The optimization problem is:

$$\widetilde{W}_{1}^{(1)}(A), \widetilde{H}_{1}^{(1)}(A) = \operatorname{argmin}_{W_{1}^{(1)}, H_{1}^{(1)}} L_{1,1}^{(tr)}(A, W_{1}^{(1)}, H_{1}^{(1)}, D_{1}^{(\mathrm{tr})}). \tag{2}$$

 $\widetilde{W}_1^{(1)}(A)$ denotes that the optimal weights $\widetilde{W}_1^{(1)}$ depend on A since 1) the training loss is a function of A and 2) \widetilde{W}_1 depends on the training loss.

3.1.2 A MIDDLE STAGE

At any other learning stage, e.g., the k-th stage in the m-th round, we learn the model weights of the k-th task in the m-th round, including encoder weights $W_k^{(m)}$ and task head $H_k^{(m)}$, by solving two optimization problems. In the first problem, we train $W_k^{(m)}$ by transferring knowledge from

 $\widetilde{W}_{k-1}^{(m)}(A)$. Previous knowledge transfer approaches (Rajeswaran et al., 2019; Romero et al., 2014; Chen et al., 2020a) have two major limitations. First, they impose strong restrictions on model parameters (Rajeswaran et al., 2019) and data embeddings (Romero et al., 2014), for example, making the weight parameters of two models have a small L2 distance (Rajeswaran et al., 2019). Second, they are limited to capturing low-order (e.g., pairwise (Chen et al., 2020a)) relationships between data examples, which cannot capture the global properties of a dataset sufficiently and therefore hinders knowledge transfer through this dataset. To address the limitations of existing methods, we propose a new knowledge transfer approach based on distribution matching, which offers two benefits. First, it is more flexible in the sense that it does not impose strong restrictions that weights or embeddings need to be close in terms of L2 distance as previous methods do. Second, it measures similarity at the distribution level instead of individual example level, therefore can capture high-order (> 3) relationships between data examples and more holistic properties of an entire data distribution, which facilitates more effective knowledge transfer.

We generate sets of examples where examples in the same set are from the same distribution, then compare the distributions of two example-sets and see whether they are the same. To create a set of examples that are from the same distribution, we resort to data augmentation, based on the intuition that augmented examples created from the same original example can be considered as being from the same distribution. From each input training example x_i (excluding its output label) in task k and k-1, data augmentation is applied to it to generate a set of C augmented examples $\mathcal{A}_i = \{a_n^{(i)}\}_{n=1}^C$, which are considered as samples drawn from the same distribution since they all originate from x_i .

Given two augmentation sets \mathcal{A}_i and \mathcal{A}_j , we encourage $W_k^{(m)}$ and $\widetilde{W}_{k-1}^{(m)}(A)$ to be consistent in predicting whether the distributions of \mathcal{A}_i and \mathcal{A}_j are the same. We first use the optimal weights $\widetilde{W}_{k-1}^{(m)}$ of t_{k-1} to label whether \mathcal{A}_i and \mathcal{A}_j are from the same distribution in the following way. We use $\widetilde{W}_{k-1}^{(m)}$ to encode examples in \mathcal{A}_i and \mathcal{A}_j , and measure the maximum mean discrepancy (MMD) (Gretton et al., 2012) on the encodings. Let $z_n^{(i)}$ denote the embedding of $a_n^{(i)}$ by $\widetilde{W}_{k-1}^{(m)}$ and $f(\cdot,\cdot)$ be a kernel function. The MMD between \mathcal{A}_i and \mathcal{A}_j can be calculated as:

$$\frac{1}{C(C-1)} \sum_{n=1}^{C} \sum_{m \neq n}^{C} f(z_n^{(i)}, z_m^{(i)}) + \frac{1}{C(C-1)} \sum_{n=1}^{C} \sum_{m \neq n}^{C} f(z_n^{(j)}, z_m^{(j)}) - \frac{2}{C^2} \sum_{n=1}^{C} \sum_{m=1}^{C} f(z_n^{(i)}, z_m^{(j)}).$$
(3)

If the MMD is less than a threshold τ , A_i and A_j are considered to be from the same distribution.

Given these binary labels (regarding whether two augmentation sets are from the same distribution), we learn $W_k^{(m)}$ of t_k by fitting these labels. Specifically, we use $W_k^{(m)}$ to embed \mathcal{A}_i and \mathcal{A}_j , and calculate the MMD $d(\mathcal{A}_i,\mathcal{A}_j;W_k^{(m)})$. If \mathcal{A}_i and \mathcal{A}_j are labeled as being from the same distribution by $\widetilde{W}_{k-1}^{(m)}$, then we encourage $d(\mathcal{A}_i,\mathcal{A}_j;W_k^{(m)})$ to be less than τ . We use hinge loss to encourage the MMD calculated by $W_k^{(m)}$ and $\widetilde{W}_{k-1}^{(m)}(A)$ to be consistent: if they are both larger than τ or smaller than τ , there are no penalty. The optimization problem is:

$$\begin{split} \widetilde{W}_k^{(m)}(A) &= \mathrm{argmin}_{W_k^{(m)}} \ \sum_{\mathcal{A}_i, \mathcal{A}_j} \mathrm{max} \bigg(0, - \big(d(\mathcal{A}_i, \mathcal{A}_j; \widetilde{W}_{k-1}^{(m)}(A)) - \tau \big) \big(d(\mathcal{A}_i, \mathcal{A}_j; W_k^{(m)}) - \tau \big) \bigg), \end{split} \tag{4}$$

where the loss is defined on pairs of augmentation sets. In practice, we randomly sample a small number (specifically, 20) of input training examples to generate augmentation sets. Half of these examples are from task k-1 and the other half are from task k. The computation cost on the 20 sets is not high. When k=1 and m>1 (at the beginning of a round), the optimal encoder weights of the previous task is $\widetilde{W}_K^{(m-1)}(A)$ (at the end of previous round). The optimal encoder weights $\widetilde{W}_k^{(m)}$ depend on A. A is not updated at this learning stage, for the same reason described above.

In the second optimization problem, given the trained encoder $\widetilde{W}_k^{(m)}(A)$, we train the head $H_k^{(m)}$ by minimizing a training loss $L_{k,m}^{(tr)}$ defined on the model of task k in round m (including $\widetilde{W}_k^{(m)}(A)$, $H_k^{(m)}$, architecture A) and the training data $D_k^{(\mathrm{tr})}$ of task k. The second optimization problem is:

$$\widetilde{H}_{k}^{(m)}(A) = \operatorname{argmin}_{H_{k}^{(m)}} L_{k,m}^{(tr)}(A, \widetilde{W}_{k}^{(m)}(A), H_{k}^{(m)}, D_{k}^{(\operatorname{tr})}). \tag{5}$$

3.1.3 THE FINAL STAGE

At the final stage, each task k evaluates the loss $L_k^{(val)}$ of its model learned in the final round M (including encoder $\widetilde{W}_k^{(M)}(A)$, head $\widetilde{H}_k^{(M)}(A)$, architecture A) on its validation set $D_k^{(\mathrm{val})}$. A is learned by minimizing the validation losses of all tasks:

$$\min_{A} \sum_{k=1}^{K} L_{k}^{(val)}(A, \widetilde{W}_{k}^{(M)}(A), \widetilde{H}_{k}^{(M)}(A), D_{k}^{(val)}). \tag{6}$$

In experiments, these losses for different tasks are normalized to have similar scales. Note that our method focuses on alleviating the burden of tuning the weights of tasks' training losses. As for the validation losses, we follow the standard practice of the multi-task learning literature, which treats each task's validation performance (after normalization) equally.

3.1.4 MULTI-LEVEL OPTIMIZATION FRAMEWORK

To this end, we are ready to formulate the IMTNAS problem as a multi-level optimization framework. From bottom to top, the K tasks perform M rounds of interleaving learning. After the interleaving process is finished, the shared architecture A is optimized by minimizing validation losses.

$$\begin{split} & \underset{s.t.}{\min_{A}} \ \sum_{k=1}^{K} L_{k}^{(val)}(A, \widetilde{W}_{k}^{(M)}(A), \widetilde{H}_{k}^{(M)}(A), D_{k}^{(val)}) \\ & s.t. \\ & \cdots \\ & \widetilde{H}_{k}^{(m)}(A) = \underset{H_{k}}{\operatorname{argmin}}_{H_{k}^{(m)}} \ L_{k,m}^{(tr)}(A, \widetilde{W}_{k}^{(m)}(A), H_{k}^{(m)}, D_{k}^{(\operatorname{tr})}) \\ & \widetilde{W}_{k}^{(m)}(A) = \underset{W_{k}}{\operatorname{argmin}}_{W_{k}^{(m)}} \ \sum_{\mathcal{A}_{i}, \mathcal{A}_{j}} \max \bigg(0, - \big(d(\mathcal{A}_{i}, \mathcal{A}_{j}; \widetilde{W}_{k-1}^{(m)}(A)) - \tau \big) \Big(d(\mathcal{A}_{i}, \mathcal{A}_{j}; \widetilde{W}_{k}^{(m)}) - \tau \big) \bigg) \\ & \cdots \\ & \widetilde{W}_{1}^{(1)}(A), \widetilde{H}_{1}^{(1)}(A) = \underset{W_{1}^{(1)}, H_{1}^{(1)}}{\operatorname{argmin}}_{W_{1}^{(1)}, H_{1}^{(1)}} L_{1,1}^{(\operatorname{tr})}(A, W_{1}^{(1)}, H_{1}^{(1)}, D_{1}^{(\operatorname{tr})}) \end{split}$$

The optimization algorithm for solving the problem in Eq.(7) is deferred to the appendix.

Differentiable NAS. In differentiable NAS (Liu et al., 2019a), the search space consists of many candidate building blocks, whose outputs are multiplied with architecture variables. The search process learns these variables using gradient methods. After learning, architecture blocks with largest variable values are retained to form the final architecture.

Reduce computation and memory costs. To reduce computation and memory costs, we adopt a hypernetwork (Ha et al., 2016) based approach. A hypernetwork (with weight parameters V and a fixed architecture) takes a task number k, a round number r, and an architecture A as inputs, and generates the weight parameters $W_k^{(m)}$ for A. In this way, we do not need to store $W_k^{(m)}$ (for $m=1,\cdots,M$ and $k=1,\cdots,K$) in memory, but only need to store V, which is much smaller. The hypernetwork is designed as follows. We put all weight parameters into a long vector (with dimension D), then partition it into D/100 sub-vectors, each with a dimension of 100. The hypernetwork is a feedforward network which takes the architecture A, task index k, round index r, sub-vector index d as inputs and produces a 100-dimensional sub-vector. All the produced sub-vectors are concatenated to form the weight parameter vector. The hypernetwork has two hidden layers, with 30 and 50 hidden units respectively. The architecture A is represented as a continuous vector. The dimension of the vector is the number of operations in the search space. A_i denotes the importance weight of the i-th operation (i.e., whether it should be selected).

4 EXPERIMENTS

In this section, we report experimental results. We experimented with two multi-task settings: 1) two heterogeneous tasks: image classification and object detection; 2) three homogeneous tasks: image classification on three different datasets. In all experiments, in Eq.(3), the number C of augmentations in each set is set to 20. τ is set to 5. The kernel in MMD is set to a radial basis function kernel with the scale parameter set to 0.1. We use random crop (He et al., 2019), flip, rotation, and color jitter to generate augmentations for images. Please refer to the appendix for detailed hyperparameter settings and additional results, such as ablation study on the number C of augmented examples, memory costs, increasing the number of rounds, how the MMD threshold τ affects experimental results, etc.

4.1 DATASETS

For the image classification task, we used the following datasets: CIFAR-10 (Krizhevsky & Hinton, 2010), CIFAR-100 (Krizhevsky et al., 2009), and ImageNet (Deng et al., 2009). For the object

detection task, we used the COCO (Lin et al., 2014) dataset. CIFAR-10 contains 60K images from 10 classes. CIFAR-100 contains 60K images from 100 classes. For CIFAR-10 and CIFAR-100, each of them is split into train/validation/test sets with 25K/25K/10K images respectively. ImageNet has 1.2M training images and 50K test images, from 1000 classes. COCO is broadly used for common object detection. It contains a training set (train2017) of 118K images and a test set (val2017) of 5K images, from 80 classes. Following many previous NAS methods (e.g., DARTS, PDARTS, PCDARTS) which split the training data of CIFAR-10 and CIFAR-100 into a new training set and a validation set with a ratio of 1:1, we randomly split the 118K COCO training set into a new training and validation set with a ratio of 1:1. Baseline methods (when applicable) also use the same split.

4.2 Baselines

We compare with the following baselines. 1) Multi-task learning (MTL) with soft weight sharing (MTL-SWS) (Maninis et al., 2019): tasks are encouraged to have similar encoder weights via L2 regularization. 2) MTL with hard weight sharing (MTL-HWS) (Kokkinos, 2017): different tasks share the same encoder weights. 3) MTL sharing architecture only (MTL-SAO): encoder weights of tasks are independent. 4) MTL-NAS (Gao et al., 2020): task-agnostic NAS for general-purpose MTL. 5) BMTAS (Bruggemann et al., 2020): branched multi-task architecture search. 6) Minibatch interleaving (MI) (Doire & Okubadejo, 2019): the model is the same as MTL-SWS; in the training process, parameters of different tasks are updated using their minibatches alternatively. 7) Separate interleaving (SI) (Tian et al., 2021): we perform task 1 first, then use the architecture and weights of its encoder to initialize that of task 2, which are then finetuned; this process applies to other tasks as well and interleaves for multiple rounds. 8) Performing two tasks independently: SingPathNAS (Stamoulis et al., 2019), MobileNetV3 (Howard et al., 2019), MnasNet-A2 (Tan et al., 2019), MixNet-M (Tan & Le, 2019), and FairNAS (Chu et al., 2019).

4.3 MTL ON IMAGE CLASSIFICATION AND OBJECT DETECTION

In this section, the two tasks that are interleaved over are: image classification on ImageNet and object detection on COCO.

4.3.1 EXPERIMENTAL SETTINGS

For encoder architecture, its search space is the same as that in ProxylessNAS (Cai et al., 2019), which contains 19 layers and each layer contains 7 choices. For the objection detection head (ODH), we use the one in RetinaNet (Lin et al., 2017). Hyperparameter settings for the ODH are the same as those in (Lin et al., 2017). The classification head for ImageNet is a 1000-way linear classifier. We use 3 interleaving rounds, with an order of "cls,det,cls,···", where "cls" and "det" denote classification and detection tasks respectively. For image classification, we use top-1 and top-5 errors for evaluation. For object detection, we use Average Precision (AP) (Ren et al., 2015) as an evaluation metric, across IoU thresholds from 0.5 to 0.95 with an interval of 0.05 and different scales (small, medium, large). The model was trained for 40 epochs. Network weights were trained using SGD, with an initial learning rate of 0.02 (with cosine decay). Architecture variables were trained using Adam (Kingma & Ba, 2014) with a learning rate of 5e-4. Batch size was 1024 for ImageNet and 12 for COCO.

4.3.2 RESULTS AND ANALYSIS

Table 1 shows object detection results on the COCO test set and classification results on the ImageNet test set. From this table, we make the following observations. **First**, IMTNAS performs better than MTL-SWS. In MTL-SWS, object detection and image classification are performed at the same time by minimizing the weighted sum of their losses. More loss decrease of one task leads to less loss decrease of the other task. To verify this, we plot how test performances of these two tasks (mAP of the detection task on the left axis and accuracy of the classification task on the right axis) vary with the weight β of the classification task loss (weight of the detection loss is set to 1) in Figure 1(Left). As β increases (more attention is paid to the classification task), top-1 test accuracy on ImageNet increases while test mAP on COCO decreases. This demonstrates that improving the performance of one task leads to the performance degradation of the other task. In contrast, each loss in our method focuses on learning one task at a time, thus can avoid this problem.

On the other hand, the interleaving mechanism in IMTNAS enables each task to transfer its knowledge to any other task along the interleaving loop. Figure 1(Right) shows the test performance of models in a 3-round interleaving process. As the number of rounds increases, detection models and classification models perform better. This demonstrates that via interleaving, each task continuously improves by leveraging knowledge from previous tasks. From Figure 1(Right), we see that our method avoids

Table 1: Object detection on COCO test set and classification accuracy on ImageNet test set. APS, APM, APL: AP at small, medium, large scale. AP50 and AP75 are AP with IoU thresholds of 0.5 and 0.75. Acc is accuracy.

		COCO						Imag	ImageNet	
	Acc	AP	AP50	AP75	APS	APM	APL	Top-1	Top-5	(M)
SingPathNAS (Stamoulis et al., 2019)	75.0	30.7	49.8	32.2	15.4	33.9	41.6	75.0	92.2	365
MobileNetV3 (Howard et al., 2019)	75.2	29.9	49.3	30.8	14.9	33.3	41.1	75.2	92.2	219
MnasNet-A2 (Tan et al., 2019)	75.6	30.5	50.2	32.0	16.6	34.1	41.1	75.6	92.7	340
MixNet-M (Tan & Le, 2019)	77.0	31.3	51.7	32.4	17.0	35.0	41.9	77.0	93.3	360
FairNAS-C (Chu et al., 2019)	76.7	31.2	50.8	32.7	16.3	34.4	42.3	76.7	93.3	325
MTL-SAO	76.8	31.0	51.2	33.4	16.1	34.6	41.6	76.3	92.8	379
MTL-HWS (Kokkinos, 2017)	77.0	31.5	51.5	33.2	16.5	34.8	41.7	76.8	93.1	384
MTL-SWS (Maninis et al., 2019)	77.3	31.9	51.8	33.8	16.6	35.1	42.1	77.1	93.4	395
MTL-NAS (Gao et al., 2020)	77.2	31.3	51.1	33.5	16.1	34.4	41.7	76.5	93.0	495
BMTAS (Bruggemann et al., 2020)	76.3	30.5	51.0	33.1	16.2	34.3	41.3	76.3	92.6	432
MI (Doire & Okubadejo, 2019)	77.1	31.4	51.5	33.4	16.4	34.7	41.6	76.8	93.2	389
SI (Tian et al., 2021)	77.1	31.1	50.6	32.7	15.8	34.1	41.4	76.5	92.7	371
IMTNAS (ours)	77.7	32.5	53.1	34.1	17.1	36.8	43.4	77.8	93.9	381

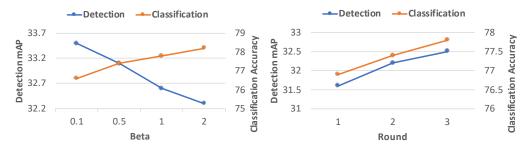


Figure 1: (Left) In MTL-SWS, improving one task leads to performance degradation of the other task. (Right) In our method, how task performance varies with the number of rounds.

catastrophic forgetting (Kirkpatrick et al., 2017). This is because in our method, all tasks are learned jointly in a unified framework where every task can transfer knowledge to any other task and all tasks learn the shared encoder architecture together.

Second, our method performs better than MTL-HWS. It is because MTL-HWS requires two tasks to share exactly the same encoder, which is not good for task-specific representation learning. The datasets of different tasks have different properties. We need to use task-specific encoders with different weight parameters to learn task-specific representations for these datasets. MTL-HWS, which makes the encoders of all tasks share the same weights, compromises this purpose. Third, our method outperforms MTL-SAO. In MTL-SAO, two tasks only share encoder architecture and do not share commonality in encoder weights. Weights also carry important knowledge, which should be transferred between tasks. Fourth, our method performs better than MTL-NAS and BMTAS. Similar to MTL-SWS, in these two methods, more decrease of the loss in one task renders less decrease of the loss in the other task. Fifth, our method achieves better performance than SI. In SI, different tasks are trained separately while our method performs two tasks in an end-to-end framework. Sixth, our method outperforms baselines that perform two tasks independently. These methods do not conduct knowledge transfer between tasks. Seventh, the computational demand (the number of composite multiply-accumulate (×+) operations for a single image) of our method is comparable with baselines.

4.4 MULTI-TASK CLASSIFICATION ON THREE TASKS

In this section, we report experimental results on three tasks: image classification on CIFAR-100, image classification on CIFAR-10, and image classification on ImageNet.

4.4.1 EXPERIMENTAL SETTINGS

Following the experimental protocol in (Liu et al., 2019a), each experiment consists of an architecture search phase and an architecture evaluation phase. In the search phase, an optimal architecture cell is searched by minimizing a validation loss. In the evaluation phase, a larger network is created by stacking multiple copies of an optimally searched cell. This new network is re-trained from scratch and evaluated on a test set. For ImageNet, following (Xu et al., 2020), we randomly sample 10%

Table 2: Classification errors (%) on test sets of CIFAR-100, CIFAR-10, and ImageNet, number of model parameters (millions), and search cost (GPU days on a Nvidia 1080Ti, sum of costs of all tasks). Results on ImageNet are under mobile settings. IMTNAS-Darts2nd: our IMTNAS is applied to DARTS-2nd. Results marked with * are taken from DARTS $^-$ (Chu et al., 2020), DrNAS (Chen et al., 2020b), β -DARTS (Ye et al., 2022), and AGNAS (Sun et al., 2022). Methods marked with † were re-run 10 times with random initialization.

	CIFAR100 Error(%)	CIFAR10 Error(%)	Imag Top-1	geNet Top-5	Param (M)	Cost (GPU days)
*Darts2nd (Liu et al., 2019a)	20.58±0.44	2.76±0.09	26.7	8.6	3.1/3.3/4.7	4.0/4.0/4.0
†Pdarts (Chen et al., 2019)	17.42 ± 0.14	2.54 ± 0.04	24.4	7.3	3.6/3.5/4.9	0.3/0.3/0.3
*R-DARTS (Zela et al., 2020)	17.98 ± 0.21	2.92 ± 0.24	-	-	-/-/-	1.6/1.6/-
†Prdarts (Zhou et al., 2020)	16.48 ± 0.06	2.37 ± 0.03	24.1	7.3	3.4/3.5/5.0	0.2/0.2/0.2
*ISTA-NAS (Yang et al., 2020)	-	2.51 ± 0.04	24.0	7.1	-/3.3/5.7	-/0.1/4.2
*MiLeNAS (He et al., 2020)	-	2.49 ± 0.08	24.7	7.6	-/3.9/5.3	-/0.3/0.3
*PDARTS-ADV (Chen & Hsieh, 2020)	_	2.44 ± 0.05	24.2	7.2	-/3.4/-	-/1.1/-
*DARTS+PT (Wang et al., 2021)	-	2.61 ± 0.08	25.5	8.0	-/3.0/4.6	-/0.8/-
*GAEA (Li et al., 2021)	-	2.50 ± 0.06	24.0	7.3	-/-/-	-/0.1/3.8
*DOTS (Gu et al., 2021)	16.48 ± 0.13	2.49 ± 0.06	24.0	7.2	4.1/3.5/5.3	0.3/0.3/1.3
* β -DARTS (Ye et al., 2022)	16.24 ± 0.22	2.53 ± 0.08	23.9	7.0	3.8/3.8/5.5	0.4/0.4/0.4
*AGNAS (Sun et al., 2022)	-	$2.53 \pm .003$	23.4	6.8	-/3.6/6.7	-/0.4/3.3
†Pcdarts (Xu et al., 2020)	17.96±0.15	2.57 ± 0.07	24.2	7.3	4.3=(3.9+3.6+5.3)/3	4.0=0.1+0.1+3.8
MTL-NAS-Pcdarts (Gao et al., 2020)	19.38 ± 0.20	2.81 ± 0.13	26.4	8.6	4.9	12.9
BMTAS-Pcdarts (Bruggemann et al., 2020)	20.14 ± 0.26	2.78 ± 0.11	27.1	9.2	4.3	10.5
MTL-SAO-Pcdarts	18.09 ± 0.27	2.68 ± 0.07	24.4	7.5	4.5	4.2
MTL-HWS-Pcdarts (Kokkinos, 2017)	18.10 ± 0.11	2.69 ± 0.14	24.5	7.6	4.1	4.4
MTL-SWS-Pcdarts (Maninis et al., 2019)	17.93±0.36	$2.54{\pm}0.08$	24.5	7.4	4.3	4.2
MI-Pcdarts (Doire & Okubadejo, 2019)	18.15±0.19	2.68 ± 0.10	24.6	7.7	4.0	4.1
SI-Pcdarts (Tian et al., 2021)	17.96 ± 0.09	2.66 ± 0.07	25.3	8.1	4.2	4.5
IMTNAS-Pcdarts (ours)	16.08 ±0.05	2.31 ± 0.03	23.1	6.3	4.1=(3.7+3.7+5.0)/3	4.0

Table 3: Ablation on knowledge transfer approaches.

Method	CIFAR-100 Error (%)	CIFAR-10 Error (%)	ImageNet Top-1/Top-5 Error
L2-Weights	17.41 ± 0.07	2.55±0.05	24.4/7.6
L2-Embedding	17.28 ± 0.10	2.51 ± 0.08	24.1/7.3
Example-Similarity	17.12 ± 0.15	2.49 ± 0.09	23.9/7.1
Ours	16.08 ±0.05	2.31 ±0.03	23.1/6.3

of the original 1.3M training images as a new training set and randomly sample another 2.5% of the 1.3M images as a validation set. IMTNAS was applied to PC-DARTS in consideration of its computational efficiency. The classification head for CIFAR-10, CIFAR-100, and ImageNet is a 10-way, 100-way, and 1000-way linear classifier respectively. We set the number of interleaving rounds to 2, with the following task order: CIFAR-100, CIFAR-10, ImageNet, CIFAR-100, CIFAR-10, ImageNet. During architecture search, network weights were optimized using the SGD optimizer with a batch size of 64, an initial learning rate of 0.025, a learning rate scheduler of cosine decay, a weight decay of 3e-4, a momentum of 0.9, and an epoch number of 50 (only for our method; the number of epochs for baselines are given in the appendix). Architecture variables were optimized using the Adam (Kingma & Ba, 2014) optimizer with a learning rate of 3e-4 and a weight decay of 1e-3. The rest of hyperparameters follow those in PC-DARTS. The experiments were conducted on Nvidia 1080Ti GPU. Mean and standard deviation of classification errors obtained from 10 random runs are reported.

4.4.2 RESULTS AND ANALYSIS

Table 2 shows results on CIFAR-100, CIFAR-10, and ImageNet. When our proposed IMTNAS framework is applied to PCDARTS, the errors of PCDARTS are greatly reduced, which demonstrates the effectiveness of IMTNAS. In the interleaving process, each task transfers its knowledge to the next task. To ensure successful transfer, each task needs to learn representations not only good for itself, but also good for the next task. By doing this, the learned representations have good generalizability across tasks.

Other observations made from Table 2 are similar to those from Table 1: 1) our method works better than MTL baselines including MTL-SAO, MTL-HWS, MTL-SWS, SI, MTL-NAS, and BMTAS; 2) our method outperforms baselines which conduct different tasks separately. The analysis of reasons

Figure 2: How classification errors (%) on CIFAR-100, CIFAR-10, and ImageNet change with the number of interleaving rounds M.

Table 4: Ablation study on task order.

Order	CIFAR-100 Error	CIFAR-10 Error	ImageNet Top-1/Top-5 Error
CIFAR-100, CIFAR-10, ImageNet, CIFAR-100, CIFAR-10, ImageNet	16.08 ± 0.05	2.31 ± 0.03	23.1/6.3
CIFAR-100, ImageNet, CIFAR-10, CIFAR-100, ImageNet, CIFAR-10	16.11 ± 0.09	2.30 ± 0.04	23.3/6.4
ImageNet, CIFAR-10, CIFAR-100, ImageNet, CIFAR-10, CIFAR-100	16.10 ± 0.03	2.34 ± 0.04	23.1/6.4

is similar to that for results in Table 1. The computational costs and number of model parameters of our methods are similar to those of baselines, while our method achieves better accuracy.

4.4.3 ABLATION STUDIES

We compare our distribution matching based knowledge transfer (KT) approach with several other KT methods, including 1) L2-Weights (Rajeswaran et al., 2019): encouraging the L2 distance between $\widetilde{W}_{k-1}^{(m)}(A)$ and $W_k^{(m)}$ to be small; 2) L2-Embedding (Romero et al., 2014): encouraging the embeddings of input images extracted by $\widetilde{W}_{k-1}^{(m)}(A)$ and $W_k^{(m)}$ to have small L2 distance; and 3) Example-Similarity (Chen et al., 2020a): using $\widetilde{W}_{k-1}^{(m)}(A)$ to label whether a pair of input examples are similar and learn $W_k^{(m)}$ by fitting these labels. Table 3 shows the results. Our method outperforms L2-Weights and L2-Embedding because these two baselines impose strong restrictions that weights or embeddings need to be close in an L2 sense while our method is more flexible. Our method works better than Example-Similarity. This method is limited to capturing low-order relationships between data examples (the order is two specifically, since pairwise similarity is measured). In contrast, our method measures similarity at the distribution level instead of individual example level, therefore can capture high-order relationships between data examples and more holistic properties of an entire data distribution.

We study how the test error of the final model in an interleaving sequence changes with the number of interleaving rounds M. Figure 2 shows the results. When M increases from 1 to 2, the errors are reduced. When M=1, the interleaving effect is weak: classification on CIFAR-100 influences classification on CIFAR-10 and subsequently on ImageNet but not the other way around. When M=2, the interleaving effect is strong: all tasks influence each other. This further demonstrates the effectiveness of interleaving learning. Increasing M from 2 to 3 does not significantly reduce the errors further. This is probably because 2 rounds of interleaving have brought in a sufficient interleaving effect.

Next, we explore whether task order affects performance. We experimented with three orders (with the number of rounds set to 2): 1) CIFAR-100, CIFAR-10, ImageNet, CIFAR-100, CIFAR-10, ImageNet; 2) CIFAR-100, ImageNet, CIFAR-100, ImageNet, CIFAR-100, ImageNet, CIFAR-100, CIFAR-100, ImageNet, CIFAR-100, Table 4 shows the test errors under these different orders. As can be seen, the errors are not affected by the task order significantly. The reason is: via interleaving, each task influences the other task at some point in the interleaving sequence; therefore, it does not matter too much regarding which task should be performed first.

5 CONCLUSIONS AND DISCUSSIONS

We propose an interleaving multi-task neural architecture search framework, where multiple tasks are performed in an interleaving fashion. Via interleaving, different models transfer their knowledge to each other. Experiments on various datasets demonstrate the efficacy of our method. One major limitation of this work is that it requires the NAS methods to be differentiable. For non-differentiable NAS methods such as those based on reinforcement learning and evolutionary algorithms, our method is not applicable.

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A OPTIMIZATION ALGORITHM

We use a well-established algorithm developed in (Liu et al., 2019a) to solve the interleaving learning problem. Theoretic convergence of this algorithm has been broadly analyzed in (Ghadimi & Wang, 2018; Grazzi et al., 2020; Ji et al., 2021; Liu et al., 2021; Yang et al., 2021). For the ease of presentation, we first present the algorithm for a special case of our method where the number of tasks is two and the number of interleaving rounds is two as well. Extensions to more tasks and more rounds can be straightforwardly conducted. In the second subsection, we will introduce a general version of this algorithm. At each level of optimization problem, the optimal solution (on the left-hand side of the equal sign, marked with $\widetilde{\cdot}$), its exact value is computationally expensive to compute. To address this problem, following (Liu et al., 2019a), we approximate the optimal solution using a one-step gradient descent update and plug the approximation into the next level of optimization problem. In the sequel, $\frac{\partial \cdot}{\partial \cdot}$ denotes partial derivative. $\frac{d\cdot}{d\cdot}$ denotes an ordinary derivative.

A.1 SPECIAL CASE

We consider the following special case of our method with two tasks and two interleaving rounds.

$$\begin{split} & \min_{A} L_{1}^{(val)}(A, \widetilde{W}_{1}^{(2)}(A), \widetilde{H}_{1}^{(2)}(A), D_{1}^{(val)}) + L_{2}^{(val)}(A, \widetilde{W}_{2}^{(2)}(A), \widetilde{H}_{2}^{(2)}(A), D_{2}^{(val)}) \\ & s.t. \\ & \widetilde{H}_{2}^{(2)}(A) = \operatorname{argmin}_{H_{2}^{(2)}} L_{2,2}^{(tr)}(A, \widetilde{W}_{2}^{(2)}(A), H_{2}^{(2)}, D_{2}^{(tr)}) \\ & \widetilde{W}_{2}^{(2)}(A) = \operatorname{argmin}_{W_{2}^{(2)}} \sum_{\mathcal{A}_{i}, \mathcal{A}_{j}} \max \left(0, -\left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; A, \widetilde{W}_{1}^{(2)}(A)) - \tau \right) \left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; A, W_{2}^{(2)}) - \tau \right) \right) \\ & \widetilde{H}_{1}^{(2)}(A) = \operatorname{argmin}_{H_{1}^{(2)}} L_{1,2}^{(tr)}(A, \widetilde{W}_{1}^{(2)}(A), H_{1}^{(2)}, D_{1}^{(tr)}) \\ & \widetilde{W}_{1}^{(2)}(A) = \operatorname{argmin}_{W_{1}^{(2)}} \sum_{\mathcal{A}_{i}, \mathcal{A}_{j}} \max \left(0, -\left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; A, \widetilde{W}_{2}^{(1)}(A)) - \tau \right) \left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; A, W_{1}^{(2)}) - \tau \right) \right) \\ & \widetilde{H}_{2}^{(1)}(A) = \operatorname{argmin}_{H_{2}^{(1)}} L_{2,1}^{(tr)}(A, \widetilde{W}_{2}^{(1)}(A), H_{2}^{(1)}, D_{2}^{(tr)}) \\ & \widetilde{W}_{2}^{(1)}(A) = \operatorname{argmin}_{W_{2}^{(1)}} \sum_{\mathcal{A}_{i}, \mathcal{A}_{j}} \max \left(0, -\left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; A, \widetilde{W}_{1}^{(1)}(A)) - \tau \right) \left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; A, W_{2}^{(1)}) - \tau \right) \right) \\ & \widetilde{W}_{1}^{(1)}(A), \widetilde{H}_{1}^{(1)}(A) = \operatorname{argmin}_{W_{1}^{(1)}, H_{1}^{(1)}} L_{1,1}^{(tr)}(A, W_{1}^{(1)}, H_{1}^{(1)}, D_{1}^{(tr)}) \end{aligned} \tag{8}$$

Next, we present the algorithm for solving this special case. We approximate $\widetilde{W}_1^{(1)}(A)$ using one-step gradient descent update of $W_1^{(1)}$ w.r.t $L_{1,1}^{(tr)}(A,W_1^{(1)},H_1^{(1)},D_1^{(\mathrm{tr})})$:

$$\widetilde{W}_{1}^{(1)}(A) \approx \widehat{W}_{1}^{(1)} = W_{1}^{(1)} - \eta_{w} \nabla_{W_{1}^{(1)}} L_{1,1}^{(tr)}(A, W_{1}^{(1)}, H_{1}^{(1)}, D_{1}^{(\text{tr})})$$

$$\tag{9}$$

We update $H_1^{(1)}$ using gradient descent:

$$H_1^{(1)} \leftarrow H_1^{(1)} - \eta_h \nabla_{H_1^{(1)}} L_{1,1}^{(tr)}(A, W_1^{(1)}, H_1^{(1)}, D_1^{(tr)})$$
 (10)

We plug the approximation $\widetilde{W}_1^{(1)}(A) \approx \widehat{W}_1^{(1)}$ into $\sum_{\mathcal{A}_i,\mathcal{A}_j} \max \left(0,-\left(d(\mathcal{A}_i,\mathcal{A}_j;A,\widetilde{W}_1^{(1)}(A))-\tau\right)\right)$ and get an approximated objective. Then we approximate $\widetilde{W}_2^{(1)}(A)$ using one-step gradient descent update of $W_2^{(1)}$:

$$\widetilde{W}_{2}^{(1)}(A) \approx \widehat{W}_{2}^{(1)} = W_{2}^{(1)} - \eta_{w} \nabla_{W_{2}^{(1)}} \sum_{\mathcal{A}_{i}, \mathcal{A}_{j}} \max \left(0, -\left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; A, \widehat{W}_{1}^{(1)}) - \tau \right) \left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; A, W_{2}^{(1)}) - \tau \right) \right)$$

$$(11)$$

We plug the approximation $\widetilde{W}_2^{(1)}(A) \approx \widehat{W}_2^{(1)}$ into $L_{2,1}^{(tr)}(A,\widetilde{W}_2^{(1)}(A),H_2^{(1)},D_2^{(\mathrm{tr})})$ and get an approximated objective. Then we approximate $\widetilde{H}_2^{(1)}(A)$ using one-step gradient descent update of $H_2^{(1)}$:

$$\widetilde{H}_{2}^{(1)}(A) \approx \widehat{H}_{2}^{(1)} = H_{2}^{(1)} - \eta_{h} \nabla_{H_{2}^{(1)}} L_{2,1}^{(tr)}(A, \widehat{W}_{2}^{(1)}, H_{2}^{(1)}, D_{2}^{(tr)})$$
(12)

We plug the approximation $\widetilde{W}_2^{(1)}(A) \approx \widehat{W}_2^{(1)}$ into $\sum_{\mathcal{A}_i,\mathcal{A}_j} \max \left(0, -\left(d(\mathcal{A}_i,\mathcal{A}_j;A,\widetilde{W}_2^{(1)}(A)) - \tau\right)\right)$ and get an approximated objective. Then we approximate $\widetilde{W}_1^{(2)}(A)$ using one-step gradient descent update of $W_1^{(2)}$:

$$\widetilde{W}_{1}^{(2)}(A) \approx \widehat{W}_{1}^{(2)} = W_{1}^{(2)} - \eta_{w} \sum_{\mathcal{A}_{i}, \mathcal{A}_{j}} \max \left(0, -\left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; A, \widehat{W}_{2}^{(1)}) - \tau \right) \left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; A, W_{1}^{(2)}) - \tau \right) \right)$$
(13)

We plug the approximation $\widetilde{W}_1^{(2)}(A) \approx \widehat{W}_1^{(2)}$ into $L_{1,2}^{(tr)}(A,\widetilde{W}_1^{(2)}(A),H_1^{(2)},D_1^{(\mathrm{tr})})$ and get an approximated objective. Then we approximate $\widetilde{H}_1^{(2)}(A)$ using one-step gradient descent update of $H_1^{(2)}$:

$$\widetilde{H}_{1}^{(2)}(A) \approx \widehat{H}_{1}^{(2)} = H_{1}^{(2)} - \eta_{h} \nabla_{H_{1}^{(2)}} L_{1,2}^{(tr)}(A, \widehat{W}_{1}^{(2)}, H_{1}^{(2)}, D_{1}^{(tr)})$$
(14)

We plug the approximation $\widetilde{W}_1^{(2)}(A) \approx \widehat{W}_1^{(2)}$ into $\sum_{\mathcal{A}_i,\mathcal{A}_j} \max \left(0,-\left(d(\mathcal{A}_i,\mathcal{A}_j;A,\widetilde{W}_1^{(2)}(A))-\tau\right)\right)$ and get an approximated objective. Then we approximate $\widetilde{W}_2^{(2)}(A)$ using one-step gradient descent update of $W_2^{(2)}$:

$$\widetilde{W}_{2}^{(2)}(A) \approx \widehat{W}_{2}^{(2)} = W_{2}^{(2)} - \eta_{w} \nabla_{W_{2}^{(2)}} \sum_{\mathcal{A}_{i}, \mathcal{A}_{j}} \max \left(0, -\left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; A, \widehat{W}_{1}^{(2)}) - \tau \right) \left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; A, W_{2}^{(2)}) - \tau \right) \right)$$

$$\tag{15}$$

We plug the approximation $\widetilde{W}_2^{(2)}(A) \approx \widehat{W}_2^{(2)}$ into $L_{2,2}^{(tr)}(A,\widetilde{W}_2^{(2)}(A),H_2^{(2)},D_2^{(\mathrm{tr})})$ and get an approximated objective. Then we approximate $\widetilde{H}_2^{(2)}(A)$ using one-step gradient descent update of $H_2^{(2)}$:

$$\widetilde{H}_{2}^{(2)}(A) \approx \widehat{H}_{2}^{(2)} = H_{2}^{(2)} - \eta_{h} \nabla_{H_{2}^{(2)}} L_{2,2}^{(tr)} (A, \widehat{W}_{2}^{(2)}, H_{2}^{(2)}, D_{2}^{(tr)})$$
(16)

Finally, we plug the approximations $\widetilde{W}_1^{(2)}(A) \approx \widehat{W}_1^{(2)}$, $\widetilde{H}_1^{(2)}(A) \approx \widehat{H}_1^{(2)}$, $\widetilde{W}_2^{(2)}(A) \approx \widehat{W}_2^{(2)}$, $\widetilde{H}_2^{(2)}(A) \approx \widehat{H}_2^{(2)}$ into the validation loss and get an approximated validation loss. We update A by minimizing the approximated validation loss.

$$A \leftarrow A - \eta_a \nabla_A(L_1^{(val)}(A, \widehat{W}_1^{(2)}, \widehat{H}_1^{(2)}, D_1^{(val)}) + L_2^{(val)}(A, \widehat{W}_2^{(2)}, \widehat{H}_2^{(2)}, D_2^{(val)}))$$
(17)

 $abla_A L_2^{(val)}(A,\widehat{W}_2^{(2)},\widehat{H}_2^{(2)},D_2^{(\mathrm{val})})$ can be calculated as:

$$\nabla_{A} L_{2}^{(val)}(A, \widehat{W}_{2}^{(2)}, \widehat{H}_{2}^{(2)}, D_{2}^{(val)}) = \frac{\partial L_{2}^{(val)}(A, \widehat{W}_{2}^{(2)}, \widehat{H}_{2}^{(2)}, D_{2}^{(val)}) + d\widehat{W}_{2}^{(2)}}{\partial A} + \frac{d\widehat{W}_{2}^{(2)}}{\partial A} \frac{\partial L_{2}^{(val)}(A, \widehat{W}_{2}^{(2)}, \widehat{H}_{2}^{(2)}, D_{2}^{(val)})}{\partial \widehat{W}_{2}^{(2)}} + (18)$$

$$(\frac{\partial \widehat{H}_{2}^{(2)}}{\partial A} + \frac{d\widehat{W}_{2}^{(2)}}{dA} \frac{\partial \widehat{H}_{2}^{(2)}}{\partial \widehat{W}_{2}^{(2)}}) \frac{\partial L_{2}^{(val)}(A, \widehat{W}_{2}^{(2)}, \widehat{H}_{2}^{(2)}, D_{2}^{(val)})}{\partial \widehat{H}_{2}^{(2)}}$$

where

$$\frac{d\widehat{W}_{2}^{(2)}}{dA} = \frac{d\widehat{W}_{1}^{(2)}}{dA} \frac{\partial \widehat{W}_{2}^{(2)}}{\partial \widehat{W}_{*}^{(2)}} + \frac{\partial \widehat{W}_{2}^{(2)}}{\partial A}$$
(19)

where

$$\frac{d\widehat{W}_{1}^{(2)}}{dA} = \frac{d\widehat{W}_{2}^{(1)}}{dA} \frac{\partial \widehat{W}_{1}^{(2)}}{\partial \widehat{W}_{2}^{(1)}} + \frac{\partial \widehat{W}_{1}^{(2)}}{\partial A}$$
(20)

where

$$\frac{d\widehat{W}_{2}^{(1)}}{dA} = \frac{d\widehat{W}_{1}^{(1)}}{dA} \frac{\partial \widehat{W}_{2}^{(1)}}{\partial \widehat{W}_{1}^{(1)}} + \frac{\partial \widehat{W}_{2}^{(1)}}{\partial A}$$
(21)

where

$$\frac{d\widehat{W}_{1}^{(1)}}{dA} = -\eta_{w} \nabla_{A, W_{1}^{(1)}}^{2} L_{1, 1}^{(tr)}(A, W_{1}^{(1)}, H_{1}^{(1)}, D_{1}^{(tr)})$$
(22)

Algorithm 1 Optimization algorithm for the special case

while not converged do

Update the approximation $\widehat{W}_{1}^{(1)}$ of $\widetilde{W}_{1}^{(1)}(A)$ using Eq.(9)

Update $H_1^{(1)}$ using Eq.(10)

Update the approximation $\widehat{W}_2^{(1)}$ of $\widetilde{W}_2^{(1)}(A)$ using Eq.(11) Update the approximation $\widehat{H}_2^{(1)}$ of $\widetilde{H}_2^{(1)}(A)$ using Eq.(12) Update the approximation $\widehat{W}_1^{(2)}$ of $\widehat{W}_1^{(2)}(A)$ using Eq.(13) Update the approximation $\widehat{H}_1^{(2)}$ of $\widehat{H}_1^{(2)}(A)$ using Eq.(14) Update the approximation $\widehat{W}_2^{(2)}$ of $\widehat{W}_2^{(2)}(A)$ using Eq.(15) Update the approximation $\widehat{H}_2^{(2)}$ of $\widehat{H}_2^{(2)}(A)$ using Eq.(16) Update $\widehat{H}_2^{(2)}(A)$ using Eq.(17)

Update A using Eq.(17)

end while

 $\nabla_A L_1^{(val)}(A,\widehat{W}_1^{(2)},\widehat{H}_1^{(2)},D_1^{(val)})$ can be calculated as:

$$\frac{\partial \nabla_{A} L_{1}^{(val)}(A, \widehat{W}_{1}^{(2)}, \widehat{H}_{1}^{(2)}, D_{1}^{(val)})}{\partial A} + \frac{d\widehat{H}_{1}^{(2)}}{dA} \frac{\partial \nabla_{A} L_{1}^{(val)}(A, \widehat{W}_{1}^{(2)}, \widehat{H}_{1}^{(2)}, D_{1}^{(val)})}{\partial \widehat{H}_{1}^{(2)}} + \frac{d\widehat{W}_{1}^{(2)}}{\partial \widehat{W}_{1}^{(2)}} \frac{\partial \nabla_{A} L_{1}^{(val)}(A, \widehat{W}_{1}^{(2)}, \widehat{H}_{1}^{(2)}, D_{1}^{(val)})}{\partial \widehat{W}_{1}^{(2)}}$$
(23)

where

$$\frac{d\widehat{H}_{1}^{(2)}}{dA} = \frac{\partial \widehat{H}_{1}^{(2)}}{\partial A} + \frac{d\widehat{W}_{1}^{(2)}}{dA} \frac{\partial \widehat{H}_{1}^{(2)}}{\partial \widehat{W}_{1}^{(2)}}$$
(24)

and $\frac{d\widehat{W}_{1}^{(2)}}{dA}$ is given in Eq.(20).

The gradient descent update of A in Eq.(17) can run one or more steps. After A is updated, the one-step gradient-descent approximations (in equation 2-9), which are functions of A, change with Aand need to be re-updated. Then, the gradient of A, which is a function of one-step gradient-descent approximations, needs to be re-calculated and is used to refresh A. In sum, the update of A and the updates of one-step gradient-descent approximations mutually depend on each other. These updates are performed iteratively until convergence. Algorithm 1 shows the algorithm.

In the gradient of A calculated using chain rule, the number of chains is the same as the number of levels in our proposed multi-level optimization (MLO) formulation. This shows that this optimization algorithm preserves the multi-level nested optimization nature of the MLO formulation. In the MLO formulation, A is optimized after finishing M rounds of interleaving learning. In the optimization algorithm, A is updated iteratively, at the end of each of the M rounds. This discrepancy is due to: the optimization of A in the MLO formulation is symbolic, but the update of A in the optimization algorithm is numerical and needs to be conducted iteratively.

A.2 GENERAL ALGORITHM

Next, we describe the general algorithm that can be applied for any number of tasks and any number of rounds. The special algorithm described in the previous subsection is a special case of the general algorithm introduced in this subsection. Following (Liu et al., 2019a), we approximate $\widetilde{W}_1^{(1)}(A)$ by performing one-step gradient descent update of $W_1^{(1)}$ w.r.t $L(A,W_1^{(1)},H_1^{(1)},D_1^{(\mathrm{tr})})$:

$$\widetilde{W}_{1}^{(1)}(A) \approx \overline{W}_{1}^{(1)}(A) = W_{1}^{(1)} - \eta \nabla_{W_{1}^{(1)}} L(A, W_{1}^{(1)}, H_{1}^{(1)}, D_{1}^{(\text{tr})}). \tag{25}$$

At a middle stage, we approximate $\widetilde{W}_k^{(m)}(A)$ using a one-step gradient descent update of $W_k^{(m)}$ w.r.t. $\sum_{\mathcal{A}_i, \mathcal{A}_j} \max \left(0, -\left(d(\mathcal{A}_i, \mathcal{A}_j; \overline{W}_{k-1}^{(m)}(A)) - \tau\right) \left(d(\mathcal{A}_i, \mathcal{A}_j; W_k^{(m)}) - \tau\right)\right)$, where $\overline{W}_{k-1}^{(m)}(A)$ is

Algorithm 2 Optimization algorithm for interleaving multi-task learning

```
 \begin{tabular}{ll} \be
```

the approximation of $\widetilde{W}_{k-1}^{(m)}(A)$.

$$\begin{split} \widetilde{W}_{k}^{(m)}(A) &\approx \overline{W}_{k}^{(m)}(A) = W_{k}^{(m)}(A) \\ &- \eta \nabla_{W_{k}^{(m)}(A)} \left(\sum_{\mathcal{A}_{i}, \mathcal{A}_{j}} \max \left(0, -\left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; \overline{W}_{k-1}^{(m)}(A)) - \tau \right) \left(d(\mathcal{A}_{i}, \mathcal{A}_{j}; W_{k}^{(m)}) - \tau \right) \right) \right). \end{split}$$

$$(26)$$

Note that $\{\overline{W}_k^{(m)}(A)\}_{k,m=1}^{K,M}$ are calculated recursively, where $\overline{W}_k^{(m)}(A)$ is a function of $\overline{W}_{k-1}^{(m)}(A)$, $\overline{W}_{k-1}^{(m)}(A)$ is a function of $\overline{W}_{k-2}^{(m)}(A)$, and so on. When m>1 and k=1, $\overline{W}_{k-1}^{(m)}(A)=\overline{W}_K^{(m-1)}(A)$. For $\widetilde{H}_k^{(m)}(A)$, the approximation is:

$$\overline{H}_{k}^{(m)}(A) = H_{k}^{(m)}(A) - \eta \nabla_{H_{k}^{(m)}(A)} L(A, \overline{W}_{k}^{(m)}, H_{k}^{(m)}, D_{k}^{(\text{tr})}). \tag{27}$$

In the validation stage, we plug the approximations of $\{\widetilde{W}_k^{(M)}(A)\}_{k=1}^K$ and $\{\widetilde{H}_k^{(M)}(A)\}_{k=1}^K$ into the validation loss function, calculate the gradient of the approximated objective w.r.t the encoder architecture A, then update A via:

$$A \leftarrow A - \eta \sum_{k=1}^{K} \nabla_{A} L(A, \overline{W}_{k}^{(M)}(A), \overline{H}_{k}^{(M)}(A), D_{k}^{(\text{val})}). \tag{28}$$

We reiterate the analysis given before. After A is updated in equation 21, the one-step gradient-descent approximations (in equation 18-20), which are functions of A, need to be re-updated. The gradient update of A, which is a function of one-step gradient-descent approximations, needs to be re-calculated. In sum, the update of A and the updates of one-step gradient-descent approximations mutually depend on each other. The update steps from Eq.(25) to Eq.(28) iterate until convergence. The entire algorithm is summarized in Algorithm 2.

A.3 PSEUDO CODE

The pseudo code of our method is shown below.

While not converged

```
#Update W_1^{(1)} and H_1^{(1)}
Sample a minibatch m_1 from D_1^{(tr)}
Calculate gradient \triangle W_1^{(1)} = \nabla_{W_1^{(1)}} L_{1,1}^{(tr)}(A, W_1^{(1)}, H_1^{(1)}, m_1) Update W_1^{(1)} \leftarrow W_1^{(1)} - \eta_w \triangle W_1^{(1)} Calculate gradient: \triangle H_1^{(1)} = \nabla_{H_1^{(1)}} L_{1,1}^{(tr)}(A, W_1^{(1)}, H_1^{(1)}, m_1) Update H_1^{(1)} \leftarrow H_1^{(1)} - \eta_h \triangle H_1^{(1)}
\label{eq:potential} \begin{split} & \# \text{Update } \{W_k^{(1)}, H_k^{(1)}\}_{k=2}^K \\ & \text{For } k=2:K \end{split}
       Sample ten examples from D_{k-1}^{(\mathrm{tr})} and ten examples from D_k^{(\mathrm{tr})} For each example, construct an augmented set \mathcal{A}_i
      \begin{aligned} & \text{Calculate gradient } \triangle W_k^{(1)} = \nabla_{W_k^{(1)}} \left( \sum_{\mathcal{A}_i, \mathcal{A}_j} \max \bigg( 0, - \big( d(\mathcal{A}_i, \mathcal{A}_j; W_{k-1}^{(1)}) - \tau \big) \Big( d(\mathcal{A}_i, \mathcal{A}_j; W_k^{(1)}) - \tau \big) \right) \bigg) \\ & \text{Update } W_k^{(1)} \leftarrow W_k^{(1)} - \eta_w \bigtriangleup W_k^{(1)} \end{aligned}
Calculate gradient \bigtriangleup H_k^{(1)} = \nabla_{H_k^{(1)}} L(A, W_k^{(1)}, H_k^{(1)}, m_k) Update H_k^{(1)} \leftarrow H_k^{(1)} - \eta_h \bigtriangleup H_k^{(1)} End
\#\mbox{Update }\{W_k^{(m)},H_k^{(m)}\}_{m=2,k=1}^{M,K} For m=2:M,\;k=1:K
       Sample ten examples from D_{k-1}^{(\mathrm{tr})} and ten examples from D_{k}^{(\mathrm{tr})}
       For each example, construct an augmented set A_i
       Calculate gradient \triangle W_k^{(m)} = \nabla_{W_k^{(m)}} \left( \sum_{\mathcal{A}_i, \mathcal{A}_j} \max \left( 0, -\left( d(\mathcal{A}_i, \mathcal{A}_j; W_{k-1}^{(m)}) - \tau \right) \left( d(\mathcal{A}_i, \mathcal{A}_j; W_k^{(m)}) - \tau \right) \right) \right)
       Update W_k^{(m)} \leftarrow W_k^{(m)} - \eta_w \triangle W_k^{(m)}
       Sample a minibatch m_k from D_k^{(tr)}
      Calculate gradient \triangle H_k^{(m)} = \nabla_{H_k^{(m)}}^{\kappa} L(A, W_k^{(m)}, H_k^{(m)}, D_k^{(\text{tr})})

Update H_k^{(m)} \leftarrow H_k^{(m)} - \eta_h \triangle H_k^{(m)}
 End
 #Update A
 For k = 1 : K
       Sample a minibatch m_k from D_k^{(tr)}
Calculate gradient \triangle A = \sum_{k=1}^K \nabla_A L(A, W_k^{(M)}, H_k^{(M)}, m_k) Update A \leftarrow A - \eta_a \triangle A
```

End

B Additional ablation study results

We compare the search and memory costs of our hypernetwork (HN) based approaches with those of non-hypernetwork (NHN) approaches that store $W_k^{(m)}$ in memory. Table 5 shows the results. As can be seen, hypernetwork-based approaches have much lower search and memory costs than

Table 5: Comparison between hypernetwork (HN) based approaches and non-hypernetwork (NHN) approaches.

	Param. (M)	Search Cost (GPU days)	Memory Cost
Pcdarts (Xu et al., 2020)	4.3	4.0	1x
NHN-Pcdarts (ours)	4.0	8.5	2.4x
HN-Pcdarts (ours)	4.1	4.0	1.1x

Table 6: Test errors on CIFAR-100, CIFAR-10, and ImageNet under different values of C.

С	0	3	10	20	30	40
CIFAR-100 Error	19.96	17.29	16.51	16.08	16.06	16.09
CIFAR-10 Error	2.81	2.56	2.38	2.31	2.32	2.31
ImageNet Top-1 Error	26.2	24.8	23.6	23.1	23.1	23.2
ImageNet Top-5 Error	9.7	8.2	6.9	6.3	6.4	6.2

non-hypernetwork approaches while the test errors of hypernetwork-based approaches are similar to those of non-hypernetwork approaches as shown in Table 2.

We also investigate how the number C of augmented examples during knowledge transfer affects model performance. Table 6 shows a moderate C yields the lowest overall errors. As can be seen, when C is small, the performance is not very good. This is because under a small C, a smaller number of augmented examples cannot represent a distribution very well. When C is more than 20, the performance does not improve significantly, but the computational cost will increase. So in our experiments, we set C to 20. For C=0, no knowledge is transferred, which leads to worse performance.

We perform a study which increases the number of interleaving rounds to 20. Meanwhile, to keep the computational cost of this ablation study the same as that of the study (referred to as main study) corresponding to Table 2 where the number of rounds is 2 and the epoch number is 50, we reduce the number of epochs in this ablation study to 5. Table 7 shows the results. As can be seen, increasing the number of rounds to 20 leads to worse performance. The reason is: 1) epoch number is too small; 2) it is not very necessary to perform 20 rounds of interleaving; two rounds are sufficient for enabling each task to transfer knowledge to any other task.

Finally, we investigate how the MMD threshold τ in Eq.(4) affects performance. Table 8 shows the results. As can be seen, the value of τ does not have a significant influence on the performance. τ is like the margin in a support vector machine; its absolute value does not matter much. When τ increases or decreases, the scale of MMD can increase or decrease accordingly (by scaling the norm of latent representations $z_n^{(i)}$ and $z_m^{(j)}$ up or down accordingly) to match with the scale of τ .

C POTENTIAL NEGATIVE SOCIETAL IMPACT

One potential negative societal impact of our work is: our method currently does not take interpretability of searched architectures into account, which may yield less reliable predictions and incur trustworthiness concerns in mission-critical domains such as healthcare and finance.

D ADDITIONAL RELATED WORKS

D.1 BI-LEVEL AND MULTI-LEVEL OPTIMIZATION

A bi-level optimization (BLO) problem has two nested levels of optimization problems. At the lower level, model weights are learned by minimizing a training loss. At the upper level, meta parameters are learned by minimizing a validation loss. BLO has been applied for neural architecture search (Liu et al., 2019a; Zhang et al., 2021), hyperparameter optimization (Baydin et al., 2018; Feurer et al., 2015; Franceschi et al., 2017; 2018; Lorraine et al., 2020; Maclaurin et al., 2015), reinforcement

Table 7: Test errors under different round numbers.

-	Round number	Epoch number	Cifar100 error	Cifar10 error	ImageNet Top-1/Top-5 error
Ablation study	20	5	22.39 ± 0.15	5.19 ± 0.11	27.4/9.4
Main study	2	50	16.08 ± 0.05	2.31 ± 0.03	23.1/6.3

Table 8: Test errors on CIFAR-100, CIFAR-10, and ImageNet under different values of τ .

$\overline{\tau}$	0.1	1	5	25	100
C100 Error	16.09	16.12	16.08	16.10	16.08
C10 Error	2.32	2.31	2.31	2.33	2.31
ImageNet Top-1 Error	23.2	23.1	23.1	23.3	23.2
ImageNet Top-5 Error	6.4	6.3	6.3	6.4	6.4

learning (Hong et al., 2020; Konda & Tsitsiklis, 1999; Rajeswaran et al., 2020), data valuation (Ren et al., 2020; Shu et al., 2019; Wang et al., 2020), meta learning (Finn et al., 2017; Rajeswaran et al., 2019), and label correction (Zheng et al., 2019) where meta parameters are neural architectures, hyperparameters, importance weights of data, etc. Many optimization algorithms (Couellan & Wang, 2016; Ghadimi & Wang, 2018; Grazzi et al., 2020; Ji et al., 2021; Liu et al., 2021; Yang et al., 2021) have been proposed for solving BLO problems. Multi-level optimization (MLO) (Garg et al., 2022; He et al., 2021; Raghu et al., 2021; Somayajula et al., 2022; Such et al., 2020; Xie & Du, 2022) is an extension of BLO, with multiple levels of nested optimization problems. MLO has been applied to formulate problems that involve multiple learning stages which need to be conducted end-to-end. Sato et al. (2021) proposed a gradient-based optimization algorithm for solving MLO problems with theoretical guarantees.

D.2 TRANSFER LEARNING (TL)

Our work is also related to TL (Pratt, 1993; Mihalkova et al., 2007; Niculescu-Mizil & Caruana, 2007; Pan & Yang, 2009; Luo et al., 2017; Zhuang et al., 2020), which aims to improve performance of a target task by transferring knowledge from a source task. Various transfer strategies have been proposed, based on reweighting source data to better match target distribution (Huang et al., 2006; Jiang & Zhai, 2007; Sugiyama et al., 2007; Foster et al., 2010; Moore & Lewis, 2010; Axelrod et al., 2011; Wang et al., 2017b; Ngiam et al., 2018), transforming source and target data into a common feature space (Borgwardt et al., 2006; Pan et al., 2008; 2010; Duan et al., 2012; Long et al., 2013; Wang et al., 2017a), extracting domain-invariant representations based on adversarial learning (Ganin et al., 2016; Long et al., 2017; Tzeng et al., 2017; Hoffman et al., 2018; Shen et al., 2018; Zhang et al., 2019), regularizing target model using source model (Luo et al., 2008; Duan et al., 2009; Zhuang et al., 2009; Tommasi et al., 2010; Duan et al., 2012), etc. Unlike TL which focuses on learning a target task with a source task as auxiliary, MTL methods (including ours) aim to improve all tasks and have no source/target distinction.

Our work is also related to meta learning, which learns to transfer meta knowledge to end tasks (Andrychowicz et al., 2016; Li & Malik, 2016; Ravi & Larochelle, 2016; Finn et al., 2017; Li et al., 2017; Mishra et al., 2017; Antoniou et al., 2018; Gidaris & Komodakis, 2018; Lee & Choi, 2018; Qiao et al., 2018; Hospedales et al., 2020; Lekkala & Itti, 2020; Sun et al., 2020; Yao et al., 2020). Different from these methods, our method transfers knowledge directly between end tasks.

E NOTATIONS

Table 9 shows the notations of our method.

Table 9: Notations in our method

Notation	Meaning
\overline{K}	the number of tasks
$D_k^{(\mathrm{tr})}$	the training dataset of task k
$D_k^{(\mathrm{val})}$	the validation dataset of task k
$W_k^{(m)}$	the encoder weights of task k in round m
$H_k^{(m)}$	the head weights of task k in round m
A	meta parameters
M	the number of rounds
K	the number of tasks
C	the number of augmented data examples
$L_{k,m}^{(tr)}$	training loss defined on the model of task k in round m
$L_k^{(val)}$	validation loss of task k

F ADDITIONAL RESULTS OF BASELINES

For MTL with soft weight sharing (MTL-SWS), the formulation is:

$$\begin{aligned} & \min_{A} \ \sum_{k=1}^{K} \alpha_{k} L_{k}(A, \widetilde{W}_{k}(A), \widetilde{H}_{k}(A), D_{k}^{(\text{val})}) \\ & s.t. \ \{\widetilde{W}_{k}(A), \widetilde{H}_{k}(A)\}_{k=1}^{K} = \min_{\{W_{k}, H_{k}\}_{k=1}^{K}} \ \sum_{k=1}^{K} \beta_{k} L_{k}(A, W_{k}, H_{k}, D_{k}^{(\text{tr})}) + \gamma \sum_{1 \leq k < j \leq K} \|W_{k} - W_{j}\|_{2}^{2} \end{aligned}$$

where $\{\alpha\}_{k=1}^K$, $\{\beta\}_{k=1}^K$, and γ are tradeoff parameters.

In the experiments on CIFAR-100 and CIFAR-10, tasks 1 and 2 in MTL-SWS are classification on CIFAR-100 and CIFAR-10. α_1 , α_2 , β_1 , β_2 , and γ are set to 1, 1, 1, 1, and 100 respectively.

In the experiments on ImageNet and COCO, tasks 1 and 2 in MTL-SWS are classification and detection. α_1 , α_2 , β_1 , β_2 , and γ are set to 1, 0.5, 1, 0.5, and 10 respectively.

For MTL with hard weight sharing (MTL-HWS), the formulation is:

$$\min_{A} \sum_{k=1}^{K} \alpha_{k} L(A, \widetilde{W}(A), \widetilde{H}_{k}(A), D_{k}^{(\text{val})})
s.t. \widetilde{W}(A), \{\widetilde{H}_{k}(A)\}_{k=1}^{K} = \min_{W, \{H_{k}\}_{k=1}^{K}} \sum_{k=1}^{K} \beta_{k} L(A, W, H_{k}, D_{k}^{(\text{tr})})$$
(29)

where $\{\alpha_k\}_{k=1}^K$ and $\{\beta_k\}_{k=1}^K$ are tradeoff parameters.

For MTL sharing architecture only (MTL-SAO), the formulation is:

$$\min_{A} \sum_{k=1}^{K} \alpha_{k} L(A, \widetilde{W}_{k}(A), \widetilde{H}_{k}(A), D_{k}^{(\text{val})})
s.t. \{\widetilde{W}_{k}(A), \widetilde{H}_{k}(A)\}_{k=1}^{K} = \min_{\{W_{k}, H_{k}\}_{k=1}^{K}} \sum_{k=1}^{K} \beta_{k} L(A, W_{k}, H_{k}, D_{k}^{(\text{tr})})$$
(30)

where $\{\alpha_k\}_{k=1}^K$ and $\{\beta_k\}_{k=1}^K$ are tradeoff parameters.

For non-MTL baselines, we mostly follow their original hyperparameter settings.

Table 10 shows results of additional baselines in the ImageNet-CoCo interleaving experiment.

Table 10: Object detection on COCO test set and classification accuracy on ImageNet test set. APS, APM, APL: AP at small, medium, large scale. AP50 and AP75 are AP with IoU thresholds of 0.5 and 0.75. Acc is accuracy.

		COCO						Imag	×+	
	Acc	AP	AP50	AP75	APS	APM	APL	Top-1	Top-5	(M)
MobileNetV2 (Sandler et al., 2018)	72.0	28.3	46.7	29.3	14.8	30.7	38.1	72.0	91.0	300
FairNAS-A (Chu et al., 2019)	77.5	32.4	52.4	33.9	17.2	36.3	43.2	77.5	93.7	392
FairNAS-B (Chu et al., 2019)	77.2	31.7	51.5	33.0	17.0	35.2	42.5	77.2	93.5	349
IMTNAS (ours)	77.8	32.7	53.3	34.1	17.2	37.1	43.8	77.8	94.0	394

G DEFINITION OF MMD

Let $k(\cdot, \cdot)$ be a kernel function. Given two distributions p and q, their MMD is defined as:

$$\mathbb{E}_{x \sim p, x' \sim p}[k(x, x')] + \mathbb{E}_{y \sim q, y' \sim q}[k(y, y')] - 2\mathbb{E}_{x \sim p, y \sim q}[k(x, y)]. \tag{31}$$

Given a set of sample $\{x_i\}_{i=1}^m$ drawn from p and a set of sample $\{y_i\}_{i=1}^n$ drawn from q, the empirical MMD is calculated as:

$$\frac{1}{m(m-1)} \sum_{i=1}^{m} \sum_{j \neq i}^{m} k(x_i, x_j) + \frac{1}{n(n-1)} \sum_{i=1}^{n} \sum_{j \neq i}^{n} k(y_i, y_j) - \frac{2}{mn} \sum_{i=1}^{m} \sum_{j=1}^{n} k(x_i, y_j).$$
(32)

H EXPERIMENTAL DETAILS OF NEURAL ARCHITECTURE SEARCH

For methods based on PC-DARTS, including IMTNAS-pcdarts (ours), SI-pcdarts, MI-pcdarts, MTL-SWS-pcdarts, MTL-HWS-pcdarts, MTL-SAO-pcdarts, the experimental settings are similar. The search space of PC-DARTS follows that of DARTS. In the search spaces of PCDARTS, the candidate operations include: 3×3 and 5×5 separable convolutions, 3×3 and 5×5 dilated separable convolutions, 3×3 max pooling, 3×3 average pooling, identity, and zero. The stride of all operations is set to 1. The convolved feature maps are padded to preserve their spatial resolution. The order for convolutional operations is ReLU-Conv-BN. Each separable convolution is applied twice. The output node is the depthwise concatenation of all intermediate nodes, excluding the input nodes. We create a network by stacking 8 cells. The first and second nodes of cell k are equal to the outputs of cell k-2 and cell k-1, respectively. 1×1 convolutions are inserted when necessary. Reduction cells are located at the 1/3 and 2/3 of the total depth of the network. In reduction cells, operations adjacent to the input nodes have a stride of 2.

For architecture search on CIFAR-100 and CIFAR-10, the hyperparameter K was set to 4. The network is a stack of 8 cells. Each cell contains 6 nodes. Initial channel number is set to 16. The architecture variables are trained using the Adam optimizer. The learning rate is set to 6e-4, without decay. The weight decay is set to 1e-3. The momentum is set to (0.5, 0.999). The network weight parameters are trained using SGD. The initial learning rate is set to 0.1. Cosine scheduling is used to decay the learning rate, down to 0 without restart. The momentum is set to 0.9. The weight decay is set to 3e-4. The batch size is set to 256. Warm-up is utilized: in the first 15 epochs, architecture variables are frozen and only network weights are optimized.

The settings for architecture evaluation on CIFAR-100 and CIFAR-10 follow those of DARTS. 18 normal cells and 2 reduction cells are stacked into a large network. The initial channel number is set to 36. The stacked network is trained from scratch using SGD for 600 epochs, with batch size 128, initial learning rate 0.025, momentum 0.9, weight decay 3e-4, norm gradient clipping 5, drop-path rate 0.3, and cutout. The learning rate is decayed to 0 using cosine scheduling without restart.

We combine our method and PC-DARTS to directly search for architectures on ImageNet. The stacked network starts with three convolution layers which reduce the input image resolution from 224×224 to 28×28 , using stride 2. After the three convolution layers, 6 normal cells and 2 reduction cells are stacked. Each cell consists of N=6 nodes. The sub-sampling rate was set to 0.5. Architecture variables are trained using Adam. The learning rate is fixed to 6e-3. The weight decay is set to 1e-3. The momentum is set to 1e-3. The momentum is set to 1e-3. The initial learning rate is set to 1e-3. It is decayed to 0 using cosine scheduling without restart. Momentum is set to 1e-3. Weight decay is set to 1e-3. The batch-size is set to 1e-3. Epoch number is set to 1e-3. Eight Tesla V100 GPUs were used.

For architecture evaluation on ImageNet, the stacked network starts with three convolution layers which reduce the input image resolution from 224×224 to 28×28 , using stride 2. After the three convolution layers, 12 normal cells and 2 reduction cells are stacked. Initial channel number is set to 48. The network is trained from scratch using SGD for 250 epochs, with batch size 1024, initial learning rate 0.5, weight decay 3e-5, and momentum 0.9. For the first 5 epochs, learning rate warm-up is used. The learning rate is linearly decayed to 0. Label smoothing and auxiliary loss tower is used.

H.1 HYPERPARAMETER TUNING STRATEGY

To tune the interleaving round number K, we randomly sample 2.5K data from the 25K training set and sample 2.5K data from the 25K validation set. Then we use the 5K sampled data as a hyperparameter tuning set. K is tuned in $\{2,3,4\}$. For each configuration of K, we use the remaining 22.5K training data and 22.5K validation data to perform architecture search and use their combination to perform architecture evaluation (retraining a larger stacked network from scratch). Then we measure the performance of the stacked network on the 5K sampled data. The value of K yielding the best performance on the 5K sampled data is selected. For other hyperparameters, they mostly follow those in PC-DARTS (Xu et al., 2020).

H.2 IMPLEMENTATION DETAILS

We use PyTorch to implement all models. The version of Torch is 1.4.0 (or above). We build our method upon the official python package for PC-DARTS¹.

I SIGNIFICANCE TEST RESULTS

To check whether the performance of our proposed methods are significantly better than baselines, we perform a statistical significance test using a double-sided T-test. We use the function in the python package "scipy.stats.ttest_1samp" and report the average results over 10 different runs. Table 11 and 12 show the results.

Table 11: Significance test results on CIFAR-100

Our method	Baseline	p-value
IMTNAS-Pcdarts	SI-Pcdarts	7.02e-4
IMTNAS-Pcdarts	MI-Pcdarts	2.59e-5
IMTNAS-Pcdarts	MTL-SWS-Pcdarts	9.33e-4
IMTNAS-Pcdarts	MTL-HWS-Pcdarts	3.37e-5
IMTNAS-Pcdarts	MTL-SAO-Pcdarts	1.71e-5
IMTNAS-Pcdarts	Pcdarts	4.08e-4

Table 12: Significance test results on CIFAR-10

Our method	Baseline	p-value
IMTNAS-Pcdarts	SI-Pcdarts	5.96e-7
IMTNAS-Pcdarts	MI-Pcdarts	3.19e-7
IMTNAS-Pcdarts	MTL-SWS-Pcdarts	4.62e-6
IMTNAS-Pcdarts	MTL-HWS-Pcdarts	3.05e-7
IMTNAS-Pcdarts	MTL-SAO-Pcdarts	8.93e-7
IMTNAS-Pcdarts	Pcdarts	7.31e-6

From these two tables, we can see that the p-values are small between baselines methods and our methods, which demonstrate that the errors of our methods are significantly lower than those of baselines.

J FULL LISTS OF HYPERPARAMETER SETTINGS

Table 13, 14, and 15 show the hyperparameter settings of IMTNAS.

K VISUALIZATION OF SEARCHED ARCHITECTURES

Figure 3 visualizes the normal cell and reduction cell searched by our method.

¹https://github.com/yuhuixu1993/PC-DARTS/

Table 13: Hyperparameter settings in IMTNAS-PCDARTS during architecture search. For MTLSWS-PCDARTS, most hyperparameter settings are the same as those in this table.

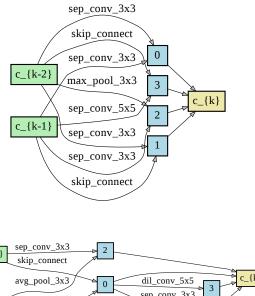
Name	Value
Optimizer for architecture variables	Adam
Learning rate for architecture variables	3e-4
Weight decay for architecture variables	0.001
Optimizer for network weights	Momentum SGD
Learning rate for network weights	0.025
Minimum learning rate for network weights	0.0
Momentum for network weights	0.9
Weight decay for network weights	3e-4
Epochs	50
Batch size	64
Number of layers	8
Number of initial channels	16
Max gradient norm	5.0

Table 14: Hyperparameter settings of IMTNAS-PCDARTS during architecture evaluation on CI-FAR10 and CIFAR100. For MTLSWS-PCDARTS, most hyperparameter settings are the same as those in this table.

Name	Value
Optimizer	Momentum SGD
Learning rate	0.025
Epochs	600
Weight decay	3e-4
Weight momentum	0.9
Auxiliary weight	0.4
Cutout length	16
Path dropout probability	0.3
Batch size	96
Number of layers	20
Number of initial channels	36
Max gradient norm	5.0

Table 15: Hyperparameter settings of architecture evaluation on ImageNet

Name	Value
Optimizer	Momentum SGD
Learning rate	0.5
Learning rate decay gamma	0.97
Epochs	250
Weight decay	3e-5
Weight momentum	0.9
Auxiliary weight	0.4
Path dropout probability	0
Label smoothing epsilon	0.1
Batch size	1024
Number of layers	14
Number of initial channels	48
Max gradient norm	5.0



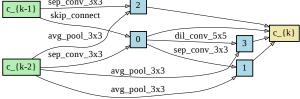


Figure 3: (Left) Normal cell searched by our method. (Right) Reduction cell searched by our method.