Low-Rank Knowledge Decomposition for Medical Foundation Models

A. Three Pre-training Datasets

- The dataset RadImageNet [12] consists of 1.35 million images, covering 11 tasks and 3 common modalities. The distribution of diseases is shown in Figure 1. In our experiments, we decompose the pre-trained models on RadImageNet [12] (which have been publicly released by the authors) into 11 lightweight expert models corresponding to task IDs.
- The dataset MedMnist is selected from MedMnistV2 [21], consisting of 705,689 images, covering 10 tasks and 7 different modalities. The distribution of diseases is shown in Figure 2. In our experiments, we decompose the fully pre-trained models on MedMnist into 10 lightweight expert models corresponding to task IDs.
- The dataset Med-ML is a multi-task dataset we constructed, consisting of 119,655 images, covering 8 tasks and 5 different modalities, including APTOS [3], ISIC [4], BUSI [7], Kvasir [15], Shenzhen X-ray [6], Shoulder X-ray [5], VinDr [14] and Bone [8]. The distribution of diseases is shown in Figure 3. In our experiments, we decompose the fully pre-trained models on Med-ML into 8 lightweight expert models corresponding to task IDs.



Figure 1. Disease distribution in Radimagenet.



Figure 2. Disease distribution in MedMnist.



Figure 3. Disease distribution in Med-MT.

B. Seven Downstream Datasets

In the experiments, the downstream datasets we used include COVID [20], BTC [17], AD [1], Mura [16], AUITD [2], HAM10000 [18], and DET10 [11]. These datasets cover five common modalities and are used to thoroughly validate the effectiveness and generalization of our method. The description of these datasets is shown in Figure 4.



Figure 4. Seven downstream datasets used in our paper.

C. Correspondence between experts and datasets

For STL-based methods, because they train models independently for each task in the pre-training dataset, the trained models can be considered as "expert models" lacking general knowledge. In this case, fine-tuning is only performed when the downstream task matches the model. In the main text, the "-" symbol is used to indicate whether there is a match.

For MTL-based methods, we fine-tune their shared encoders on all downstream datasets since MTL-based methods do not generate task-specific experts.

For KF and our method LoRKD, we fine-tune the corresponding expert models on each downstream dataset. The correspondence between expert models and downstream datasets can be seen in Table 1. The symbol \dagger indicates the absence of a corresponding expert model (due to task or modality mismatch). Following the work of [22], in such cases, we fine-tune a shared backbone that incorporates general knowledge learned from multiple tasks.

Table 1. Correspondence between expert models and downstream datasets.

Pre-trained data	COVID [20]	BTC [17]	AD [1]	Mura_s [16]	AUTID [2]	HAM10000 [18]	DET10[11]
Radimagenet MedMnist Med-MT	Expert_1 Expert_10 †	Expert_10 † †	Expert_10 † †	Expert_6 †	Expert_3 Expert_5	† Expert_7 Expert_2	Expert_1 Expert_10 †

D. Efficiency Analysis

The role of our EKS conv is to construct personalized low-rank adapters for each sample in the mini-batch, and there is not a unique way to achieve this goal. Thus, to demonstrate the efficiency advantage of EKS conv, we follow FLoRA [19] to compare the computational costs of different methods [9, 19] from a theoretical perspective (as shown in Table. 2). Following [19], we denote b and l as the batch size and the maximum sequence length in the input batch, and $W_0 \in \mathbb{R}^{d \times k}$, $B_i \in \mathbb{R}^{d \times r}$, $A_i \in \mathbb{R}^{r \times k}$. In addition, c_1 and c_2 represent the computational coefficients of batched matmuls (bmm, " φ ") and matrix multiplication (" \circ ") respectively. We also omit the cost of " \circ " and set d = k as [19], and T is the number of tasks.

	able 2. Efficiency comparison of	different methods for constructi	ng personalized low-ran	k experts for each	sample in a mini-batch.
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Method	Improved Operation	Computational Cost
LoRA [9, 19]	$ $ $\mathbf{Y} = \mathbf{X}W_0 + \varphi(\varphi(\mathbf{X}, \mathbf{B}), \mathbf{A})$	$2c_1(dblr) + c_2(bld^2)$
FLoRA [19]	$ \qquad \mathbf{Y} = \mathbf{A} \circ ((\mathbf{B} \circ \mathbf{X}) \mathbf{W}_0)$	$c_2(rbld^2)$
EKS conv (ours)	$ \mathbf{Y} = \mathbf{X}(\mathbf{W}_{0} + \sum_{i=1}^{T} (\widetilde{\mathbf{BA}} \odot \mathbf{M})_{i})$	$Tc_2(rd^2) + c_2(bld^2)$

Since both our EKS conv and FLoRA aim to construct personalized low-rank adapters for each sample in a mini-batch, but their principles for improving efficiency are different. Specifically, FLoRA replaces expensive batched matmuls (bmm) with cheap element-wise multiplications, while we perform parameter fusion before the forward pass of DNNs. If we say the efficiency of EKS conv is better than that of FLoRA, the condition in the following must be satisfied:

$$\frac{rbld^2}{Td^2r + bd^2l} \ge 1 \Longrightarrow \frac{Tr}{bl} + 1 \le r$$

Note that this inequality holds true in most real-world cases, as bl > Tr and r > 2 are common training settings. In addition, using broadcasting to improve efficiency as [19] cannot be widely generalized to convolution operations, while our method is not subject to this limitation.

E. Results on Larger foundation models

Considering that larger foundation models may be encountered in the real world, here we add LVM-Med [13] and BioMed-CLIP [23] to further validate the effectiveness of our LoRKD on dataset Med-MT.

Table 3. Comparison of larger foundation models with the best baseline on the Med-MT dataset in terms of decomposition performance.

Table I	Method	Retinal	Skin	Breast	GI tract	Lung	Shoulder	Lung	Bone	Avg
LVM-Med	best baseline	78.14	78.57	77.85	87.94	69.91	79.81	64.37	49.41	73.25
	LoRKD	79.64	82.42	78.76	88.25	75.45	82.69	64.87	53.94	75.75
BioMedCLIP	best baseline	78.14	78.57	77.85	87.94	69.91	79.81	64.37	49.41	73.25
	LoRKD	79.64	80.52	76.89	89.19	77.88	85.58	65.12	52.53	75.92

Table 4. Comparison of larger foundation models with the best baseline on the Med-MT dataset in terms of transferability.

Table II	Method	COVID	BTC	AD	Mura_s	AUITD	HAM10000	DET10	Avg
LVM-Med	best baseline	82.76	76.65	77.48	77.09	97.49	74.92	87.15	81.93
	LoRKD	84.24	79.70	77.23	74.96	97.77	77.28	87.23	82.63
BioMedCLIP	best baseline	82.76	76.65	77.48	77.09	97.49	74.92	87.15	81.93
	LoRKD	84.24	78.68	77.98	76.91	97.49	77.28	87.34	82.96

We can summarize two following points: 1) Compared with results in submission, decomposing the larger foundation models achieves the better decomposition and transferring performance. 2) Compared with results in these two tables, the superiority of our method over best baselines still holds, which confirms the advantage of our method.

F. Discussion of Comparison methods

Knowledge decomposition of foundation models to save cost during serving is relatively a new topic, especially in the medical area. The only directly correlated and available baseline is KF [22], which is proposed and verified in natural domains. Thus, we try to verify the effectiveness of LoRKD as much as possible in the following perspectives for comprehensive comparison.

Туре	Comparison Methods
Pre-training baseline	Foundation models
Single-task direct training baseline	STL
SOTA Multi-task training baselines	MTL, MoCo-MTL, Aligned-MTL
Knowledge Distillation (KD)	STL-KD, MTL-KD
Knowledge Decomposition (KDe)	KF

Table 5. Different types of comparison methods

Generally, we would like to form three following points by comparison: 1) By means of both pre-training models and small models decomposed from pre-training models, the downstream task performance should be better than directly training or multi-task collaborative training on the narrow downstream data. 2) Small models decomposed from pre-training models will maintain and even outperform the performance of pre-training models in specific tasks, due to the merits of distillation on pre-training model. 3) Naive distillation from pre-training models to a specific model is not better than the distillation from pre-training models to our mixture of Low-rank Expert modules as we consider the heterogeneity harmony and the task collaboration benefits in design.

G. Knowledge Disentanglement



Figure 5. The CKA feature similarity matrices of MTL and Ours.

Figure 5 shows the Centered Kernel Alignment (CKA) feature similarity matrices [10] of our method and MTL on Radimagenet dataset. It is evident that our method exhibits significantly lower CKA feature similarity between different tasks compared to MTL, which confirms the knowledge disentanglement ability of our method. This phenomenon can be attributed to our low-rank expert modules being embedded at the convolutional level, which facilitates the simultaneous decomposition of shallow knowledge and deep knowledge. Meanwhile, our proposed efficient knowledge decomposition convolution ensures that this knowledge decomposition pattern can be achieved at a low cost.

H. Notation table

Notation	Description	Shape
W_0	Shared weight in backbone	$\mathbb{R}^{C^{\text{out}} \times C^{\text{in}} \times k \times k}$
B_t	Low rank factors	$\mathbb{R}^{C^{\text{out}}k \times rk}$
A_t	Low rank factors	$\mathbb{R}^{rk \times C^{\text{in}}k}$
h_t	Input features	$ \mathbb{R}^{B \times C^{\text{in}} \times H \times W} $
g_t	Output features	$\mathbb{R}^{B \times C^{\text{out}} \times H \times W}$
o_{ij}	Output feature unit	$\mathbb{R}^{B \times C^{\text{out}}}$
$h_{(i)(j)}$	Input feature unit	$\mathbb{R}^{B \times C^{\text{in}}}$
ω	Convolution weight unit	$\mathbb{R}^{C^{\text{in}} \times C^{\text{out}}}$
М	Task label	$\mathbb{R}^{B \times T}$
W'	Aggregated weight	$ \mathbb{R}^{B \times C^{\text{out}} \times C^{\text{in}} \times k \times k} $

We add a notation table here to ease reading which is summarized as below.

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