

Maestro: Uncovering Low-Rank Structures via Trainable Decomposition

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Efficient Deep Learning

- One of the **key challenges** of deploying deep learning models in practice is **the increasing costs of large-scale model training and deployment**.
- Popular remedy: **compress the network**.
- Techniques: i) quantization, ii) pruning, iii) **low-rank approximation**.
- Main goals:** i) obtain a lower footprint model, ii) avoid the overhead during training time and the accuracy degradation, iii) avoid introducing multiple hyperparameters and the need to fine-tune to recover the lost accuracy.

Trainable Decomposition

Low-rank approximation.

Best l -rank approximation:

$$\min_{U \in \mathbb{R}^{m \times l}, V \in \mathbb{R}^{n \times l}} \|\sum_{i=1}^l u_i v_i^\top - A\|_F^2 \quad (1)$$

The best rank approximation across all the ranks:

$$\min_{U \in \mathbb{R}^{m \times r}, V \in \mathbb{R}^{n \times r}} \frac{1}{r} \sum_{b=1}^r \|U_{:b} V_{:b}^\top - A\|_F^2, \quad (2)$$

Data-dependent low-rank approximation.

$$\min_{U \in \mathbb{R}^{m \times r}, V \in \mathbb{R}^{n \times r}} \mathbf{E}_{x,y \sim \mathcal{X}} \left[\sum_{b=1}^r \frac{1}{r} \|U_{:b} V_{:b}^\top x - y\|^2 \right]. \quad (3)$$

This formulation is the first building block of **Maestro**, where we propose to use **double sampling** to solve this objective, by sampling **data**, i.e., $x, y \sim \mathcal{D}$, and **rank**, i.e., $b \sim \mathcal{D}(\{1, 2, \dots, r\})$.

DNN low-rank approximation. We seek to uncover the optimal ranks for a set of d linear mappings $W^1 \in \mathbb{R}^{m_1 \times n_1}, \dots, W^d \in \mathbb{R}^{m_d \times n_d}$, where W^i 's are model parameters, e.g., **linear layer weights** (also transformers or convolutions), by decomposing them as $W^i = U^i (V^i)^\top$. To adapt **double sampling** and preserve structure, we sample **one layer and its rank** in each step.

Rank extraction via hierarchical group-lasso.

$$\lambda_{gl} \sum_{i=1}^d \sum_{b=1}^{r_i} (\|U_{:b}^i\| + \|V_{:b}^i\|), \quad (4)$$

This penalty encourages that **unimportant ranks become zero** and **can be effectively removed from the model**. For each layer we remove $V_{:b}^i$ and $U_{:b}^i$ if $\|V_{:b}^i\| \|U_{:b}^i\| \leq \epsilon_{ps}$ (numerical zero, e.g., $\epsilon_{ps} = 1e-5$).

Maestro as SVD

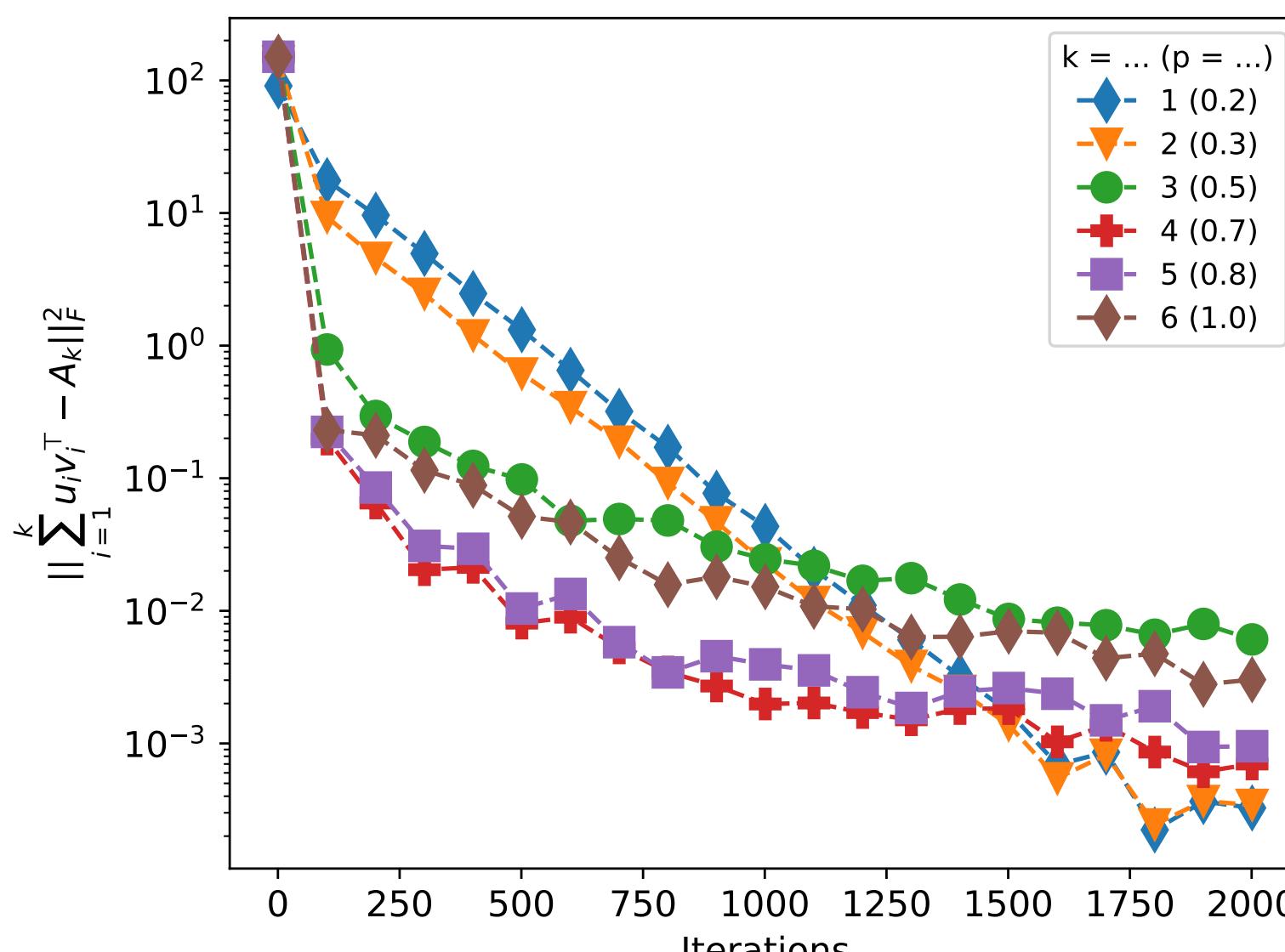


Figure 1: L2 distance between the best rank k and MAESTRO's approximation of mapping A .

If data are uniform and $y = Ax$, then
Maestro recovers SVD of A .

Maestro as PCA

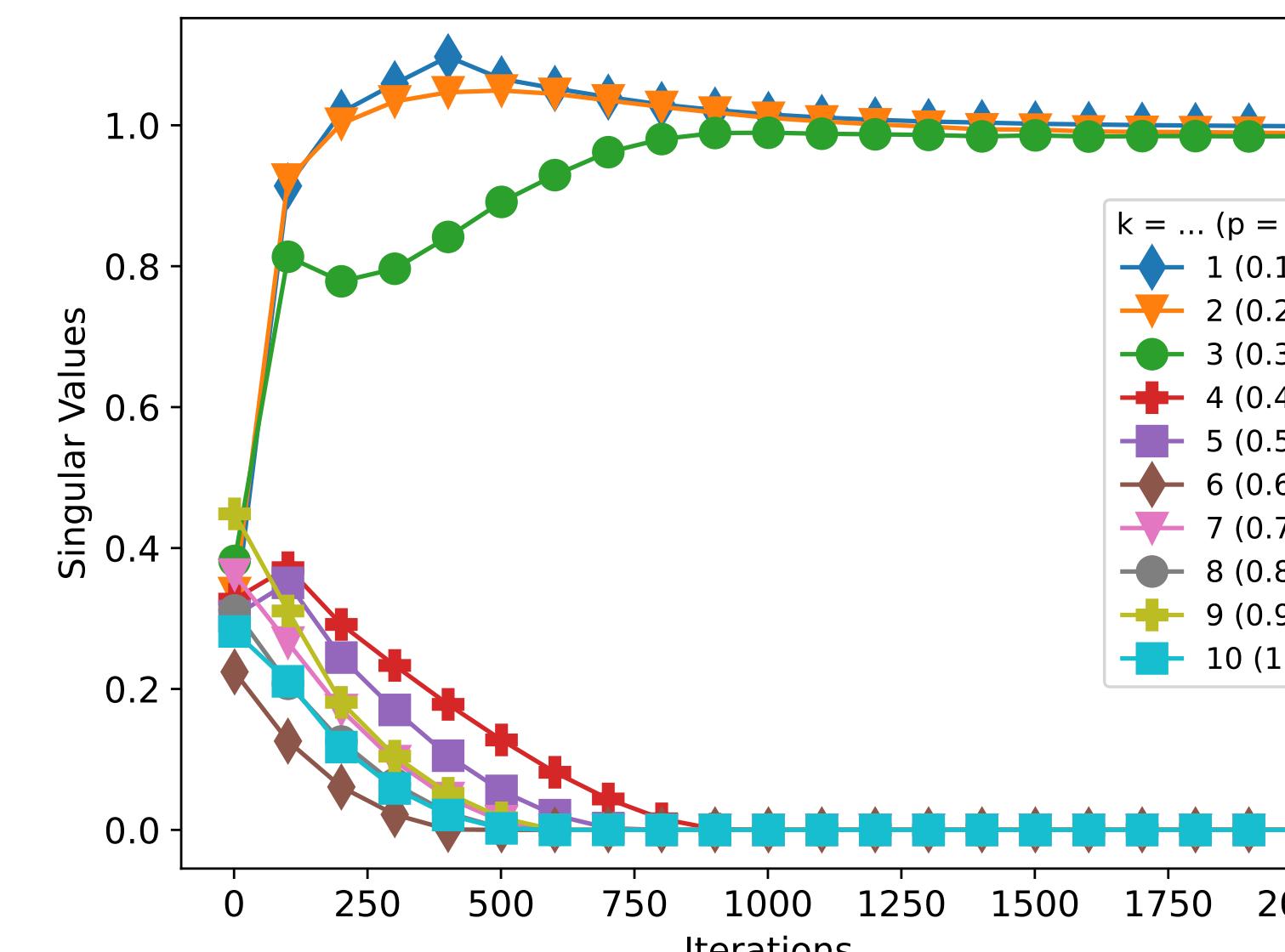


Figure 2: The plot displays the estimates of singular values.

If $y = Ix = x$, then
Maestro recovers PCA.

Experiments (ResNet18 on CIFAR10)

Table 1: MAESTRO vs. baselines

Variant	Acc. (%)	GMACs	Params. (M)
Non-factorized	93.86 ± 0.20	0.56	11.17
Pufferfish	94.17	0.22	3.336
Cuttlefish	93.47	0.3	3.108
MAESTRO [†] ($\lambda_{gp} = 16e-6$)	94.19 ± 0.07	0.39 ± 0.00	4.08 ± 0.02
MAESTRO [†] ($\lambda_{gp} = 64e-6$)	93.86 ± 0.11	0.15 ± 0.00	1.23 ± 0.00

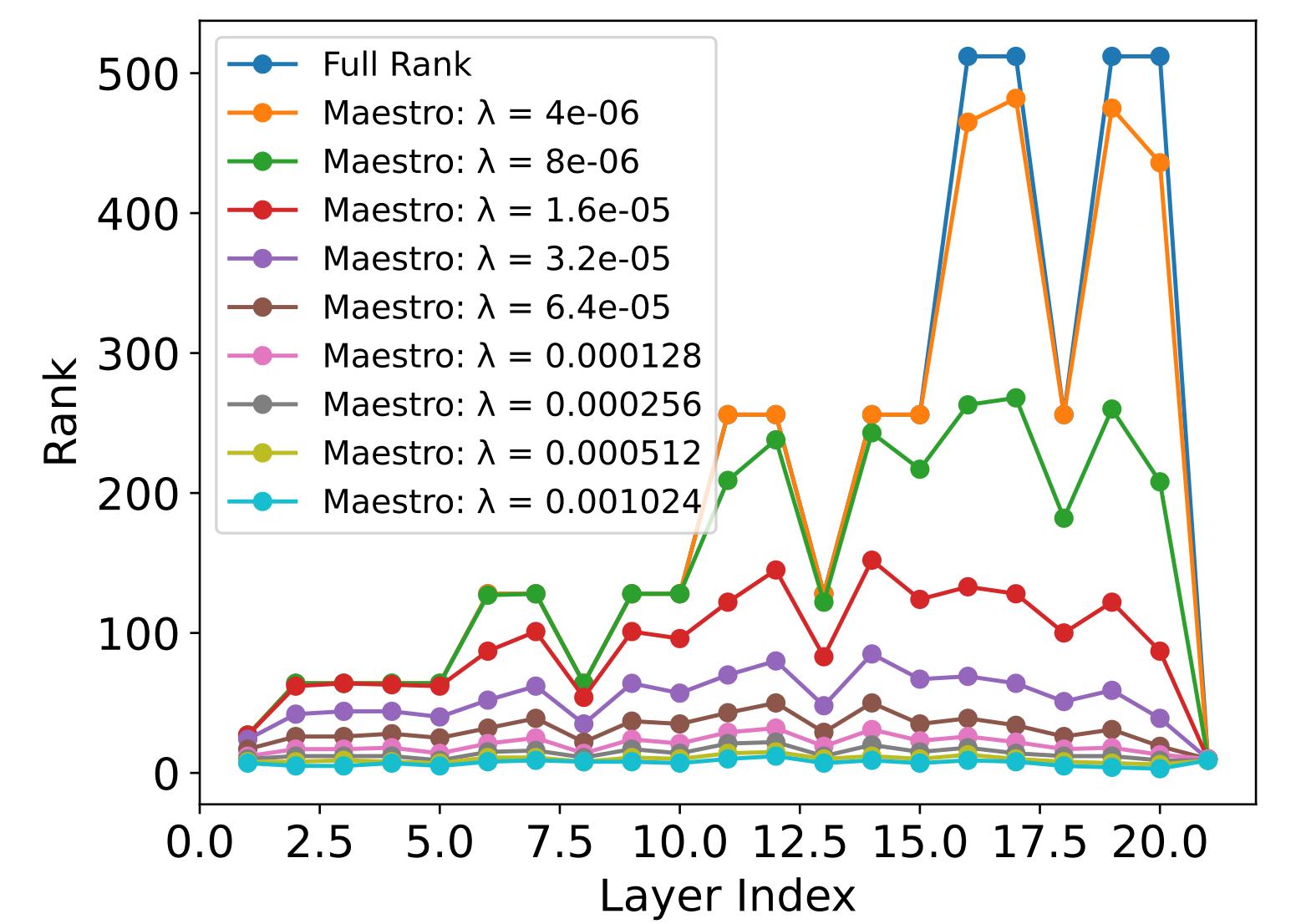
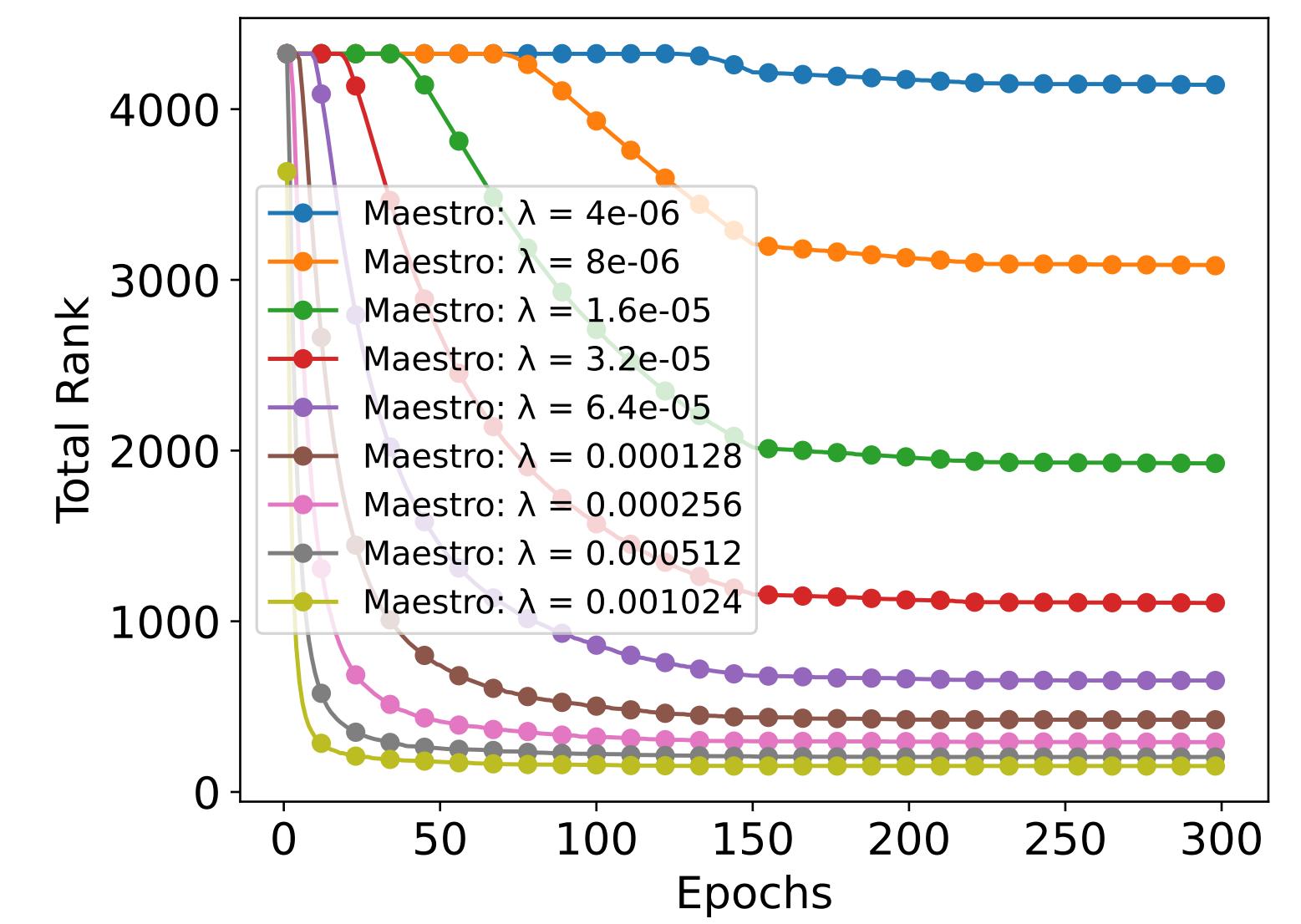


Figure 3: Training dynamics of MAESTRO. **Upper:** Total rank ($\sum_{i=1}^d r_i$). **Bottom:** Ranks r_i 's after training.

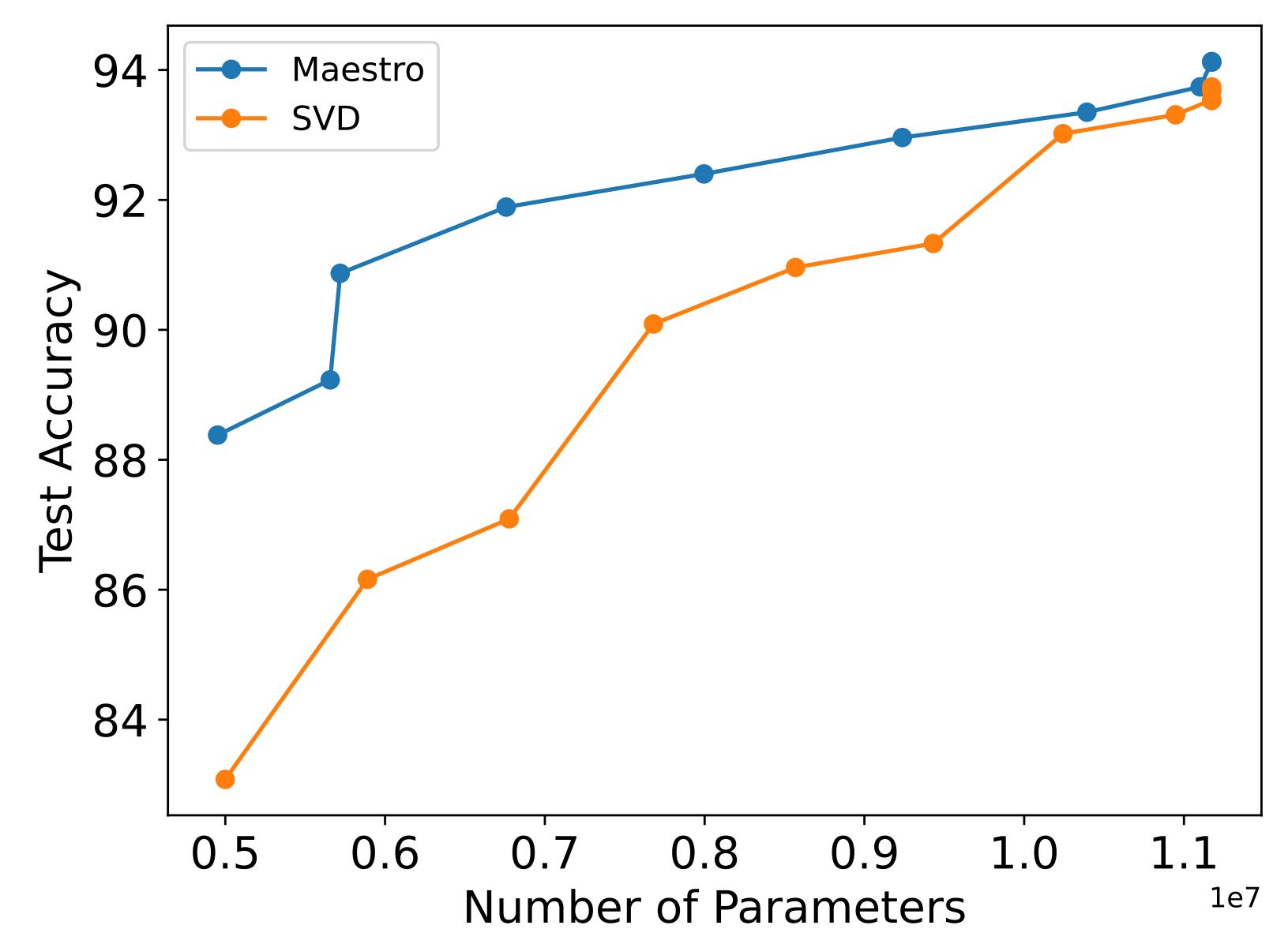
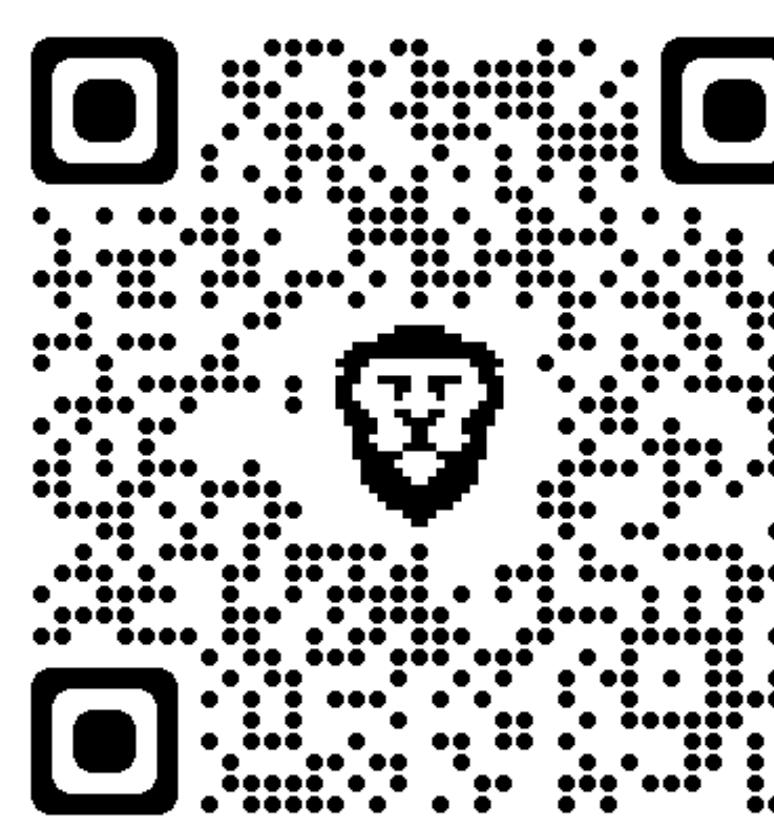


Figure 4: MAESTRO compared to SVD.

Results with extra models (VGG, Transformers) and datasets (Multi30k, TinyImageNet) in the paper.



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