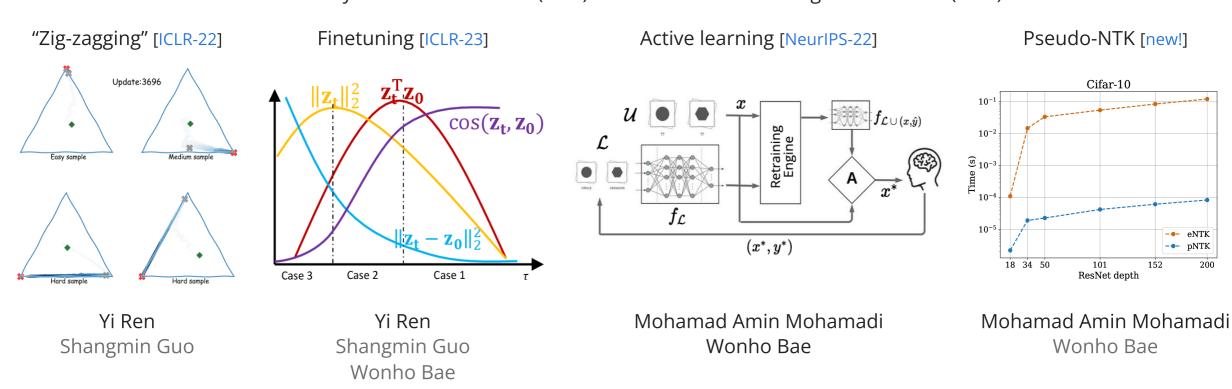
# In Defence of (Empirical) Neural Tangent Kernels

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•  $\ell'_y(\hat{y}) = \hat{y} - y$  for square loss,  $\hat{y}_y - \log \sum_{j=1}^k \exp(\hat{y}_j)$  for cross-entropy

• Full-batch GD:

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• Observation II: As f becomes "infinitely wide" with any usual architecture+init\* [Yang 2019], eNTK<sub>w0</sub>  $(x_1, x_2) \longrightarrow NTK(x_1, x_2)$ , independent of the random w<sub>0</sub>

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  - Good results in statistical testing [Jia+ 2021], dataset distillation [Nguyen+ 2021], clustering for active learning batch queries [Holzmüller+ 2022], ...

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    - Internal activations in the networks don't change much [Chizat+ 2019] [Yang/Hu 2021]

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• We now know many problems where gradient descent on an NN  $\gg$  *any* kernel method  $\circ$  Cases where GD error  $\rightarrow 0$ , any kernel is *barely* better than random [Malach+ 2021]

### What can we learn from empirical NTKs?

#### In this talk:

- As a theoretical-ish tool for local understanding:
  - Fine-grained explanation for early stopping in knowledge distillation
  - How you should fine-tune models
- As a practical tool for approximating "lookahead" in active learning
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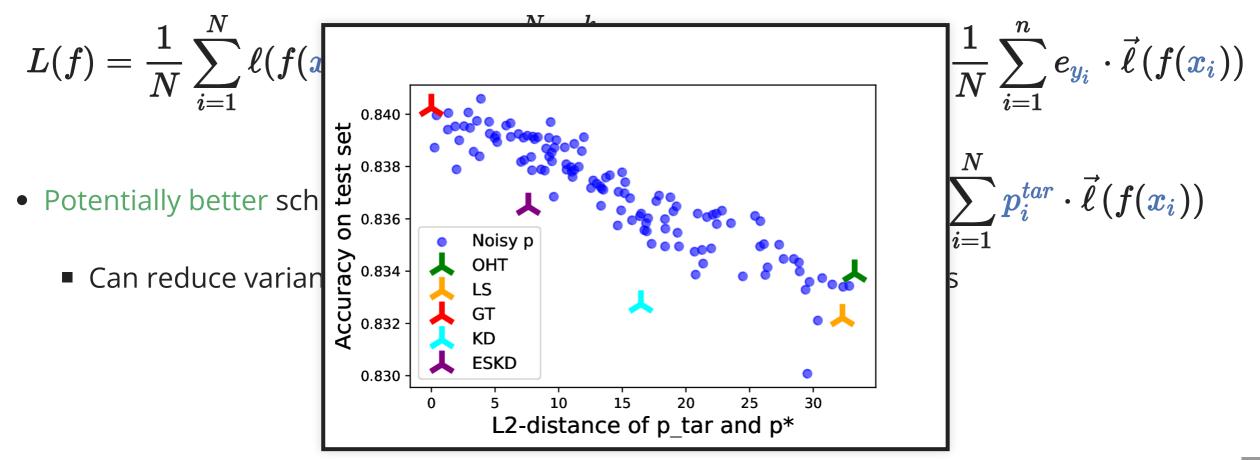
• Potentially better scheme: see  $\{(x_i, p_i^{tar})\}$ , minimize  $L^{tar}(f) = rac{1}{N} \sum_{i=1}^N p_i^{tar} \cdot \vec{\ell}(f(x_i))$ 

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  - Train a teacher  $f^{teacher}$  on  $\{(x_i, y_i)\}$  with standard ERM, L(f)
  - Train a student on  $\{(x_i, f^{teacher}(x_i))\}$  with  $L^{tar}$
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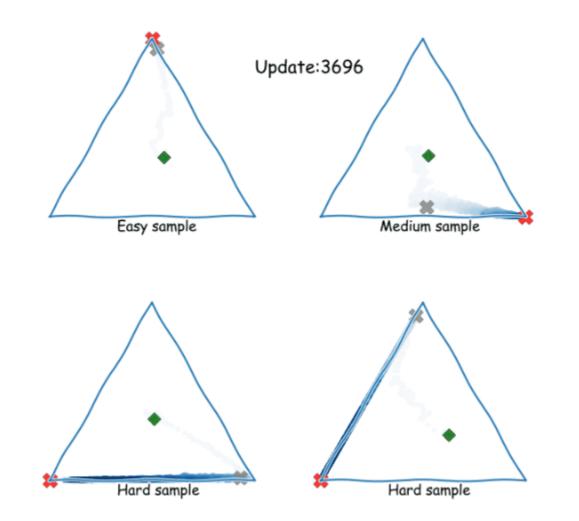
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- One possible explanation:  $f^{teacher}(x_i)$  is closer to  $p_i^*$  than sampled  $y_i$

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- One possible explanation:  $f^{teacher}(x_i)$  is closer to  $p_i^*$  than sampled  $y_i$
- But why would that be?

## Zig-Zagging behaviour in learning



Plots of (three-way) probabilistic predictions: imes shows  $p_i^*$  , imes shows  $y_i$ 

## eNTK explains it

• Let  $q_t( ilde{x}) = ext{softmax}(f_t( ilde{x})) \in \mathbb{R}^k$ ; for cross-entropy loss, one SGD step gives us

$$q_{t+1}( ilde{x}) - q_t( ilde{x}) = \eta \ \mathcal{A}_t( ilde{x}) \ ext{eNTK}_{\mathbf{w}_t}( ilde{x}, x_i) \left( p_i^{tar} - q_t(x_i) 
ight) + \mathcal{O}(\eta^2)$$

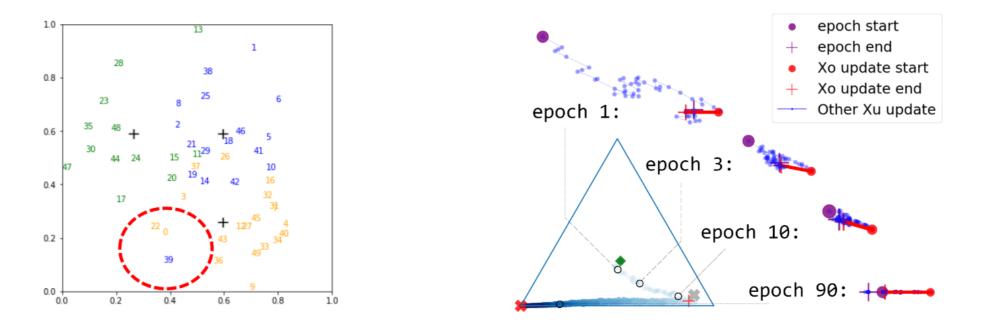
 $\mathcal{A}_t(\tilde{\pmb{x}}) = ext{diag}(q_t(\tilde{\pmb{x}})) - q_t(\tilde{\pmb{x}})q_t(\tilde{\pmb{x}})^{\mathsf{T}}$  is the covariance of a  $ext{Categorical}(q_t(\tilde{\pmb{x}}))$ 

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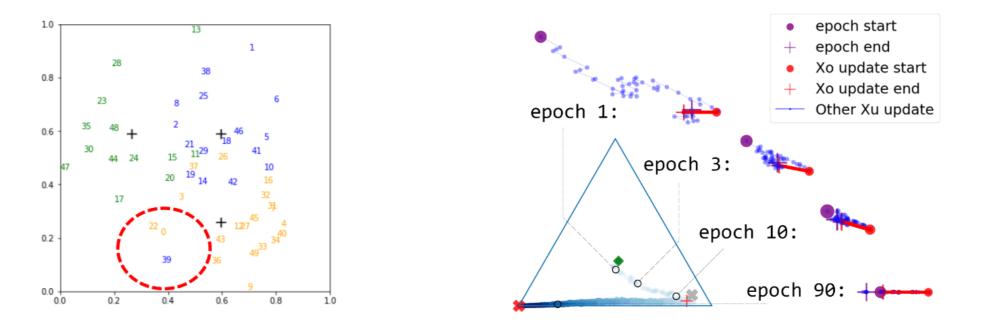


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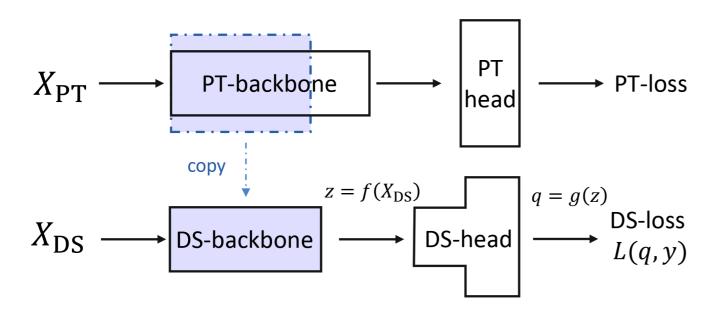


• Improves distillation (esp. with noisy labels) to take moving average of  $q_t(x_i)$  as  $p_i^{tar}$ 

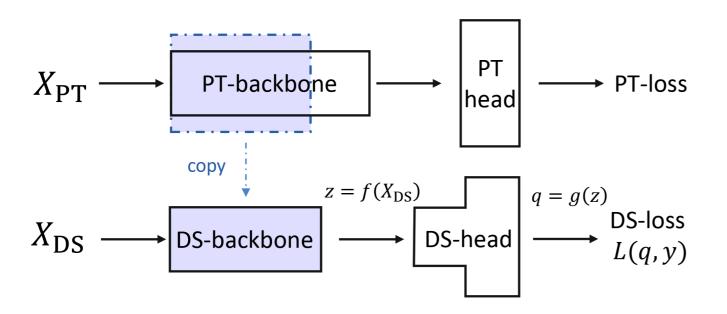
## What can we learn from empirical NTKs?

#### In this talk:

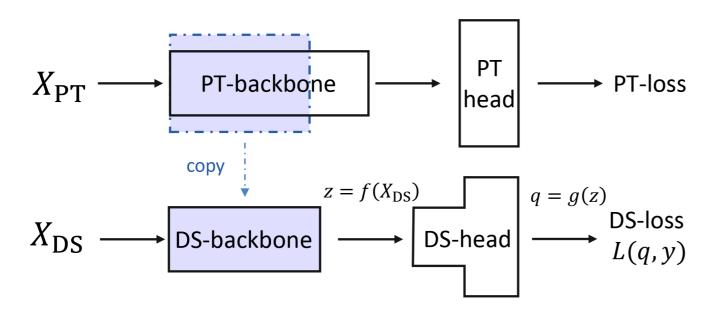
- As a theoretical-ish tool for local understanding:
  - Fine-grained explanation for early stopping in knowledge distillation
  - How you should fine-tune models
- As a practical tool for approximating "lookahead" in active learning
- Plus: efficiently approximating  ${
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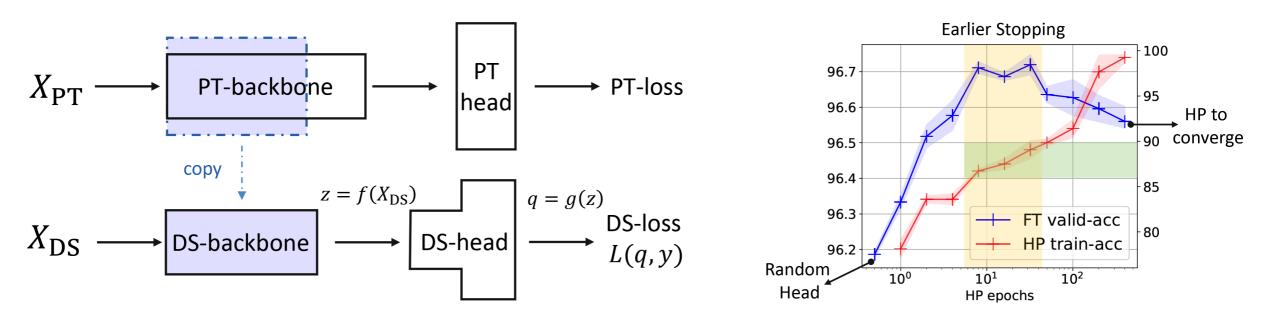
- Pretrain, re-initialize a random head, then adapt to a downstream task. Two phases:
  - Head probing: only update the head g(z)
  - Fine-tuning: update head g(z) and backbone z = f(x) together



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• Same kind of decomposition with backbone features z = f(x), head  $q = \operatorname{softmax}(g(z))$ :

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- Specializing to simple linear-linear model, can get insights about trends in  $oldsymbol{z}$
- Recommendations from paper:
  - Early stop during head probing (ideally, try multiple lengths for downstream task)
  - Label smoothing can help; so can more complex heads, but be careful

### How good will our fine-tuned features be? [Wei/Hu/Steinhardt 2022]

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• Rank-one updates for efficient computation: schema  $\Box + \Box \Box \times \Box = \begin{pmatrix} -1 \\ - \end{bmatrix} \end{pmatrix}$ 

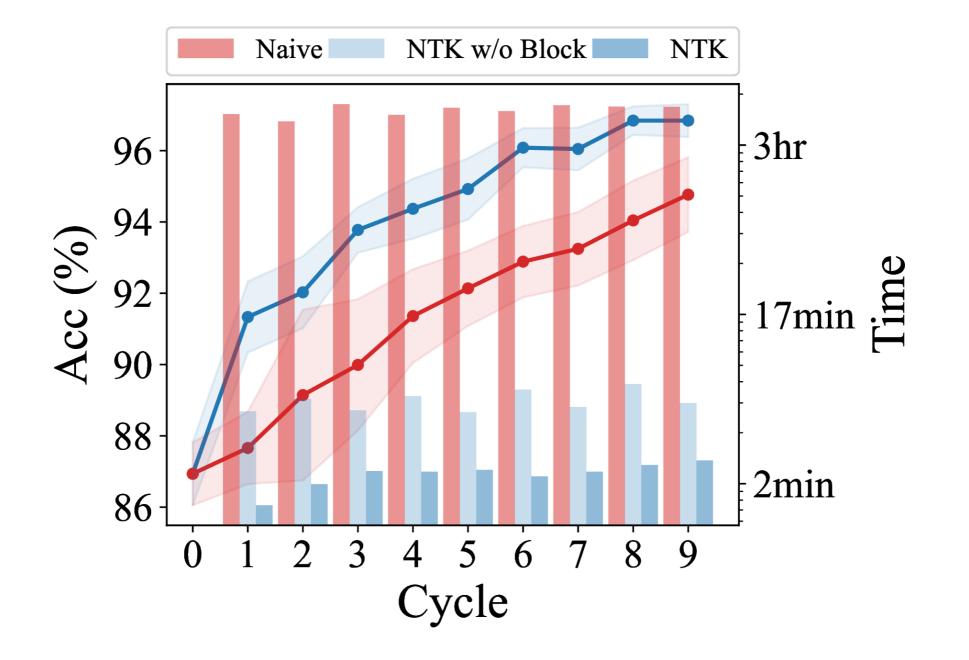
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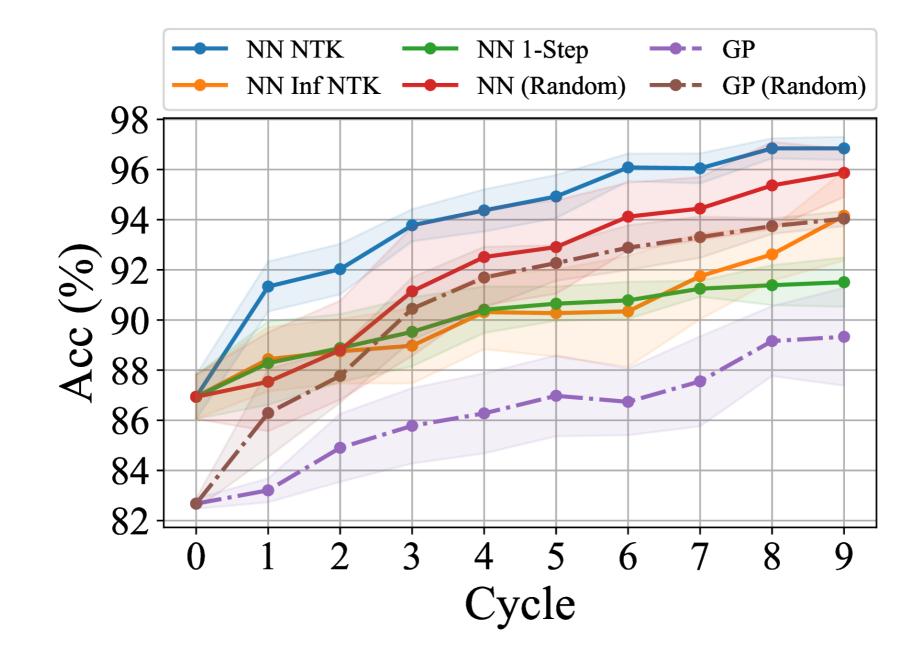
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- Local approximation with eNTK "should" work much more broadly than "NTK regime"

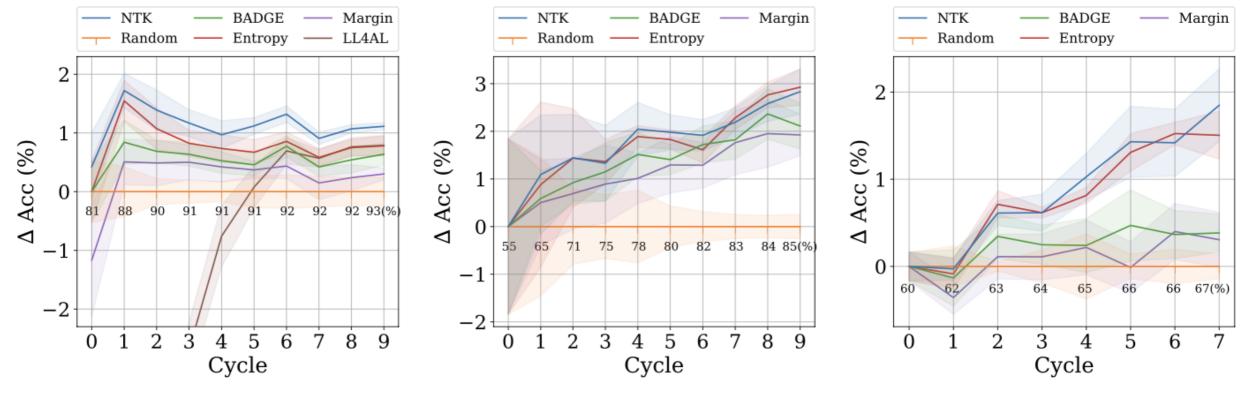
### **Much faster than SGD**



#### Much more effective than infinite NTK and one-step SGD



## Matches/beats state of the art



(a) SVHN: 1-layer WideResNet (b) CIFAR10: 2-layer WideResNet

(c) CIFAR100: ResNet18

Figure 2: Comparison of the-state-of-the-art active learning methods on various benchmark datasets. Vertical axis shows difference from random acquisition, whose accuracy is shown in text.

Downside: usually more computationally expensive (especially memory)

## **Enables new interaction modes**

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- Lots of work (including above) has used  $\mathbf{pNTK}$  instead of  $\mathbf{eNTK}$ 
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- Can we justify this more rigorously?

• Say  $f(x)=V\phi(x)$ ,  $\phi(x)\in \mathbb{R}^h$ , and  $V\in \mathbb{R}^{k imes h}$  has rows  $v_j\in \mathbb{R}^h$  with iid entries

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  - Want  $v_1^\mathsf{T} A v_1$  and  $v_j^\mathsf{T} A v_j$  to be close, and  $v_j^\mathsf{T} A v_{j'}$  small, for random v and fixed A

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  - Want  $v_1^\mathsf{T} A v_1$  and  $v_j^\mathsf{T} A v_j$  to be close, and  $v_j^\mathsf{T} A v_{j'}$  small, for random v and fixed A

• Using Hanson-Wright: 
$$\frac{\left\| \text{eNTK} - \text{pNTK} I \right\|_{F}}{\left\| \text{eNTK} \right\|_{F}} \leq \frac{\left\| \text{eNTK}^{\phi} \right\|_{F} + 4\sqrt{h}}{\text{Tr}(\text{eNTK}^{\phi})} k \log \frac{2k^{2}}{\delta}$$

- Say  $f(x)=V\phi(x)$ ,  $\phi(x)\in \mathbb{R}^h$ , and  $V\in \mathbb{R}^{k imes h}$  has rows  $v_j\in \mathbb{R}^h$  with iid entries
- If  $v_{j,i} \sim \mathcal{N}(0,\sigma^2)$ , then  $v_1$  and  $rac{1}{\sqrt{k}}\sum_{j=1}^k v_j$  have same distribution

$$egin{aligned} \mathrm{eNTK}_{\mathbf{w}}(x_1,x_2)_{jj'} &= v_j^\mathsf{T} \ \mathrm{eNTK}_{\mathbf{w} ackslash V}^\phi(x_1,x_2) \ v_{j'} + \mathbb{I}(j=j') \phi(x_1)^\mathsf{T} \phi(x_2) \ \mathbf{v}_1 \mathbf{w}_{\mathbf{w} ackslash V}(x_1,x_2) &= v_1^\mathsf{T} \ \mathrm{eNTK}_{\mathbf{w} ackslash V}^\phi(x_1,x_2) \ v_1 + \phi(x_1)^\mathsf{T} \phi(x_2) \end{aligned}$$

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• Fully-connected ReLU nets at init., fan-in mode: numerator  $\mathcal{O}(h\sqrt{h})$ , denom  $\Theta(h^2)$ 

## **pNTK's Frobenius error**

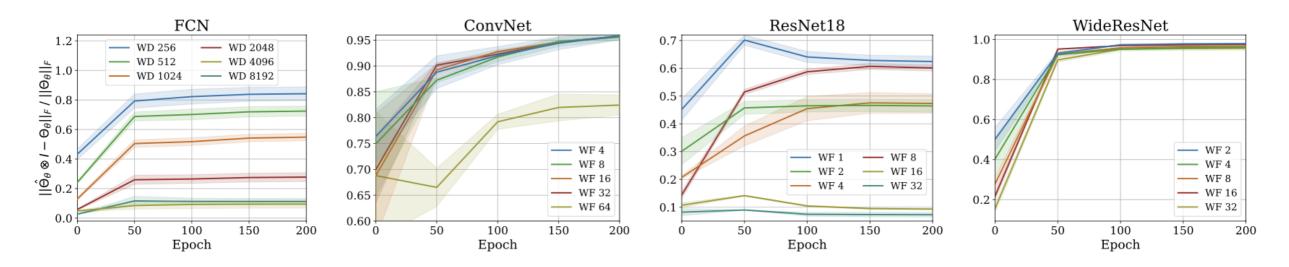


Figure 3: Evaluating the relative difference of Frobenius norm of  $\Theta_{\theta}(\mathcal{D}, \mathcal{D})$  and  $\hat{\Theta}_{\theta}(\mathcal{D}, \mathcal{D}) \otimes I_O$  at initialization and throughout training, based on  $\mathcal{D}$  being 1000 random points from CIFAR-10. Wider nets have more similar  $\|\Theta_{\theta}\|_F$  and  $\|\hat{\Theta}_{\theta} \otimes I_O\|_F$  at initialization.

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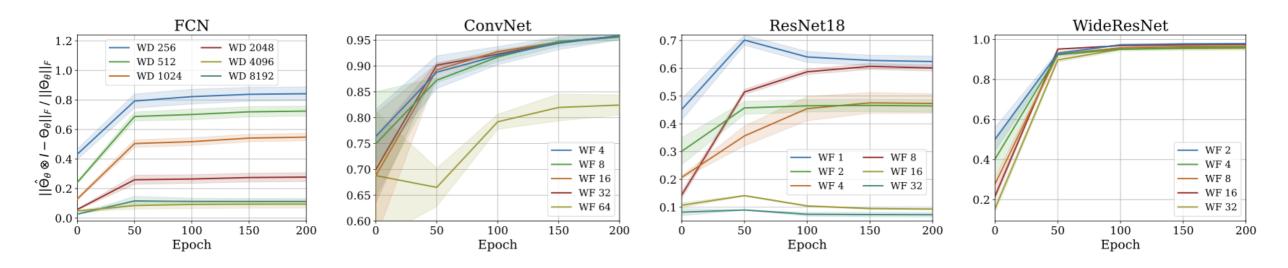


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Same kind of theorem / empirical results for largest eigenvalue, and empirical results for  $\lambda_{\min}$ , condition number

### Kernel regression with pNTK

• Reshape things to handle prediction appropriately:

$$\underbrace{f_{\text{eNTK}}(\tilde{\boldsymbol{x}})}_{k \times 1} = \underbrace{f_0(\tilde{\boldsymbol{x}})}_{k \times 1} + \underbrace{eNTK_{\mathbf{w}_0}(\tilde{\boldsymbol{x}}, \mathbf{X})}_{k \times kN} \underbrace{eNTK_{\mathbf{w}_0}(\mathbf{X}, \mathbf{X})^{-1}}_{kN \times kN} \underbrace{(\mathbf{y} - f_0(\mathbf{X}))}_{kN \times 1}$$
$$\underbrace{f_{\text{pNTK}}(\tilde{\boldsymbol{x}})}_{k \times 1} = \underbrace{f_0(\tilde{\boldsymbol{x}})}_{k \times 1} + \underbrace{(\underbrace{pNTK_{\mathbf{w}_0}(\tilde{\boldsymbol{x}}, \mathbf{X})}_{1 \times N} \underbrace{pNTK_{\mathbf{w}_0}(\mathbf{X}, \mathbf{X})^{-1}}_{N \times N}}_{N \times N} \underbrace{(\mathbf{y} - f_0(\mathbf{X}))}_{N \times k})^{\mathsf{T}}$$

• We have  $\|f_{ ext{eNTK}}( ilde{x}) - f_{ ext{pNTK}}( ilde{x})\| = \mathcal{O}(rac{1}{\sqrt{h}})$  again

• If we add regularization, need to "scale"  $\lambda$  between the two

## **Kernel regression with pNTK**

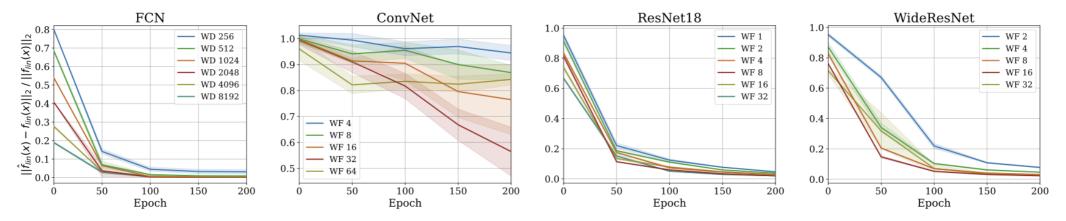


Figure 7: The relative difference of kernel regression outputs, (4) and (5), when training on  $|\mathcal{D}| = 1000$  random CIFAR-10 points and testing on  $|\mathcal{X}| = 500$ . For wider NNs, the relative difference in  $\hat{f}^{lin}(\mathcal{X})$  and  $f^{lin}(\mathcal{X})$  decreases at initialization. Surprisingly, the difference between these two continues to quickly vanish while training the network.

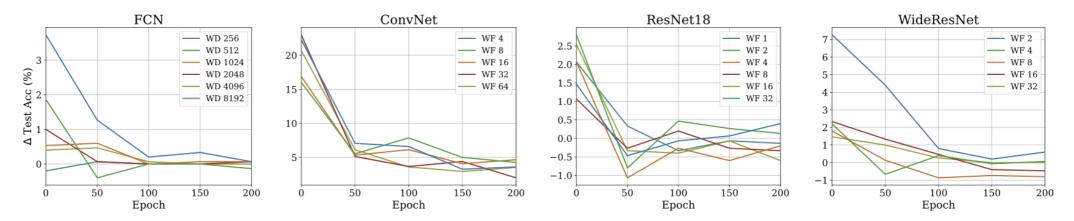


Figure 8: Using pNTK in kernel regression (as in Figure 7) almost always achieves a higher test accuracy than using eNTK. Wider NNs and trained nets have more similar prediction accuracies of  $\hat{f}^{lin}$  and  $f^{lin}$  at initialization. Again, the difference between these two continues to vanish throughout the training process using SGD.

#### pNTK speed-up

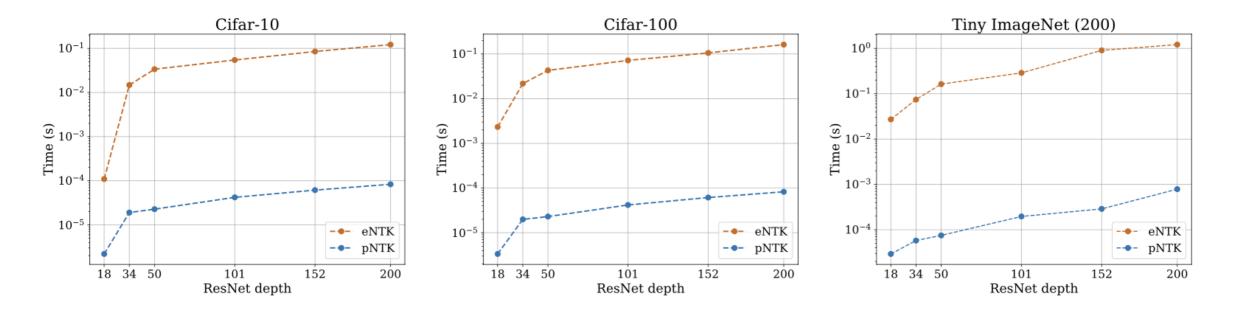
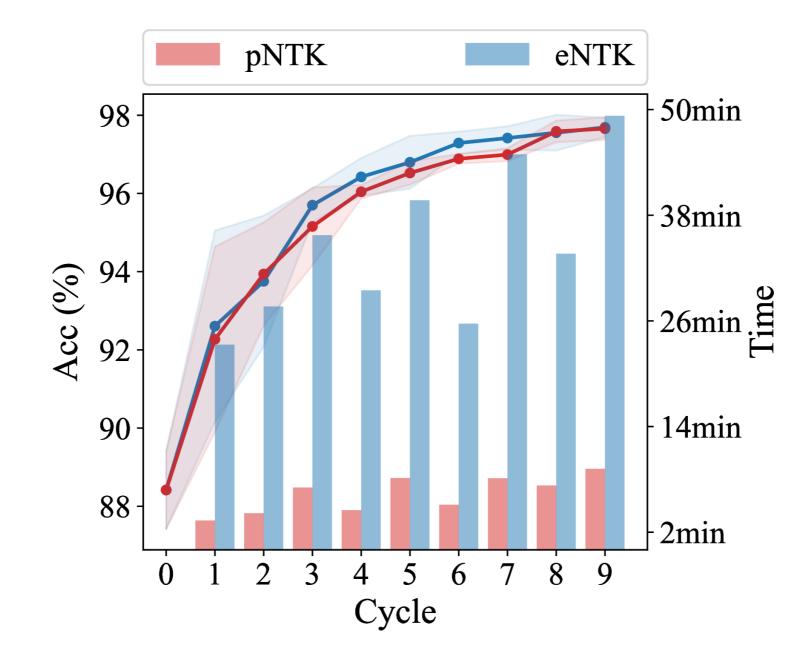


Figure 1: Wall-clock time to evaluate the eNTK and pNTK for one pair of inputs, across datasets and ResNet depths.

### pNTK speed-up on active learning task



## pNTK for full CIFAR-10 regression

- eNTK(X, X) on CIFAR-10: 1.8 terabytes of memory
- pNTK(X, X) on CIFAR-10: 18 gigabytes of memory

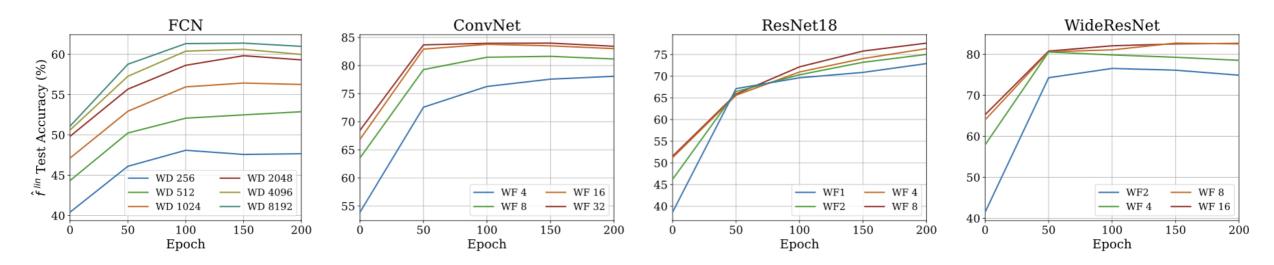


Figure 9: Evaluating the **test accuracy of kernel regression predictions using pNTK as in** (5) **on the full CIFAR-10 dataset**. As the NN's width grows, the test accuracy of  $\hat{f}^{lin}$  also improves, but eventually saturates with the growing width. Using trained weights in computation of pNTK results in improved test accuracy of  $\hat{f}^{lin}$ .

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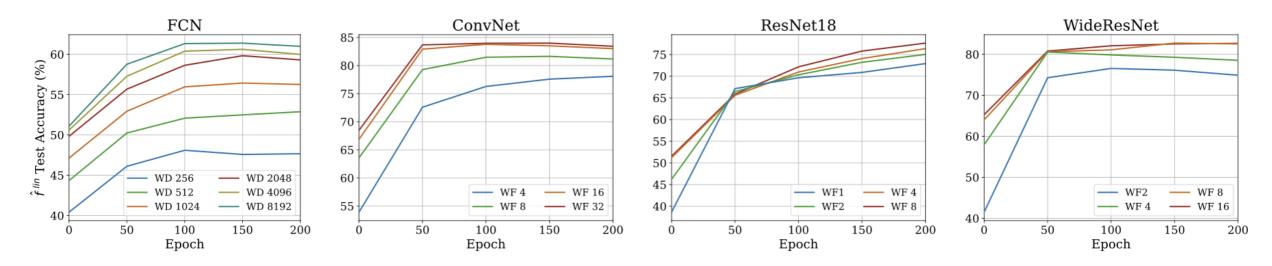


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• Worse than infinite NTK for FCN/ConvNet (where they can be computed, if you try hard)

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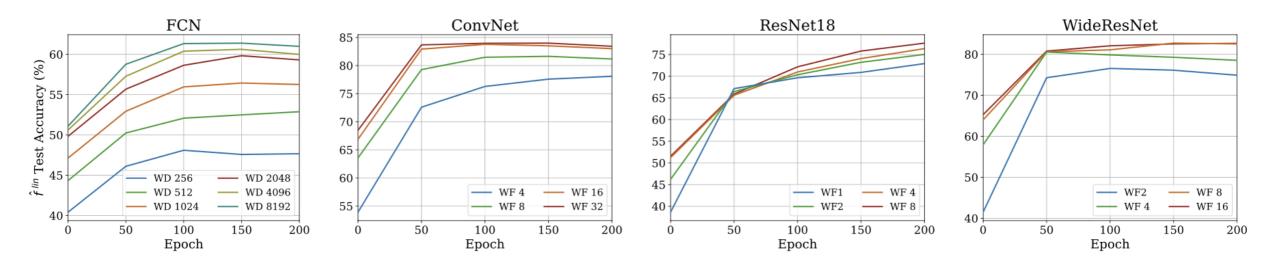


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- Worse than infinite NTK for FCN/ConvNet (where they can be computed, if you try hard)
- Way worse than SGD

### Recap

#### eNTK is a good tool for intuitive understanding of the learning process

Ren, Guo, S. Better Supervisory Signals by Observing Learning Paths

Ren, Guo, Bae, S. How to prepare your task head for finetuning

eNTK is practically very effective at "lookahead" for active learning

Mohamadi\*, Bae\*, S. Making Look-Ahead Active Learning Strategies Feasible with Neural Tangent Kernels

You should probably use pNTK instead of eNTK for high-dim output problems:

Mohamadi, Bae, S. A Fast, Well-Founded Approximation to the Empirical Neural Tangent Kernel