

# Scaling Laws in Linear Regression: Compute, Parameters, and Data

Lin et al.

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Machine Learning and Economics Reading and Discussion Group

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# Motivation: Why Don't Big Models Overfit?

- Empirical observation: test error improves polynomially with model size ( $M$ ) and data size ( $N$ ):

$$\mathcal{R}(M, N) \approx \mathcal{R}^* + \frac{c_1}{M^{a_1}} + \frac{c_2}{N^{a_2}}$$

for irreducible risk  $\mathcal{R}^* > 0$ , constants  $a_1, a_2, c_1, c_2 > 0$  independent of  $M, N$

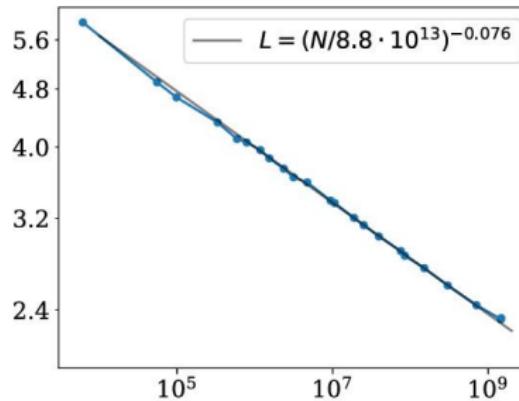
- But from statistical theory:

$$MSE = \text{Bias}^2 + \text{Variance}$$

- Bias decreases as  $M \uparrow$  (better approximation)
- Variance increases as  $M \uparrow$  (overfitting, memorising noise)

# Theory Versus Practice

- **Conflict:** theory predicts “U-shaped” curve:
  - Model improves with size initially
  - Eventually model gets too big for data  $\Rightarrow$  performance crashes
  - But in modern neural networks, we never see the crash



**Figure:** Scaling law for language models: test loss vs. number of parameters (Kaplan et al., 2020).

# Lin et al.: Simplify for Precise Analysis

- Big neural networks are complicated (e.g., transformers)
- **Solution:** focus on simple linear regression case:
  - Input: data  $x$  live in infinite-dimensional space ("true", complex world)
  - Model: can only use  $M < \infty$  covariates  $\Rightarrow$  compress infinite world into  $M$  features via Gaussian sketching<sup>1</sup>
  - Training: one-pass SGD

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<sup>1</sup> $\tilde{x} = Sx$ , where  $S$  is  $M \times \infty$  random Gaussian matrix

# Error Decomposition

- **Decomposition:** risk (error) broken down into:
  - **Approximation Error:** model is too small to represent complexity of data (decreasing in  $M$ )
  - **Bias Error:** insufficient data to converge to best solution (decreasing in  $N$ )
  - **Variance Error:** “memorising” noise in the specific training samples (usually increasing in  $M$ )

# Disappearing Variance

- In one-pass SGD, variance term is higher-order  $\Rightarrow$  effectively vanishes
- Why?
  - Implicit regularisation!
  - SGD prefers “minimum norm” solution even in the absence of explicit penalty term in loss (e.g., Ridge/LASSO)<sup>2</sup>
  - So, model can avoid memorising noise even when capacity  $M$  is sufficient to do so

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<sup>2</sup>Zou et al. (2021).

# Disappearing Variance

- Effective production function:

$$\mathcal{R}(M, N) = \mathcal{R}^* + \Theta\left(\frac{1}{M^{a-1}}\right) + \tilde{\Theta}\left(\frac{1}{(N\gamma)^{(a-1)/a}}\right)$$

  
leading order given by the sum of Approx and Bias

$$\text{Var} = \tilde{\Theta}\left(\frac{\min\{M, (N\gamma)^{1/a}\}}{N}\right)$$

  
higher order, thus unobservable

- Economic interpretation:

- Returns to  $M$  never negative  $\Rightarrow$  only limited by approximation (need bigger  $M$ ), bias (need bigger  $N$ ), or compute ( $C \approx MN$ )

# Allocating Resources

- Optimisation problem:
  - Budget of compute is  $C$
  - Compute cost  $C \approx M \times N$
  - How to choose  $M, N$  to minimise risk?
- Solution:
  - Optimal ratio:  $M \propto C^{\frac{1}{b+1}}, N \propto C^{\frac{b}{b+1}}$ <sup>3</sup>
- Comparison with **Chinchilla**<sup>4</sup>:
  - Famous paper suggesting that  $N$  and  $M$  should scale equally
  - This paper: optimal ratio depends on structure of data through spectral decay  $b$

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<sup>3</sup> $b > 1$  controls the decay of the optimal model parameter  $\mathbf{w}^*$  ( $=$  signal, since  $y = \mathbf{x}^T \mathbf{w}^* + \epsilon$ ), and so is a measure of the difficulty of the task where larger  $b \rightarrow$  simpler.

<sup>4</sup>Hoffman et al. (2022).

# Takeaways

## ■ Key points:

- Variance term negligible  $\Rightarrow$  want to uniformly increase parameters/data
- However, “optimal” AI production function depends on the structure of the data:
  - Harder problems require larger  $N$  (input of intermediate goods?) relative to  $M$  (capital?)

# Key Limitations

## 1 Linear Model:

- Focus on linear setting is tractable, shows scaling applies even in simple settings
- But most models (NNs) are non-linear: features  $\neq$  weights
- Feature learning ("grokking") is arguably where a lot of interesting stuff happens

## 2 One-pass SGD:

- Seeing data once brings theoretical neatness
- But real models train for multiple epochs
- Authors admit: multi-pass SGD may cause variance to return, but their theory can't handle extra complexity

## 3 Data assumptions:

- Assume data have Gaussian distribution, power-law spectrum
- Real data could be meaningfully different (heavy-tailed, structured)  $\Rightarrow$  exponents in scaling law may not hold

# Future Weeks

- Tuesday 10th February (2:30pm): double descent in linear models, presented by Max Kasy
- Tuesday 24th February (2:30pm): empirical scaling laws in LLMs, presented by Thomas Foster
- Tuesday 10th March (2:30pm): scaling and the means of prediction, presented by Aarushi Kalra